657 File: 541.460.000n C.H.



# THE MEMPHIS DEPOT **TENNESSEE**

# **ADMINISTRATIVE RECORD COVER SHEET**

AR File Number 65



STL Pittsburgh 450 William Pitt Way Pittsburgh, PA 15238

Tel 412 820 8380 Fax 412 820 2080 www.stl-inc.com

## ANALYTICAL REPORT

PROJECT NO. UXB 7512-060

Dunn Field, Def Depot Memphis

Lot #: COR130194

Frank Johnson

Memphis Depot

SEVERN TRENT LABORATORIES, INC.

Dave Dunlap Project Manager

# CASE NARRATIVE UXB International Inc. Dunn Field

## LOT # C0E130194

## Sample Receiving:

Samples were received on May 13, 2000. The COC listed only one sample, although four sets of sample bottles were received. It was agreed with the client that the bottles were to be broken out as individual samples and labeled as the sample listed on the COC, with each sample having a suffix to differentiate between them.

## Volatiles:

There were no problems associated with the analyses.

## Semivolatiles:

There were no problems associated with the analyses.

## Pesticides:

Several of the calibration verification standards had compounds outside of the  $\pm 15\%$  difference criteria. The average % difference of all compounds in the calibration verification standards were within the  $\pm 15\%$  difference criteria. Compounds exceeding these criteria that were not detected in the samples. The following table lists the calibration verifications and compounds outside of the  $\pm 15\%$  difference criteria.

Column; #RTX-CLP2			
Calibration Verification Date & Times	Affected Compound	<sup>n</sup> / <sub>e</sub> Difference	Average % Difference of All Compounds in Standard
5/19/00 @ 09:50/10:16	alpha-BHC gamma-BHC	-16.7 -19.8	11
5/19/00 @ 14:10/14:36	Tetrachloro-m- xylene	-16.3	13
	alpha-BHC	-18.3	
	gamma-BHC	-20.6	
1	heptachlor	-16.4	
	endrin	-15.1	

## Herbicides:

There were no problems associated with the analyses.

## Metals:

There were no problems associated with the analyses.

## **TCLP Metals:**

The serial dilution percent difference was outside control limits for barium. All associated results were flagged with an "E" qualifier.

# CASE NARRATIVE UXB International Inc. Dunn Field

## LOT # C0E130194

## General Chemistry:

The reactive cyanide and reactive sulfide analyses were completed at the STL North Canton, OH laboratory.

## C0R130194

PARAMETER	METHOD	PREPARATION METHOD
Chlorinated Herbicides by GC	SW846 8151A	SW846 1311/8150
Ignitability	SW846 SECTION 7	SW846 SECTION 7
Inductively Coupled Plasma (ICP) Metals	SW846 6010B	SW846 1311/3010
Mercury in Liquid Waste (Manual Cold-Vapor)	SW846 7470A	SW846 1311/7470
Organochlorine Pesticides	SW846 8081A	SW846 1311/3510
Reactive Cyanide	SW846 7.3.3	SW846 7.3.3
Reactive Sulfide	SW846 7.3.4	SW846 7.3.4
Semivolatile Organic Compounds by GC/MS	SW846 8270C	SW846 1311/3520
Soil and Waste pH	SW846 9045C	SW846 9045C
Total Residue as Percent Solids	MCAWW 160.3 MOD	MCAWW 160.3 MOD
Trace Inductively Coupled Plasma (ICP) Metals	SW846 6010B	SW846 3050B
Volatile Organics by GC/MS	SW846 8260B	SW846 1311/5030

## References:

MCAWW	"Methods for Chemical Analysis of Water and Wastes", EPA-600/4-79-020, March 1983 and subsequent revisions.
SW846	"Test Methods for Evaluating Solid Waste, Physical/Chemical

"Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 and its updates.

## **SAMPLE SUMMARY**

## C0B130194

WO # 5	SAMPLE#	CLIENT SAMPLE ID	DATE
DD6A4	001	DF/S1/0133/SDC/001 A	05/12/00
DD6A5	002	DF/S1/0133/SDC/001 B	05/12/00
DD6A6	003	DF/S1/0133/SDC/001 C	05/12/00
DD6A7	004	DF/S1/0133/SDC/001 D	05/12/00
NOTE (S'	) <u>-</u>		

- The analytical results of the samples listed above are presented on the following pages.
- All calculations are performed before rounding to avoid round-off errors in calculated results.
- Results noted as "ND" were not detected at or above the stated limit.
- This report must not be reproduced, except in full, without the written approval of the laboratory.
- Results for the following parameters are never reported on a dry weight basis: color, corrosivity, density, flashpoint, ignitability, layers, odor, paint filter test, pH, porosity pressure, reactivity, redox potential, specific gravity, spot tests, solids, solubility, temperature, viscosity, and weight.

Chain of
Custody Record
34878

Quanterra, Inc. - Pittsburgh PA Lab 450 William Pitt Way Pittsburgh PA 15238



Client TASI TASE	Project Manager	CONTROL	Date / CA CO	Chain Of Custody Number 67179
6141 OLD POPLAR PIFC	121	VFax Number	Lab Number	
Memoris State Zacode 119		)		Analysis
Project Name  DUNN FILLO.  Contract/Purchase Order/Quote No	Carner/Waybill Number		TCLP(R	
Sample I D No and Description Date	Time Sample Type Volume	Containers Type No. Preservative		
DF/S1/0133/SOC/001 S/12/001043	043 50(L 4,8,326 GLASS	<del></del>	×	
Special Instructions				
Possible Hazard Identification    Non-Hazard   Flammable   Skin Imitant   P	Polson B Unknown	Sample Disposal  Return To Client	☐ Disposa{ By Lab ☐ Archive For	or Months
	OC Level	Project Specific (Specify)		Date Time
2. Relinquished By	S)12/00 12/16	22	***************************************	5/3/B 13 W
3 Relinquished By	Date	3. Received By		Date
L. Comments				
DISTRIBUTION. WHITE - Stays with Sample, CANARY - Returned to Client with Report; PINK - Field Copy	nt with Report; PINK - Field Copy			

## SBCCOM Monitoring Branch Laboratory

## **CLEARANCE REPORT**

May-12, 2000

Dunn Field, Memphis Defense Depot

Results for CWM Composite Soil Sample Analysis

Analyst: Shawn Heinlein

Sample#		1,4- Thioxane	1,4- Dithiane	TDG	Mustard	Lewisite
DF/S1/01 33/SDC/ 001		ND	ND	N/A	ND	ND
						\ 1
			<u></u>			
					74,	
	,					•
			11-44-y-A-			,
			,	<u> </u>		
			Ť			

ND= Not detected at or above the method detection limit (MDL)

MDL= 200 ppb

BDL= Below detection limit, results > 100ppb, but < 200 ppb

MS= matrix spike

MSD= matrix spike duplicate

DUP= duplicate

657 8

# Cool r Receipt Form STL Pittsburgh

Client:	Project:	_ Quote: _	-A	<del></del>
	Rec'd & Opened for Temp. Check on: 5/13/co		V M	<u> </u>
Cooler	s Opened and Unpacked on:	_ By: <u> </u>	(Signa	ture)
car D	ittsburgh Lot Number:		(Digua	tuic)
21LF	misourgh Lot Number:		Yes	No
			/ /	110
1.	Were custody seals on the outside of the cooler?		<u></u>	<del></del>
	If YES, how many and where? Quantity Location/h	<del> </del>		
	Were signatures and date correct?			<del></del>
2.	Were custody papers included inside the cooler?			
3.	Were custody papers properly filled out (ink, signed, match labels)?			<del></del>
4.	Did you sign the custody papers in the appropriate place?		/	
5.	Was shippers packing slip attached to this form?			
6.	Were packing materials used?		-,-	
	If YES, what type?			_
7.	Were the samples chilled? (Record temperatures on reverse side.)			
8.	Were the samples appropriately preserved?		M	
9.	Were all bottles sealed in separate plastic bags?		_	
10.	Did all bottles arrive in good condition (unbroken)?		-/-	
11.	Were all bottle labels complete (sample ID, preservatives, etc.)?	<del> := : : :</del>		
12.	Did all bottle labels and/or tags agree with custody papers?			
13.	Were correct bottles used for tests indicated?			
14.	Were all VOA vials checked for the presence of air bubbles?	<del>,</del> -	_	
15.	Was a sufficient amount of sample sent in each bottle?			
16.	Samples received by: FEDEX UPS CLIENT DROP-OFF OTHER	AIRBOI		
Expl	ain any discrepancies:			
Leve	2 Review	to resolve d	liscrena	ncies
Was	contacted on by	_ 10 1620146 G	reserve	

STL PT/May-00/96-005/COOLDOC

Page 1 of 2

STL Pittsburgh

P: Preserved UP: Unpreserved

Sample ID	TMET PH<2	DMET PH<2	HG PH<2	NUT(1) PH<2	CN PH ≥12	OG TPHC PH<2	PHEN PH<2	SULF PH ≥12	TOC PH<2	TOX PH<2	VOA P/UP	hrdnss PH<2			
															T
<u></u>		<del> </del>													1
		1	1				_								7
		<del>                                     </del>								Ī					T
															7
															T
															I
							L								
										<u> </u>			<u> </u>		
									<u> </u>			<u> </u>	<u> </u>	<u> </u>	
						<u> </u>			<u> </u>	ļ <u>.</u>	<u> </u>			<del></del>	
						<u> </u>			<u> </u>	<u> </u>	<u> </u>		<u> </u>	<u> </u>	
		<u> </u>	<u> </u>			ļ		ļ	<u> </u>	<u> </u>		ļ	ļ	<del> </del>	_
						ļ	ļ		<u> </u>	<b> </b>		<u> </u>	<b> </b>	╀	4
			<del> </del>	<u> </u>		ļ			<del>                                     </del>	<u> </u>	<del>                                     </del>	<del> </del>		<del></del>	4
		<b></b>	<del>                                     </del>		ļ	<del> </del>	<u> </u>		<del> </del>	<del>                                     </del>	<del> </del>	-	<del> </del>	<b>↓</b>	4
			<u> </u>		-	<del> </del>	<del>                                     </del>	<del> </del>	<del> </del>	<b> </b>	<del> </del>		<del> </del>	<del> </del>	_
			1	<b></b>		<del>  </del>	<del> </del>		ļ	<del> </del> _			ļ	<del> </del>	_

Comments:

Cooler Number	Temperature		Bottle Type	Lot Number*
	3'			
		1 1		
	<u> </u>	╣		
		1		

\* Please use an asterisk if bottle lot number was covered by the label.

	Packages up to 150 fbs.  Devery commons may be use in some seen.  Hard because alemone devery by a see it come a seen.  Hard because alemone devery be selective companies.  Fedex Englises Saver*  Tedex Companies on the selective companies on the selective companies.	Second barrages der Treight  Telex 20 ay Freight	Pecker Park & Octave Prog.	Sunday Dalivery HOLD Wheekday HOLD Saturday Assatoris redes Proch 1 at Fed Ex Location Assatoris resecutive Comparts and Fed Ex Location Researched Fed Ex Location Researched Fed Ex Demography Comparts and Fed Ex Bongs Regerous goods?	Ves		a heyber valve. See back for details:  mitted obtaining agrains.  Twithout dibtaining a segment.  Twithout dibtaining a segment.  Twithout dibtaining a segment.  Twithout dibtaining a segment.
Fedia madeleral months of the second	Sender's Jim / Monte Lo I Flore To I US -U 999   Sender's Day.	Company CXXS 4b Express Freight Service  Address 6141 CLO POLAR ARC Costs Communication Costs Communication Costs Communication Costs Costs Confined Costs Confined Costs Confined Costs Confined Costs Confined Costs Confined Costs Co	<sub>[],</sub>	To Secret Handling Asturday Delivery Recipients Rustry Vicinic Phone 412 820-4380 Destuday Delivery Name Destuday Delivery Name Destuday Delivery Destuday Destuday Delivery Delivery Destuday Delivery Destuday Delivery Destuday Delivery Destuday Delivery D	No 1/8s section 2/8s section 2/	917 Sale P.A. 219 15238	8 Release Signature spousorous dolongy as through the Stocker that shipmen without a mind to the Stocker that shipmen to describe the Stocker that shipmen to describe the Stocker that shipmen the Stocker that ship the Stocker t

DATA SUMMARY PACKAGE

## **GC/MS VOLATILE SUMMARY**

657 13

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID

Lab Sample ID: C0E130194 001

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD6A4101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

Moisture %:10

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 A

CAS NO.	COMPOUND (ug/L c	r ug/kg) mg/L	0
71-43-2	Benzene	0.050	U
78-93-3	2-Butanone	0.027	J
56-23-5	Carbon tetrachloride	0.050	i
108-90-7	Chlorobenzene	0.050	<u>י</u>
67-66-3	Chloroform	0.050	Ū
107-06-2	1,2-Dichloroethane	0.050	<u>י</u>
75-35-4	1,1-Dichloroethene	0.050	Ū
127-18-4	Tetrachloroethene	0.050	U
79-01-6	Trichloroethene	0.050	ט
75-01-4	Vinyl chloride	0.050	ט

## UXB INTERNATIONAL MATRIX SPIKE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 001

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD6A410P Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

Moisture %:10

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 A

CAS NO.	COMPOUND (ug/L o	r ug/kg) mg/L Q
71-43-2	Benzene	0.544
78-93-3	2-Butanone	0.532
56-23-5	Carbon tetrachloride	0.563
108-90-7	Chlorobenzene	0.546
67-66-3	Chloroform	0.536
107-06-2	1,2-Dichloroethane	0.552
75-35-4	1,1-Dichloroethene	0.518
127-18-4	Tetrachloroethene	0.555
79-01-6	Trichloroethene	0.545
75-01-4	Vinyl chloride	0.635

# UXB INTERNATIONAL MATRIX SPIKE DUPLICATE COMPOUNDS

657 15

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 001

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD6A410Q Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

Moisture %:10

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 A

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L .Q
71-43-2	Benzene	0.552
78-93-3	2-Butanone	0.526
56-23-5	Carbon tetrachloride	0.571
108-90-7	Chlorobenzene	0.550
67-66 <b>-</b> 3	Chloroform	0.541
107-06-2	1,2-Dichloroethane	0.561
75-35-4	1,1-Dichloroethene	0.522
127-18-4	Tetrachloroethene	0.567
79-01-6	Trichloroethene	0.551
75-01-4	Vinyl chloride	0.638

## 657 16

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 002

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD6A5101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

Moisture %:8.6

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 B

CAS NO.	COMPOUND (ug/L c	r ug/kg) mg/L	0
71-43-2	Benzene	0.050	Ū
<u>78-93-3</u>	2-Butanone	0.027	J
56-23-5	Carbon tetrachloride	0.050	U
108-90-7	Chlorobenzene	0.050	U
67-66-3	Chloroform	0.050	Ū
107-06-2	1,2-Dichloroethane	0.050	U
75-35-4	1,1-Dichloroethene	0.050	U
127-18-4	Tetrachloroethene	0.050	U
79-01-6	Trichloroethene	0.050	
75-01-4	Vinyl chloride	0.050	U

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:COE130194 003

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD6A6101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

Moisture %:10

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 C

CAS NO.	COMPOUND (ug/L c	or ug/kg) mg/L	0
71-43-2	Benzene	0.050	ال
78-93-3	2-Butanone	0.023	J
56-23-5	Carbon tetrachloride	0.050	Ū
108-90-7	Chlorobenzene	0.050	U
67-66-3	Chloroform	0.050	U
107-06-2	1,2-Dichloroethane	0.050	Ū
75-35-4	1,1-Dichlorcethene	0.050	יט ו
127-18-4	Tetrachloroethene	0.050	Ū
79-01-6	Trichloroethene	0.050	וֹט
75-01-4	Vinyl chloride	0.050	i u

## 657 18

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 004

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD6A7101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

Moisture %:11

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 D

CAS NO.	COMPOUND (ug/L c	or ug/kg) mg/L	Q
71-43-2	Benzene	0.050	ו
78-93-3	2-Butanone	0.039	J
56-23-5	Carbon tetrachloride	0.050	U
108-90-7	Chlorobenzene	0.050	<u>ט</u>
<u>67-66-3</u>	Chloroform_	0.050	ן ש
107-06-2	1,2-Dichloroethane	0.050	U
75-35-4	1,1-Dichloroethene	0.050	ט ו
127-18-4	Tetrachloroethene	0.050	
79-01-6	Trichloroethene	0.050	<u></u>
75-01-4	Vinyl chloride	0.050	<u>ט</u>

657 19

## UXB INTERNATIONAL CHECK SAMPLE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:COE170000 262

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DDA08101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

Moisture %:NA

QC Batch: 0138262

Client Sample Id: CHECK SAMPLE

CAS NO.	COMPOUND (ug/L or	<u>rug/kg)</u> mg/L Q
71-43-2	Benzene	[0.546 ]
78-93-3	2-Butanone	0.503
56-23-5	Carbon tetrachloride	0.567
108-90-7	Chlorobenzene	0.543
67-66-3	Chloroform	0.540
107-06-2	1,2-Dichloroethane	0.552
75-35-4	1,1-Dichloroethene	0.510
127-18-4	Tetrachloroethene	0.566
79-01-6	Trichloroethene	0.545
75-01-4	Vinyl chloride	0.627

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

QESSDG:

Lot #: C0E130194

Ī	CLIENT ID.	SRG01	SRG02	SRG03	SRG04	TOT OUT
		======	======	======	=======	=====
	DF/S1/0133/SDC/001 A	110	103	93	_103	00
02	DF/S1/0133/SDC/001 B	108	99	_89	104	00
03	DF/S1/0133/SDC/001 C	106	106	93	102	00
04	DF/S1/0133/SDC/001 D	119	96	91	107	00
05	METHOD BLK. DD95L101	104	100	90	98	_00
,	LCS DDA08101	103	104	103	102	00
07	DF/S1/0133/SDC/001 A D	103	103	102	102	00
80	DF/S1/0133/SDC/001 A S	102	104	102	103	00

SURROGA	TES		QC LIMITS
SRG01	=	1,2-Dichloroethane-d4	(77-120)
SRG02	=	Toluene-d8	( 78-111)
SRG03	=	4-Bromofluorobenzene	( 80-114)
SRG04	=	Dibromofluoromethane	( 78-110)

FORM II

<sup>#</sup> Column to be used to flag recovery values

<sup>\*</sup> Values outside of required QC Limits

D System monitoring Compound diluted out

#### SW846 8260B METHOD BLANK SUMMARY

BLANK WORKORDER NO. DD95L101

Lab Name: Severn Trent Laboratories, Inc.

Lab Code: QESPIT

SDG Number:

Lab File ID: 4051703.d

Lot Number: COE130194

Date Analyzed: 05/17/00

Time Analyzed: 14:39

Matrix: SOLID

Date Extracted: 05/17/00

GC Column: RTX-624 ID: .18

Extraction Method: 1311/5030B

Instrument ID: HP4

Level: (low/med) LOW

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, LCS, LCSD, MS, MSD:

-		SAMPLE	LAB	DATE	TIME
1	CLIENT ID.	WORK ORDER #	FILE ID	ANALYZED	ANALYZED
		******	=======================================		_======
01	CHECK SAMPLE	DDA08101 C	4051707.d	05/17/00	17:04
02	DF-001	DD6A410P S	4051705.d	05/17/00	16:07
	DF-001	DD6A410Q D	4051706.d	05/17/00	16:35
04	DF-001	DD6A4101	4051708.d	05/17/00	17:39
05	<u>S1-001</u>	DD6A5101	4051709.d	05/17/00	18:08
06	0133/001	DD6A6101	4051712.d	05/17/00	19:29
07	SDC/001	DD6A7101	4051711.d	05/17/00	19:00
98					
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22	••				
23					
24					
25					
26					
27					
28					
29					
30					

COMMENTS:		
		_
	FORM IV	

657 22 SW846 8260B CHECK SAMPLE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Lot #: COE170000

WO #: DDA08101 BATCH: 0138262

COMPOUND	SPIKE ADDED (mg/L)	SAMPLE CONCENT. (mg/L)	% REC	QC LIMITS REC	QUAL
Benzene	0.500	0.546	109	79 - 116	,
2-Butanone	0.500	0.503	101	35- 156	
Carbon tetrachloride	0.500	0.567	113	72 - 133	
Chlorobenzene	0.500	0.543	109	81 - 115	
Chloroform	0.500	0.540	108	81- 122	
1,2-Dichloroethane	0.500	0.552	110	73 - 127	
1,1-Dichloroethene	0.500	0.510	102	65 - 119	
Tetrachloroethene	0.500	0.566	113	78 - 131	
Trichloroethene	0.500	0.545	109	80- 122	
Vinyl chloride	0.500	0.627	125	53 - 134	

* Values outside	of QC limits					
Spike Recovery:	0 out of	10 o	utside 1	limits		
COMMENTS:						
<del></del>					 	

FORM III

NOTES (S):

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A

Level: (low/med) LOW

Lot #: C0E130194

WO #: DD6A410P BATCH: 0138262

] ]	SPIKE	SAMPLE	MS	MS		
l dosenorma	ADDED	CONCENT.	CONCENT.	*	LIMITS	
COMPOUND	(mg/L )	(mg/L)	(mg/L)	REC	REC	QUAL
	=======	=======	========	=====		
Benzene	0.500	ND	0.544	109	73 - 123	Ì
2-Butanone	0.500	0.027	0.532	101	10- 151	¦
Carbon tetrachloride	0.500	ND	0.563	113	61 - 143	<del></del>
Chlorobenzene	0.500	ND	0.546	109	70 - 122	
Chloroform	0.500	ND	0.536	107	65 - 131	
1,2-Dichloroethane	0.500	ND	0.552	110	67 - 132	<u> </u>
1,1-Dichloroethene	0.500	ND	0.518	104	57- 138	! <del></del>
Tetrachloroethene	0.500	ND	0.555	111	70 - 130	
Trichloroethene	0.500	ND	0.545	109	58 - 141	! 
Vinyl chloride	0.500	ND	0.635	127	51 - 133	

Column	to	be	used	to	flag	recovery	and	RPD	values	with	an	asterisk	

RPD: Spike	0 out Recovery:	of <u>0</u> 0 out	outside lin	mits outside	limits	
COMME	NTS:					
				-		- <u>, , </u>

FORM III

NOTES (S):

SW846 8260B MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

657

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A

Level: (low/med) LOW

Lot #: C0E130194

WO #: DD6A410Q BATCH: 0138262

	SPIKE	MSD	MSD				
l	ADDED	CONCENT.	8	ફ	QC I	LIMITS	ļ
COMPOUND	(mg/L)	(mg/L)	REC	RPD	RPD	REC	QUAL
	22222222		=====	======	====	=======	
Benzene	0.500	0.552	110	1.5	20	73 - 123	
2-Butanone	0.500	0.526	100	1.1	34	10- 151	
Carbon tetrachloride	0.500	0.571	114_	1.5	20	61- 143	
Chlorobenzene	0.500	0.550	110	0.69	20	70 - 122	
Chloroform	0.500	0.541	108	0.79	20	65 - 131	
1,2-Dichloroethane	0.500	0.561	112	1.6	20	67- 132	
1,1-Dichloroethene	0.500	0.522	104	0.78	20	57- 138	
Tetrachloroethene	0.500	0.567	113	2.2	20	70- 130	
Trichloroethene	0.500	0.551	110	1.1	20	58- 141	i
Vinyl chloride	0.500	0.638	128	0.50	20	51- 133	

NOTES (S):					
	 	 	 	 	-

# Column to be used to flag recovery and RPD values with an asterisk \* Values outside of QC limits

RPD:	0	out	of	_	10	outside	li	mits	
Spike	Recove	erv:		0	out	of	10	outside	1 im

ke Recovery: 0 out of 10 outside limits

COMMENTS:

FORM III

## SW846 8260B METHOD BLANK SUMMARY

BLANK WORKORDER NO.

657 2

Lab Name: Severn Trent Laboratories, Inc.

Lab Code: QESPIT

SDG Number:

Lab File ID: 4051703.d

Lot Number: COE130194

Date Analyzed: 05/17/00

Time Analyzed: 14:39

Matrix: SOLID

Date Extracted:05/17/00

GC Column: RTX-624 ID: .18

Extraction Method: 1311/5030B

Instrument ID: HP4

Level: (low/med) LOW

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, LCS, LCSD, MS, MSD:

Ī		SAMPLE	LAB	DATE	TIME
	CLIENT ID.	WORK ORDER #	FILE ID	ANALYZED	ANALYZED
	=======================================		=========		=======
	CHECK SAMPLE	DDA08101 C	4051707.d	05/17/00	17:04
	DF/S1/0133/SDC/001 A	DD6A410P S	4051705.d	05/17/00	16:07
	DF/S1/0133/SDC/001 A	DD6A4100 D	4051706.d	05/17/00	16:35
	DF/S1/0133/SDC/001 A	DD6A4101	4051708.d	05/17/00	17:39
	DF/S1/0133/SDC/001 B	DD6A5101	4051709.d	05/17/00	18:08
	DF/S1/0133/SDC/001 C	DD6A6101	4051712.d	05/17/00	19:29
	DF/S1/0133/SDC/001 D	DD6A7101	4051711.d	05/17/00	19:00
80					
09				l <u></u>	
10					
11					
12				<u></u>	
13					i
14	•			<u> </u>	
15				l	
16					
17					
18 19					
20					
21			<u> </u>		<u> </u>
22			<del></del>		ļ
23			ļ		
24			ļ	<u> </u>	
25			<u> </u>		
26			<u> </u>	<u> </u>	[I
27			l i	ļ——	<u></u> !
28					ļ
29			<u> </u>		<del></del>
30				l ————————————————————————————————————	
			1·		1

COMMENTS:									
		<del></del>	<del></del> -	 					<del>-</del>
	<del></del>			 				·····	
					1	ORM	IV		

# UXB INTERNATIONAL METHOD BLANK COMPOUNDS

657 26

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E170000 101

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD95L101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

Moisture %:NA

QC Batch: 0138262

Client Sample Id: INTRA-LAB BLANK

CAS NO.	COMPOUND (ug/L o	r ug/kg) mg/L Q	!
71-43-2	Benzene	0.050	וס
78-93-3	2-Butanone	0.050	<u>י</u>
<u>56-23-5</u>	Carbon tetrachloride	0.050	וס
108-90-7	Chlorobenzene	0.050	<u>ס</u>
67-66-3	Chloroform	0.050	ט
107-06-2	1,2-Dichloroethane	0.050	ָט
75-35-4	1,1-Dichloroethene	0.050	
127-18-4	Tetrachloroethene	0.050	וט
79-01-6	Trichloroethene	0.050	U
75-01-4	Vinyl chloride	0.050	ַ

Lab Name: STL-PITTSBURGH Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: C0E130194

Lab File ID (Standard): CC40517 Date Analyzed: 05/17/00

Instrument ID: HP4 Time Analyzed: 1108

GC Column: DB 624 ID: 0.18 (mm) Heated Purge: (Y/N) N

		IS1(CBZ)		IS2 (DCB)		IS3	
,		AREA #	RT #	AREA #	RT #	AREA #	RT #
		=========	======	=======	======		======
	12 HOUR STD	149779	10.12	227254	12.46	660677	6.99
	UPPER LIMIT	299558	10.32	454508	12.66	1321354	7.19
	LOWER LIMIT	74890	9.92	113627	12.26	330338	6.79
	EPA SAMPLE	=======	======	========	=====	=======	======
	NO.						
			======		======	========	======
01	INTRA-LAB BL	154881	10.12	210752	12.46	792190	7.01
DF 51 0133 90002	DF-001 A MS	159002	10.12	229690	12.46	736767	7.00
	DF-001 & MS		10.13	232577	12.46	739353	7.00
, , , , 04	INTRA-LAB CH	163960	10.13	229414	12.46	743134	7.00
क्रिडा विद्यानिक वर्ष	DF-00₽ A	131368	10.13	173160	12.46	666337	7.01
06		135805	10.13	180580	12.46	677588	7.01
07	SDC/001 D	123751	10.13	170776	12.46	585666	7.01
V 08	0133 <del>/001</del> C	142009	10.13	196867	12.46	736808	7.01
10							<del></del>
11							
12							
13						•	,
14		<del> </del>					
15							
16							
17							ļ <u></u>
18						ļ	
19							
20 21							
21 22			<u> </u>	ļ	<del></del>		
22	l		l	1	l	i	l

IS1 (CBZ) = Chlorobenzene-d5

IS2 (DCB) = 1,4-Dichlorobenzene-d4

IS3 = Fluorobenzene

AREA UPPER LIMIT = +100% of internal standard area AREA LOWER LIMIT = - 50% of internal standard area

RT UPPER LIMIT = + 0.20 minutes of internal standard RT RT LOWER LIMIT = - 0.20 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

page 1 of 1

FORM VIII VOA

OLM03.0

GC/MS SEMIVOLATILE SUMMARY

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 001

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DD6A4102 Date Extracted:05/18/00 Dilution factor. 1 Date Analyzed: 05/22/00

Moisture %:10

QC Batch: 0140269

Client Sample Id: DF/S1/0133/SDC/001 A

CAS NO.	COMPOUND (ug/L or	r ug/kg) mg/L	Q
106-46-7	1,4-Dichlorobenzene	0.050	الا
121-14-2	2,4-Dinitrotoluene	0.050	וֹט
118-74-1	Hexachlorobenzene	0.050	ן ט
87-68-3	Hexachlorobutadiene	0.050	i
67-72-1	Hexachloroethane	0.050	ן ט
98-95-3	Nitrobenzene	0.050	
87-86-5	Pentachlorophenol	0.25	U
110-86-1	Pyridine	0.10	וט
95-95-4	2,4,5-Trichlorophenol	0.050	U
88-06-2	2,4,6-Trichlorophenol	0.050	וֹט
1319-77-3	Cresols (total)	0.050	ט

# UXB INTERNATIONAL MATRIX SPIKE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:COE130194 001

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol 200 / mL Date Received: 05/13/00 Work Order DD6A411G Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

Moisture %:10

QC Batch: 0140269

Client Sample Id: DF/S1/0133/SDC/001 A

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L Q
106-46-7	1,4-Dichlorobenzene	0.127
121-14-2	2,4-Dinitrotoluene	0.118
118-74-1	Hexachlorobenzene	0.174
87-68-3	Hexachlorobutadiene	0.130
67-72-1	Hexachloroethane	0.118
98-95-3	Nitrobenzene	0.131
87-86 <b>-</b> 5	Pentachlorophenol	0.138
110-86-1	Pyridine	0.157
95-95-4	2,4,5-Trichlorophenol	0.142
88-06-2	2,4,6-Trichlorophenol	0.140
1319-77-3	Cresols (total)	0.375

# UXB INTERNATIONAL MATRIX SPIKE DUPLICATE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 001

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DD6A411H Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

Moisture %.10

QC Batch: 0140269

Client Sample Id: DF/S1/0133/SDC/001 A

CAS NO.	COMPOUND (ug/L o	rug/kg) mg/L Q
106-46-7	1,4-Dichlorobenzene	0.131
121-14-2	2,4-Dinitrotoluene	0.120
118-74-1	Hexachlorobenzene	0.181
87-68-3	Hexachlorobutadiene	0.135
67-72-1	Hexachloroethane	0.122
98-95-3	Nitrobenzene	0.136
87-86-5	Pentachlorophenol	0.143
110-86-1	Pyridine	0.164
95-95-4	2,4,5-Trichlorophenol	0.146
88-06-2	2,4,6-Trichlorophenol	0.143
1319-77-3	Cresols (total)	0.391

657 32

#### UXB INTERNATIONAL

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 002

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DD6A5102 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

Moisture %:8.6

QC Batch: 0140269

Client Sample Id DF/S1/0133/SDC/001 B

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L	Q
106-46-7	1,4-Dichlorobenzene	0.050	ש
121-14-2	2,4-Dinitrotoluene	0.050	U
118-74-1	Hexachlorobenzene	0.050	U
87-68-3	Hexachlorobutadiene	0 050	ש
67-72-1	Hexachloroethane	0.050	U
98-95-3	Nitrobenzene	0.050	ַ ָ
87-86-5	Pentachlorophenol	0.25	U
110-86-1	Pyridine	0.10	U
95-95-4	2,4,5-Trichlorophenol	0.050	ט
88-06-2	2,4,6-Trichlorophenol	0.050	ט
1319-77-3	Cresols (total)	0.050	U

Lab Name · Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 003

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order DD6A6102 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

Moisture %:10

QC Batch: 0140269

Client Sample Id: DF/S1/0133/SDC/001 C

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L	Q
106-46-7	1,4-Dichlorobenzene	0.050	ן די
121-14-2	2,4-Dinitrotoluene	0.050	U
118-74-1	Hexachlorobenzene	0.050	ן ט
87-68-3	Hexachlorobutadiene	0.050	ן ט
67-72-1	Hexachloroethane	0.050	ן ט
98-95-3	Nitrobenzene	0.050	ן די
87-86-5	Pentachlorophenol	0.25	U
110-86-1	Pyridine	0.10	וט
95-95-4	2,4,5-Trichlorophenol	0.050	ָט
88-06-2	2,4,6-Trichlorophenol	0.050	ן ט
1319-77-3	Cresols (total)	0.050	ן ט

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix (soil/water) SOLID Lab Sample ID:COE130194 004

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DD6A7102 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

Moisture %:11

QC Batch: 0140269

Client Sample Id: DF/S1/0133/SDC/001 D

CAS NO.	COMPOUND (ug/L or	r ug/kg) mg/L (	2
106-46-7	1,4-Dichlorobenzene	0.050	ן ט
121-14-2	2,4-Dinitrotoluene	0.050	U
118-74-1	Hexachlorobenzene	0.050	U
87-68-3	Hexachlorobutadiene	0.050	ַן
67-72-1	Hexachloroethane	0.050	U
98-95-3	Nitroben <b>zene</b>	0.050	U
87-86 <b>-</b> 5	Pentachlorophenol	0.25	ט
110-86-1	Pyridine	0.10	<u> </u>
95-95-4	2,4,5-Trichlorophenol	0.050	<u>  U</u>
88-06-2	2,4,6-Trichlorophenol	0.050	ט
1319-77-3	Cresols (total)	0.050	<u>                                     </u>

# UXB INTERNATIONAL CHECK SAMPLE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID.C0E190000 269

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DDF2D102 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

Moisture % · NA

QC Batch: 0140269

Client Sample Id: CHECK SAMPLE

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L Q
106-46-7	1,4-Dichlorobenzene	0.173
121-14-2	2,4-Dinitrotoluene	0.136
118-74-1	Hexachlorobenzene	0.191
87-68-3	Hexachlorobutadiene	0.179
67-72-1	Hexachloroethane	0.170
98-95-3	Nitrobenzene	0.168
87-86-5	Pentachlorophenol	0.162
110-86-1	Pyridine	0.272
95-95-4	2,4,5-Trichlorophenol	0.171
88-06-2	2,4,6-Trichlorophenol	0.173
1319-77-3	Cresols (total)	0.489

### SW846 8270C SURROGATE RECOVERY

Lab Name: Severn Trent Laboratories, Inc Client: UXB INTERNATIONAL

Lab Code: QESPIT

QESSDG:

Lot #: C0E130194

Ī	CLIENT ID.	SRG01	SRG02	SRG03	SRG04	SRG05	SRG06	TOT OUT
		======	======	======	======		======	=====
01	DF/S1/0133/SDC/001 A	51	56	78	45	47	56	00 1
02	DF/S1/0133/SDC/001 B	55	57	75	49	51	59	00
03	DF/S1/0133/SDC/001 C	47	46	66	42	45	50	00
04	DF/S1/0133/SDC/001 D	53	54	67	48	51	59	00
05	METHOD BLK. DDF2D101	64	68	67	63	66	66	00
06	LCS DDF2D102	65	72	69	61	63	67	00
07	DF/S1/0133/SDC/001 A D	54	58	73	47	51	63	00
08	DF/S1/0133/SDC/001 A S	52	57	71	46	49	60	00

SURROGATES		ES		QC LIMITS
	SRG01	=	Nitrobenzene-d5	( 32-112)
	SRG02	=	2-Fluorobiphenyl	( 30-110)
	SRG03	=	Terphenyl-d14	( 10-144)
	SRG04	=	2-Fluorophenol	( 13-110)
	SRG05	=	Phenol-d5	( 10-113)
	SRG06	=	2,4,6-Tribromophenol	(21-122)

FORM II

<sup>#</sup> Column to be used to flag recovery values

<sup>\*</sup> Values outside of required QC Limits

D System monitoring Compound diluted out

### SW846 8270C CHECK SAMPLE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Lot #: C0E190000

WO #: DDF2D102 BATCH: 0140269

COMPOUND	SPIKE ADDED (mg/L )	SAMPLE CONCENT. (mg/L )	% REC	QC LIMITS REC	    QUAL
		==============	=====		  =========
1,4-Dichlorobenzene	0.250	0.173	69	28- 110	,
2,4-Dinitrotoluene	0.250	0.136	54	47- 131	¦
Hexachlorobenzene	0.250	0.191	77	57- 128	¦ <del></del>
Hexachlorobutadiene	0.250	0.179	72	36- 116	¦———
Hexachloroethane	0.250	0.170	68	30- 110	<del></del>
Nitrobenzene	0.250	0.168	67	45- 130	¦ —
Pentachlorophenol	0.250	0.162	65	10- 140	¦
Pyridine	0.250	0.272	109	10- 148	!
2,4,5-Trichlorophenol	0.250	0.171	69	41- 125	¦
2,4,6-Trichlorophenol	0.250	0.173	69	46- 135	! <del></del>
Cresols (total)	0.750	0.489	65	29- 144	!

Values outside of QC limits	
olke Recovery:0 out of11 outside limits	
OMMENTS:	

NOTES (S):

## 38 SW846 8270C MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name · Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code · QESPIT

SDG No:

Matrix Spike ID. DF/S1/0133/SDC/001 A Level: (low/med) LOW

Lot #: C0E130194

WO #: DD6A411G BATCH: 0140269

COMPOUND	SPIKE ADDED (mg/L)	SAMPLE CONCENT. (mg/L)	MS CONCENT. (mg/L)	MS % REC	LIMITS REC	QUAL
1,4-Dichlorobenzene	10,250	ND	0.127	]====== 51	18- 110	====================================
2,4-Dinitrotoluene	0.250	ND	0.118	47	31- 131	! ——-
Hexachlorobenzene	0.250	ND	0.174	70	36- 132	
Hexachlorobutadiene	0.250	ND	0 130	52	18- 116	
Hexachloroethane	0.250	ND	0 118	47	18- 110	
Nitrobenzene	0.250	ND	0.131	52	10- 211	
Pentachlorophenol	0.250	ND	0.138	55	10- 140	
Pyridine	0.250	ND	0.157	63	10- 148	
2,4,5-Trichlorophenol	0.250	ND	0.142	57	24- 143	
2,4,6-Trichlorophenol	0.250	ND	0.140	56	36- 135	
Cresols (total)	0.750	ND	0.375	50	25- 144	

NOTES(S):		

RPD: 0 out of 0 outside limits Spike Recovery: \_\_\_0 out of \_\_\_11 outside limits COMMENTS:

<sup>#</sup> Column to be used to flag recovery and RPD values with an asterisk

<sup>\*</sup> Values outside of QC limits

## SW846 8270C MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A Level: (low/med) LOW

Lot #: C0E130194 WO #. DD6A411H

BATCH: 0140269

	SPIKE ADDED	MSD CONCENT.	MSD %	*	OC 1	LIMITS	   
COMPOUND	(mg/L)	(mg/L )	REC	RPD	RPD	REC	QUAL
	========		=====	======	====	========	=======
1,4-Dichlorobenzene	0.250	0.131	53	3.2	36	18- 110	İ
2.4-Dinitrotoluene	0.250	0.120	4.8	1.4	32	31- 131	
Hexachlorobenzene	0.250	0.181	72	4.1	22	36- 132	
Hexachlorobutadiene	0.250	0.135	54	3.6	32	18- 116	
Hexachloroethane	0.250	0.122	49	3.4	33	18- 110	
Nitrobenzene	0.250	0.136	54	3.8	50	10- 211	
Pentachlorophenol	0.250	0.143	57	3.6	56	10- 140	
Pyridine	0.250	0.164	66	4.5	65	10- 148	
2,4,5-Trichlorophenol	0.250	0.146	58	2.4	22	24- 143	
2,4,6-Trichlorophenol	0.250	0.143	57	1.7	27	36- 135	· ····
Cresols (total)	0.750	0.391	52	4.2	33	25- 144	 

NOTES	(5)	
MOTES	101	

\* Values outside of QC limits

RPD 0 out of 11 outside limits Spike Recovery: 0 out of 11 outside limits

COMMENTS:

<sup>#</sup> Column to be used to flag recovery and RPD values with an asterisk

#### SW846 8270C METHOD BLANK SUMMARY

Lab Name: Severn Trent Laboratories, Inc.

BLANK WORKORDER NO.

Lab Code: QESPIT SDG Number:

Lab File ID S0522006. Lot Number: C0E130194

Date Analyzed: 05/22/00 Time Analyzed: 12:05

Matrix: SOLID Date Extracted:05/18/00

GC Column · DB5MS ID: .25 Extraction Method: 1311/3520C

Instrument ID: 71 Level: (low/med) LOW

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, LCS, LCSD, MS , MSD:

-					
		SAMPLE	LAB	DATE	TIME
	CLIENT ID	WORK ORDER #	FILE ID	ANALYZED	ANALYZED
	######################################	==========	=======================================	=========	=========
01	CHECK SAMPLE	DDF2D102 C	S0522007.	05/22/00	12:41
02	DF/S1/0133/SDC/001 A	DD6A4102	S0522008.	05/22/00	13:17
03	DF/S1/0133/SDC/001 A	DD6A411G S	S0522009.	05/22/00	13:53
04	DF/S1/0133/SDC/001 A	DD6A411H D	S0522010.	05/22/00	14:29
05	DF/S1/0133/SDC/001 B	DD6A5102	S0522011.	05/22/00	15:05
06	DF/S1/0133/SDC/001 C	DD6A6102	S0522012.	05/22/00	15:41
07	DF/S1/0133/SDC/001 D	DD6A7102	S0522013.	05/22/00	16:17
08					
09					
10					
11					I
12					
13					
14					
15					
16					
17					
18					
19					
20				·	
21					
22					
23			<del></del>		
24					
25				<del></del>	
26					
27				·	
28			·	<del></del>	
29					
30	<u> </u>				
1					

COMMENTS:	
	FORM IV

# UXB INTERNATIONAL METHOD BLANK COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix (soil/water) SOLID Lab Sample ID:C0E190000 269

Method. SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol· 200 / mL Date Received: 05/13/00 Work Order: DDF2D101 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

Moisture %:NA

QC Batch: 0140269

Client Sample Id. INTRA-LAB BLANK

CAS NO.	COMPOUND (ug/L or	r ug/kg) mg/L	0
106-46-7	1,4-Dichlorobenzene	0.050	ן טו
121-14-2	2,4-Dinitrotoluene	0.050	ן טו
118-74-1	Hexachlorobenzene	0.050	i Ui
87-68-3	Hexachlorobutadiene	0.050	Ü
67-72-1	Hexachloroethane	0.050	וֹט
98-95-3	Nitrobenzene	0.050	ָט
87-86-5	Pentachlorophenol	0.25	i vi
110-86-1	Pyridine	0.10	i ui
95-95-4	2,4,5-Trichlorophenol	0 050	U
88-06-2	2,4,6-Trichlorophenol	0.050	ן ט
1319-77-3	Cresols (total)	0.050	יט י

#### 657 42 FORM 8 SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Contract:

Lab Code: SAS No.: Case No.: SDG No.: C0E130194

Lab File ID (Standard): S0522CCC Date Analyzed: 05/22/00

Instrument ID: 71 Time Analyzed: 1127

		TC1 (DCD)		IS2 (NPT)	·	TCO (A NITT)	
		IS1(DCB) AREA #	RT #		א שמ	IS3 (ANT)	70m 44
		AREA #	RT #	AREA #	RT #	AREA #	RT #
===		245510	==== <b>=</b> =	========	======	0.55054	======
	2 HOUR STD	145518	5.03	535078	6.56	265064	9.60
	PPER LIMIT	291036	5.53	1070156	7.06	530128	10.10
1 17	OWER LIMIT	72759	4.53	267539	6.06	132532	9.10
[===		=========	=====	========	======	=======	======
	CLIENT						
SF	AMPLE NO.						
~ ===		=========	======				======
~ —   —	TRA-LAB BL	161868	5.03	630828	6.57	338411	9.60
	TRA-LAB CH	161298	5.04	635000	6.56	326610	9.60
	/S1/0133/S	171047	5.03	654778	6.56	330853	9.60
	/S1/0133/S		5.03	633807	6.57	317069	9.60
	/S1/0133/S		5.03	635621	6.57	322693	9.60
	/S1/0133/S	162741	5.03	618658	6.56	312452	9.60
	/S1/0133/S	171377	5.03	653524	6.56	328496	9.59
	/S1/0133/S	155705	5.03	593533	6.56	302566	9.59
09							
10							
11							
12							
13							
14							
15							
16					<del></del>		
17							
18							
19							<del></del>
20							
21	<del></del>						
22		l	l	l <u></u>	l		

(DCB) = 1,4-Dichlorobenzene-d4 IS1

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = + 0.50 minutes of internal standard RT RT LOWER LIMIT = - 0.50 minutes of internal standard RT

page 1 of 1

FORM VIII SV

<sup>#</sup> Column used to flag internal standard area values with an asterisk. \* Values outside of QC limits.

## FORM 8 SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Contract:

Lab Code: Case No.: SAS No.: SDG No.: C0E130194

Lab File ID (Standard): S0522CCC Date Analyzed: 05/22/00

Instrument ID: 71 Time Analyzed: 1127

	1	<del></del>					
	i	IS4 (PHN)		IS5 (CRY)	,	IS6 (PRY)	
		AREA #	RT #	AREA #	RT #	AREA #	RT #
	=========	========	======	=========	======	======================================	#
	12 HOUR STD	427212	12.90	446724	19.53	599590	
	UPPER LIMIT	854424	13.40	893448	20.03		22.89
	LOWER LIMIT	213606	12.40	223362		1199180	23.39
	==========	213000	12.40	223362	19.03	299795	22.39
	CLIENT				======	=======	=======
	SAMPLE NO.						
	SAMPLE NO.						
0.1	TATOON	========	======	=========	======	========	======
01	INTRA-LAB BL	564305	12.90	527745	19.53	648566	22.89
02	INTRA-LAB CH	529319	12.90	491727	19.52	616566	22.89
03	DF/S1/0133/S	540847	12.90	508153	19.53	598482	22.89
04	DF/S1/0133/S	MS 528651	12.90	549885	19.53	656369	22.90
05	DF/S1/0133/S	MSD 522771	12.90	519636	19.53	634789	22.90
06	DF/S1/0133/S	506789	12.89	498685	19.52	606512	22.88
07	DF/S1/0133/S	524299	12.89	498960	19.52	611014	22.88
08	DF/S1/0133/S	486337	12.89	495522	19.52	603824	
09	,,, -		12.00	773322	19.52	603824	22.88
10		i					
11							
12				<del></del>			
13							
14							
15							
16							
17		<u> </u>					
18							
19							
20						<del></del>	<del></del>
21				<del></del>		·	
22						<del></del>	i
		****					

IS4 (PHN) = Phenanthrene-d10 IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area AREA LOWER LIMIT = - 50% of internal standard area RT UPPER LIMIT = + 0.50 minutes of internal standard RT RT LOWER LIMIT = - 0.50 minutes of internal standard RT

# Column used to flag internal standard area values with an asterisk.
\* Values outside of QC limits.

page 1 of 1

FORM VIII SV

PESTICIDE SUMMARY

#### UXB INTERNATIONAL

Lab Name · Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATEL Lab Sample ID.COE130194 001

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A4103 Date Extracted:05/18/00 Dilution factor 1 Date Analyzed: 05/19/00

QC Batch · 0139460

Client Sample Id: DF/S1/0133/SDC/001 A

CAS NO.	COMPOUND (ug/L o	r ug/kg) mg/L (	2
57-74-9	Chlordane (technical)	0.0050	<u>U</u>
72-20-8	Endrin	0.00050	<u>  U</u>
76-44-8	Heptachlor	0.00050	<u>U</u>
1024-57-3	Heptachlor epoxide	0.00050	<u>U</u>
58-89-9	Lindane	0.00050	<u>  U</u>
72-43-5	Methoxychlor	0.0010	<u>  U</u>
8001-35-2	Toxaphene	0.020	<u>  U</u>

## UXB INTERNATIONAL MATRIX SPIKE COMPOUNDS

Lab Name:Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATER Lab Sample ID:COE130194 001

Method. SW846 8081A unsb3/co

Pesticides (8081A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A411A Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/19/00

QC Batch: 0139460

Client Sample Id: DF/S1/0133/SDC/001 A

CAS NO	COMPOUND (ug/L	or ug/kg) mg/L Q
72-20-8	Endrin	0.00212
76-44-8	Heptachlor	0 00236
1024-57-3	Heptachlor epoxide	0.00201
58-89-9	Lindane	0.00172
72-43-5	Methoxychlor	0.00289

### UXB INTERNATIONAL MATRIX SPIKE DUPLICATE COMPOUNDS

Lab Name.Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WARL

Method. SW846 8081A

Lons 163 160

Pesticides (8081A)

Sample WT/Vol: 100 / mL Work Order: DD6A411C Dilution factor: 1

Date Received: 05/13/00 Date Extracted:05/18/00

Date Analyzed: 05/19/00

Lab Sample ID:COE130194 001

QC Batch: 0139460

Client Sample Id: DF/S1/0133/SDC/001 A

CAS NO.	COMPOUND (ug/L	or ug/kg) mg/L Q
72-20-8	Endrin	0.00213
76-44-8	Heptachlor	0.00238
1024-57-3	Heptachlor epoxide	0.00199
58-89-9	Lindane	0.00172
72-43-5	Methoxychlor	0.00291

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATEL Lab Sample ID:COE130194 002

Method: SW846 8081A LMS/23/00

Pesticides (8081A)

Sample WT/Vol. 100 / mL Date Received: 05/13/00 Work Order: DD6A5103 Date Extracted.05/18/00 Dulution factor: 1 Date Analyzed: 05/19/00

QC Batch: 0139460

Client Sample Id: DF/S1/0133/SDC/001 B

CAS NO.	COMPOUND (ug/	Lor ug/kg) mg/L	Q
57-74-9	Chlordane (technical)	0.0050	<u>  U</u>
72-20-8	Endrin	0.00050	U
76-44-8	Heptachlor	0.00050	<u>  U</u>
1024-57-3	Heptachlor epoxide	0.00050	<u></u>
58-89-9	Lindane	0.00050	<u>u</u>
72-43-5	Methoxychlor	0.0010	<u>  "</u>
8001-35-2	Toxaphene	0.020	<u> </u>

#### UXB INTERNATIONAL

Lab Name · Severn Trent Laboratories, Inc SDG Number:

Matrix: (soil/water) SOLID VARL Lab Sample ID:COE130194 003

Method: SW846 8081A

Pesticides (8081A) Um 56860

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A6103 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed · 05/19/00

QC Batch: 0139460

Client Sample Id: DF/S1/0133/SDC/001 C

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L Q	
57-74-9	Chlordane (technical)	0.0050	<u>U</u>
72-20-8	Endrin	0.00050	U
76-44-8	Heptachlor	0.00050	U
1024-57-3	Heptachlor epoxide	0.00050	ַ
58-89-9	Lindane	0.00050	U
72-43-5	Methoxychlor	0.0010	U
8001-35-2	Toxaphene	0.020	<u>U</u>

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix (soil/water) SOLID WATEL Lab Sample ID.COE130194 004

Method: SW846 8081A (m 5)3 0

Pesticides (8081A)

Sample WT/Vol: 100 / mL Date Received 05/13/00 Work Order: DD6A7103 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed 05/19/00

QC Batch: 0139460

Client Sample Id: DF/S1/0133/SDC/001 D

CAS NO.	COMPOUND (ug/L c	or ug/kg) mg/L (	
57-74-9	Chlordane (technical)	0.0050	U
72-20-8	Endrin	0.00050	<u>U</u>
76-44-8	Heptachlor	0.00050	U
1024-57-3	Heptachlor epoxide	0.00050	<u> </u>
58-89-9	Lindane	0.00050	ַ ַ ַ ַ ַ ַ
72-43-5	Methoxychlor	0.0010	<u>U</u>
8001-35-2	Toxaphene	0.020	<u> </u>

## UXB INTERNATIONAL CHECK SAMPLE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATCH Lab Sample ID:COE180000 460

Method: SW846 8081A wsp33 bo

Pesticides (8081A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DDE2E102 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed 05/19/00

QC Batch: 0139460

Client Sample Id: CHECK SAMPLE

CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L	Q
72-20-8	Endrin	0.00219	
76-44-8	Heptachlor	0.00212	11
1024-57-3	Heptachlor epoxide	0.00219	
58-89-9	Lindane	0.00189	
72-43-5	Methoxychlor	0.00277	

SW846 8081A SURROGATE RECOVERY

657 52

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

QESSDG:

Lot #: C0E130194

ī	CLIENT ID.	SRG01	SRG02	TOT OUT
i		=======	======	======
וֹנח	DF/S1/0133/SDC/001 A	92	80	00
	DF/S1/0133/SDC/001 B	93	82	00
	DF/S1/0133/SDC/001 C	93	82	00
	DF/S1/0133/SDC/001 D	97	80	00
	METHOD BLK. DDE2E101	95	58	00
		97	91	00
	LCS DDE2E102	1 95	70	00
07	DF/S1/0133/SDC/001 A D	<del>!</del>		! !
081	DF/S1/0133/SDC/001 A S	97	_69	00

SURROGAT	res	QC LIMITS
SRG01	= Decachlorobiphenyl	( 10-147)
SRG02	<pre>= Tetrachloro-m-xylene</pre>	( 39-130)

- # Column to be used to flag recovery values
- \* Values outside of required QC Limits
- D System monitoring Compound diluted out

FORM II

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Lot #: C0E180000

WO #: DDE2E102 BATCH: 0139460

     COMPOUND	SPIKE ADDED (mg/L )	SAMPLE CONCENT. (mg/L )	% REC	QC LIMITS REC	    QUAL
Lindane	0.00250	0.00189	75	49- 137	======== 
Heptachlor	0.00250	0.00212	85	57- 124	
Heptachlor epoxide	0.00250	0.00219	88	53 - 135	
Endrin	0.00250	0.00219	88	46- 137	
Methoxychlor	0.00250	0.00277	111	12- 154	

NOTES (S):			
	 <del></del>	 	 <del></del>

* Val	lues	outside	of	QC	limits					
Spike	e Red	covery:		0	out of	5	outside	limits		
COMMI	ENTS	:								
_					-				 	

SW846 8081A MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY 54

657

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID. DF/S1/0133/SDC/001 A

Level: (low/med) LOW

Lot #: C0B130194

WO #: DD6A411A BATCH: 0139460

COMPOUND	SPIKE ADDED (mg/L )	SAMPLE CONCENT. (mg/L)	MS CONCENT. (mg/L)	MS % REC	LIMITS REC	QUAL
	========	=======	========	=====	========	=======
Lindane	0.00250	ND	0.00172	69	30- 148	<u> </u>
Heptachlor	0.00250	ND	0.00236	94_	25- 135	
Heptachlor epoxide	0.00250	ND	0.00201	80	38 - 138	1
Endrin	0.00250	ND	0.00212	85	28- 148	]
Methoxychlor	0.00250	ND	0.00289	116	13 - 154	l

NOTES(S):	

# Column to be used to flag recovery and RPD values with an asterisk \* Values outside of QC limits

RPD:	0	out	of		0	outs	ide	lin	nits	
Spike	Recov	ery:		0	out	of		5	outside	limits

COMMENTS:		
	 ·····	

### SW846 8081A MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A Level: (low/med) LOW

Lot #: C0E130194 WO #: DD6A411C BATCH: 0139460

	SPIKE ADDED	MSD CONCENT.	MSD %	ક	~	LIMITS	
COMPOUND	(mg/L)	(mg/L)	REC	RPD	RPD	REC	QUAL
=======================================	=======		25000	======	====	=======	=======
Lindane	0.00250_	0.00172	69	0.40 _	22	30- 148	
Heptachlor	0.00250	0.00238	95	1.0	32	<b>25-</b> 135	1
Heptachlor epoxide	0.00250	0.00199	80	0.77	31	38 - 138	i
Endrin	0.00250	0.00213	85	0.29	40	28- 148	
Methoxychlor	0.00250	0.00291	117	0.78	29	13- 154	

	to be used to outside of (	co flag recovery and RPD values with an asterisk .
D:	0 out of	5 outside limits
ike Rec	covery:	<u> </u>
OMMENTS:	:	

NOTES(S):

ntract:

Lab Name: STL-PITTSBURGH Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: C0E130194

Lab Sample ID: DDE2E101 Lab File ID: C-A2372

Matrix (soil/water) WATER Extraction: (SepF/Cont/Sonc) SW1311

Sulfur Cleanup (Y/N) N Date Extracted: 05/18/00

Date Analyzed (1): 05/19/00 Date Analyzed (2): 05/19/00

Time Analyzed (1): 1318 Time Analyzed (2): 1318

GC Column (1): RTX-CLP ID: 0.53 (mm) GC Column (2): RTX-CLP2 ID: 0.53 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	, <del></del>				
i	EPA	LAB		DATE	DATE
	SAMPLE NO.	SAMPLE	ID	ANALYZED 1	ANALYZED 2
		=======		05/00/00	05/00/00
01	DF/S1/0133/S	DD6A4103		05/19/00	05/19/00
02	DF/S1/0133/S	DD6A411A		05/19/00	05/19/00
03	DF/S1/0133/S			05/19/00	05/19/00
04	DF/S1/0133/S			05/19/00	05/19/00
05	DF/S1/0133/S			05/19/00	05/19/00
06	DF/S1/0133/S	DD6A7103		05/19/00	05/19/00
07	LCS	DDE2E102		05/19/00	05/19/00
08					
09					
	<b></b>				
10					
11					
12					
13					
14		<del></del>			
			······································		
15					
16					
17	]				
18					
19					
20	<u> </u>				
21					
22	i				
23					
24					
	\ <del></del>	l <del></del>	<del></del>	l	
25		]		ĺ	
26		1			<u></u>

COMMENTS:	

page 1 of 1

FORM IV PEST

OLM03.0

# UXB INTERNATIONAL METHOD BLANK COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATER Lab Sample ID:COE180000 460

Method: SW846 8081A m5/23/00

Pesticides (8081A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DDE2E101 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/19/00

QC Batch: 0139460

Client Sample Id: INTRA-LAB BLANK

CAS NO.	COMPOUND (ug/L o:	r ug/kg) mg/L 💢	<u> </u>
57-74-9	Chlordane (technical)	10.0050	<u>U</u>
72-20-8	Endrin	0.00050	U
76-44-8	Heptachlor	0.00050	U
1024-57-3	Heptachlor epoxide	0.00050	
58-89-9	Lindane	0.00050	U
72-43-5	Methoxychlor	0.0010	ַ
8001-35-2	Toxaphene	0.020	Ü

HERBICIDE SUMMARY

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATER Lab Sample ID:COE130194 001

Method: SW846 8151A um 5/22/00

Herbicides (8151A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A4104 Date Extracted: 05/18/00 Dilution factor: 1 Date Analyzed: 05/20/00

Moisture %:10

QC Batch: 0139489

Client Sample Id: DF/S1/0133/SDC/001 A

	CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L	Q
Ī	94-75-7	2,4-D	0.040	ַן ַ
Ì	93-72 <b>-</b> 1	2,4,5-TP (Silvex	0.010	U

### UXB INTERNATIONAL MATRIX SPIKE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATER

Lab Sample ID:COE130194 001

Method: SW846 8151A (m Skalod Herbicides (8151A)

Sample WT/Vol: 100 / mL Work Order: DD6A411D Dilution factor: 1

Date Received: 05/13/00 Date Extracted: 05/18/00

Date Analyzed: 05/20/00

Moisture %:10

OC Batch: 0139489

Client Sample Id: DF/S1/0133/SDC/001 A

CAS NO	COMPOUND	(ug/L or ug/kg) mg/L Q
94-75-7	2,4-D	0.133
93-72-1	2,4,5-TP (Silvex)	0.0322

## UXB INTERNATIONAL MATRIX SPIKE DUPLICATE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATEL Lab Sample ID:COE130194 001

Method: SW846 8151A un 5/22/00

Herbicides (8151A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A411E Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/20/00

Moisture %:10

QC Batch: 0139489

Client Sample Id: DF/S1/0133/SDC/001 A

CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L Q	
94-75-7	2,4-D	0.144	
93-72-1	2,4,5-TP (Silvex)	0.0339	l

#### UXB INTERNATIONAL

657 62

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:COE130194 002

Method: SW846 8151A WATER LM 5/22/07

Herbicides (8151A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A5104 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/20/00

Moisture %:8.6

OC Batch: 0139489

Client Sample Id: DF/S1/0133/SDC/001 B

	CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L (	2
Ī	94-75 <b>-7</b>	2,4-D	0.040	<u>U</u>
į	93-72-1	2,4,5-TP (Silvex)	0.010	lt

#### UXB INTERNATIONAL

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) COLID WATER Lab Sample ID:COE130194 003

Method: SW846 8151A unspales

Herbicides (8151A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A6104 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/20/00

Moisture %:10

QC Batch: 0139489

Client Sample Id: DF/S1/0133/SDC/001 C

_	CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L	Q
Ī	94-75-7	2,4-D_	0.040	ט
	93-72-1	2,4,5-TP (Silvex)	0.010	Ü

#### UXB INTERNATIONAL

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Method: SW846 8151A WATER UM 1

Herbicides (8151A) 5/22/00

Date Received: 05/13/00 Sample WT/Vol: 100 / mL Date Extracted:05/18/00 Work Order: DD6A7104 Date Analyzed: 05/20/00 Dilution factor: 1

Moisture %:11

QC Batch: 0139489 Client Sample Id: DF/S1/0133/SDC/001 D

CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L	Q
94-75-7	2,4-D	0.040	ן <u>ט</u>
93-72-1	2,4,5-TP (Silve:	x) 0.010	<u>  U</u>

#### UXB INTERNATIONAL CHECK SAMPLE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOUTH WATER

Method: SW846 8151A (ms 22)

Sample WT/Vol: 100 / mL Work Order: DDE4W102 Dilution factor: 1

Date Received: 05/13/00 Date Extracted: 05/18/00 Date Analyzed: 05/20/00

Lab Sample ID:COE180000 489

Moisture %:NA

QC Batch: 0139489

Client Sample Id: CHECK SAMPLE

	CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L (	2
T_	94-75-7	2,4-D	0.133	lI
i T	93 <b>-</b> 72-1	2,4,5-TP (Silvex)	0.0323	11

#### SW846 8151A SURROGATE RECOVERY 8 B 657

Name ! Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

QESSDG:

Lot #: C0E130194

CLIENT ID.	SRG01	TOT OUT
		======
01 DF/S1/0133/SDC/001 A	51	00
02 DF/S1/0133/SDC/001 B	57	00
03 DF/S1/0133/SDC/001 C	56	00
04 DF/S1/0133/SDC/001 D	66	00
05 METHOD BLK. DDE4W101	78	00
06 LCS DDE4W102	89	00
07 DF/S1/0133/SDC/001 A D	86	00
08 DF/S1/0133/SDC/001 A S	82	00

SURROGATES = DCAA SRG01

QC LIMITS (42-125)

- # Column to be used to flag recovery values
- \* Values outside of required QC Limits
- D System monitoring Compound diluted out

FORM II

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Lot #: C0E180000

WO #: DDE4W102 BATCH: 0139489

COMPOUND	SPIKE ADDED (mg/L )	SAMPLE CONCENT. (mg/L )	% REC	QC LIMITS REC	QUAL
			====	=======================================	=========
2,4-D	0.160	0.133	83	28- 136	<u> </u>
2,4,5-TP (Silvex)	0.0400	0.0323	81	50- 128	ii

Values outside	of QC	limits					
pike Recovery:	0	out of	2	outside	limits		
COMMENTS:							

NOTES (S):

657 68 SW846 8151A MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A

Lot #: C0E130194

WO #: DD6A411D BATCH: 0139489

	SPIKE ADDED	SAMPLE CONCENT.	MS CONCENT.	MS %	LIMITS REC	OUAL
COMPOUND	(mg/L )	(mg/L )  ========	(mg/L )	REC		QUAL [
2,4-D	0.160	ND	0.133	83	35- 133	
2,4,5-TP (Silvex)	0.0400	ND	0.0322	80	50- 131	

NOTES(S):				
	· · · · · · · · · · ·	 	 	·

# Column to be used to flag recovery and RPD values with an asterisk \* Values outside of QC limits RPD: 0 out of 0 outside limits Spike Recovery: 0 out of 2 outside limits COMMENTS:

## SW846 8151A MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A

Lot #: C0E130194

WO #: DD6A411E BATCH: 0139489

	SPIKE	MSD	MSD				
	ADDED	CONCENT.	ક	ક	QC :	LIMITS	ĺ
COMPOUND	(mg/L )	(mg/L )	REC	RPD	RPD	REC	QUAL
		=======	=====		====		=======================================
2,4-D	0.160	0.144	90	8.4	20	35- 133	l
2,4,5-TP (Silv	(ex) 0.0400	0.0339	85	5.2	20	50- 131	

		flag recovery and RPD values with an asterisk
Values out	tside of QC	limits
RPD:0	out of _	2 outside limits
Spike Recove	ery: <u>0</u>	out of 2 outside limits
COMMENTS:		
	<del></del>	

NOTES(S):

657	70					SUMMARY
001	1 0	SW846	8151A	METHOD	BLANK	SUMMARY

BLANK WORKORDER NO.

Lab Name: Severn Trent Laboratories, Inc.

Lab Code: QESPIT

SDG Number:

Lab File ID: a-a20536.

Lot Number: COE130194

Matrix: SOLID WATER

Extraction Method:

colcolo m

Date Analyzed(1): 05/20/00

Date Analyzed(2): N/A

Date Extracted: 05/18/00

Time Analyzed(1): 14:27

Time Analyzed(2): N/A

Instrument ID(1): A/B

Instrument ID(2): N/A

GC Column(1): DB5/DB1701 ID: 053

GC Column(2): N/A ID:

N/A

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS, AND MSD:

1		SAMPLE	DATE	DATE	ĺ
	CLIENT ID.	WORK ORDER #	ANALYZED(1)	ANALYZED(2)	
	=======================================		= =====================================	==========	
01	CHECK SAMPLE	DDE4W102 C	05/20/00	N/A	l
	DF/S1/0133/SDC/001 A	_DD6A4104	05/20/00	_N/A	l
	DF/S1/0133/SDC/001 A	DD6A411D S	05/20/00	N/A	Ì
	DF/S1/0133/SDC/001 A	DD6A411E D	05/20/00	N/A	
	DF/S1/0133/SDC/001 B	DD6A5104	05/20/00	N/A	١
	DF/S1/0133/SDC/001 C	DD6A6104	05/20/00	N/A	
07	DF/S1/0133/SDC/001 D	DD6A7104	05/20/00	N/A	
08					
09					l
10					ļ
11					1
12					
13	]				
14					١
15		1			
16	·	1			ļ
17					1
18			1		
19	; <del></del>				Ĺ
20			_		İ
~ ~ ~	I .	1	_		•

COMMENTS:	
	FORM TV

#### UXB INTERNATIONAL METHOD BLANK COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATER Lab Sample ID:COE180000 489

Method: SW846 8151A

Herbicides (8151A)

shalou

Sample WT/Vol: 100 / mL Work Order: DDE4W101 Dilution factor: 1

Date Received: 05/13/00 Date Extracted:05/18/00 Date Analyzed: 05/20/00

Moisture %:NA

QC Batch: 0139489

Client Sample Id: INTRA-LAB BLANK

'CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L (	2
94-75-7	2,4-D	0.040	<u></u> U
93-72-1	2,4,5-TP (Silve:	x) 0.010	ا <u>ت</u> ا

## METALS SUMMARY

## Metals Data Reporting Form

Sample Results

Lab Sample ID:

DD6A4 Client ID: DF/S1/0133/SDC/001 A

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 10.39

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	189.04	0.29	1.1	7.7		1	ICPST	5/17/00	9:15

657 74

### STL-Pittsburgh

### Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A5 Client ID: DF/S1/0133/SDC/001 B

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 8.64

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	189.04	0.28	1.1	7.2		1	ICPST	5/17/00	9:32

Comments: Lot #: C0E130194 Sample #: 2

## Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A6 Client ID: DF/S1/0133/SDC/001 C

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 10.27

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	189.04	0.29	1.1	8.1			ICPST	5/17/00	9:36

## Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A7 Client ID: DF/S1/0133/SDC/001 D

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 10.81

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	189.04	0.29	1.1	7.7		1	ICPST	5/17/00	9:48

Initial Calil	bration B	lank R	esults			_						
Instrument:	ICP	ST	-				Units:		ug/L			
Chart Numb	ber: <u>T00</u> :	517A.A	RC_									
Standard So	ource:	<u></u>		_			Standar	d ID:		·		
			ICB1 5/17/00 8:52 AM									
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	Q	Found	Q	Found	Q
Arsenic	189.042	10	2.6	U		-						

## Metals Data Reporting Form

Continuing Calibration Blank Results Units: ug/L **ICPST** Instrument: Chart Number: T00517A.ARC Standard ID: Standard Source: CCB1 CCB2 5/17/00 5/17/00 9:44 AM 9:57 AM WL/ Report Found Found Found Limit Found Q Element Mass **Found** 10 2.6 U 2.6 U 189.042 Arsenic

U Result is less than the MDL

B Result is between MDL and RL

### Metals Data Reporting Form

Preparation Blank Results

Lab Sample ID: DD8AMB

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	189.042	0.26	1.0	0.26	U	1	ICPST	5/17/00	9:07

#### Metals Data Reporting Form

Matrix Spike Sample Results

Spike Sample ID: DD6A4S

Original Sample ID: DD6A4 Client ID: DF/S1/0133/SDC/001 AS

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

 Weight:
 1.0
 Volume:
 100
 Percent Moisture:
 10.39

Element	WL/ Mass	OS Conc	Q	MS Conc	Q	Spike Level	% Rec	OS DF	MS DF	Instr	OS Anal Date	OS Anal Time	MS Anal Date	MS Anal Time
Arsenic	189.0	7.7		216		223.19	93.1	1	1	ICPST	5/17/00	9:15	5/17/00	9:24

Comments: Lot #: C0E130194 Sample #: 1

Version 3.63.3

U Result is less than the MDL

Form 5A Equivalent

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

\* Duplicate analysis RPD was not within limits

STL Pittsburgh 81

### Metals Data Reporting Form

Matrix Spike Duplicate RPD Report

Matrix Spike Duplicate Sample ID: DD6A4D

Matrix Spike Sample ID: DD6A4S Client ID: DF/S1/0133/SDC/001 AD

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 10.39

Element	WL/ Mass	MS Conc	o	MSD Conc	Q	RPD	MS DF	MSD DF	Instr	MS Anal Date	MS Anal Time	MSD Anal Date	MSD Anal Time
Arsenic	189.042	216		217		0.6%	1	1	ICPST	5/17/00	9:24	5/17/00	9:28

Comments: Lot #: C0E130194 Sample #: 1

Version 3.63.3

U Result is less than the MDL

Form 6 Equivalent

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

<sup>\*</sup> Duplicate analysis RPD was not within limits

## Metals Data Reporting Form

Laboratory Control Sample Results

Lab Sample ID: DD8AMC

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: NA

Element	WL/ Mass	Spike Level	Conc	Percent Recovery	Q	Range	DF	Instr	Anal Date	Anal Time
Arsenic	189.042	200	198	98.7		80-120	1	ICPST	5/17/00	9:11

B Result is between MDL and RL

## TCLP METALS SUMMARY

## Metals Data Reporting Form

Sample Results

Lab Sample ID:

DD6A4T

Client ID:

DF/S1/0133/SDC/001 A

Matrix:

Water

Units: mg/L Prep Date: 5/18/00 **Prep Batch:** 0139097

Weight: NA Volume:

100

Percent Moisture:

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.000045	0.00020	0.000045	U	1	CVAA	5/18/00	11:03

## Metals Data Reporting Form

Prep Date:

Sample Results

Lab Sample ID: DD6A5T

Client ID: DF/S1/0133/SDC/001 B

Matrix: Water Units: mg/L

5/18/00 Prep Batch: 0139097

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.000045	0.00020	0.000045	U	1	CVAA	5/18/00	11:09

### Metals Data Reporting Form

Sample Results

DD6A6T Lab Sample ID:

Client ID:

DF/S1/0133/SDC/001 C

Matrix:

Water

mg/L Units:

Prep Date: 5/18/00 **Prep Batch:** 0139097

Weight:

NA

Volume: 100

Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0 000045	0.00020	0.000045	υ	1	CVAA	5/18/00	11:10

87

Sample Results

### STL-Pittsburgh

### Metals Data Reporting Form

Lab Sample ID: DD6A7T Client ID: DF/S1/0133/SDC/001 D

**Prep Date:** 5/18/00 **Prep Batch:** 0139097 Matrix: Water Units: mg/L

NA Weight: NA Volume: 100 Percent Moisture:

Element	WL/ Mass	MDL	Report Limit	Conc	o	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.000045	0 00020	0.000045	บ	1	CVAA	5/18/00	11:13

89

### STL-Pittsburgh

## Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A4T Client ID: DF/S1/0133/SDC/001 A

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	193.7	0.030	0.50	0.15	В	1	ICP	5/19/00	9:14
Barium	493.41	0.00041	10.0	1.2	BE	1	ICP	5/19/00	9:14
Cadmium	228.80	0.0028	0.10	0.0028	U	1	ICP	5/19/00	9:14
Chromium	267.72	0.0038	0.50	0.0038	ַ ט	1	ICP	5/19/00	9:14
Lead	220.35	0.025	0.50	0.025	U	1	ICP	5/19/00	9:14
Selenium	196.03	0.067	0.25	0.067	υ	1	ICP	5/19/00	15:29
Silver	328.07	0.0031	0.50	0.0031	ับ	1	ICP	5/19/00	9:14

Comments: C0E130194001

## Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A5T Client ID: DF/S1/0133/SDC/001 B

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	193.7	0.030	0.50	0.21	В	1	ICP	5/19/00	9:27
Barium	493.41	0.00041	10.0	1.2	BE	1	ICP	5/19/00	9:27
Cadmium	228.80	0.0028	0.10	0.0028	U	1	ICP	5/19/00	9:27
Chromium	267.72	0.0038	0.50	0.0038	U	1	ICP	5/19/00	9:27
Lead	220.35	0.025	0.50	0.025	U	1	ICP	5/19/00	9:27
Selenium	196.03	0.067	0.25	0.067	U	l	ICP	5/19/00	15:42
Silver	328.07	0.0031	0.50	0.0031	U	1	ICP	5/19/00	9:27

Comments: <u>C0E130194002</u>

U Result is less than the MDL

B Result is between MDL and RL

## Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A6T Client ID: DF/S1/0133/SDC/001 C

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	193.7	0.030	0.50	0.20	В	1	ICP	5/19/00	9:36
Barium	493.41	0.00041	10.0	1.3	BE	1	ICP	5/19/00	9:36
Cadmium	228.80	0.0028	0.10	0.0028	U	1	ICP	5/19/00	9:36
Chromium	267.72	0.0038	0.50	0.0038	บ	1	ICP	5/19/00	9:36
Lead	220.35	0.025	0.50	0.025	υ.	1	ICP	5/19/00	9:36
Selenium	196.03	0.067	0.25	0.067	บ	1	ICP	5/19/00	15:51
Silver	328.07	0.0031	0.50	0.0031	U	1	ICP	5/19/00	9:36

Comments: C0E130194003

### Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A7T Client ID: DF/S1/0133/SDC/001 D

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	193.7	0.030	0.50	0.15	В	1	ICP	5/19/00	9:39
Barium	493,41	0.00041	10.0	1.2	BE	1	ICP	5/19/00	9:39
Cadmium	228.80	0.0028	0.10	0.0028	ับ	1	ICP	5/19/00	9:39
Chromium	267.72	0.0038	0.50	0.0038	U	1	ICP	5/19/00	9:39
Lead	220.35	0.025	0.50	0.025	U	1	ICP	5/19/00	9:39
Selenium	196.03	0.067	0.25	0.067	ַ ט	1	ICP	5/19/00	15.55
Silver	328.07	0.0031	0.50	0 0031	บ	1_	ICP	5/19/00	9:39

Comments: <u>C0E130194004</u>

Initial Cali	bration B	lank R	esults									
Instrument	:CV	AA	_				Units:		ug/L	_	•	-
Chart Num	ber: <u>0518</u>	HGA.P	RN_									
Standard S	ource:		<u></u> -		<del></del>		Standar	d ID:		<del></del>		
			ICB1 5/18/0 10:51 A	0								
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	Q	Found	Q	Found	Q
Mercury	253.7	0.2	0.0	В								

# Metals Data Reporting Form

Initial Calibrati	on Blank Results		
Instrument:	ICP	Units: ug/L	
Chart Number:	J00519A.ARC		
Standard Source	<b>.</b> •	Standard ID:	

			ICB1 5/19/00 8:53 AM									
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	Q	Found	Q	Found	Q
Arsenic	193.696	500	30.3	U								
Barium	493.409	10000	0.4	U								
Cadmium	228.802	100	2.8	U								
Chromium	267.716	500	3.8	U								
Lead	220.353	500	24.6	U								
Silver	328.068	500	3.1	U							<u> </u>	

94

U Result is less than the MDL

B Result is between MDL and RL

Initial Calil	oration B	lank R	esults									
Instrument:	IC	P	<u>-</u>				Units:		ug/L			
Chart Numb	oer:J005	19B.AI	RC_									
Standard So	urce:						Standar	d ID:				_
			ICB1 5/19/00 3:08 PN									
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	Q	Found	Q	Found	Q
Selenium	196.026	250	67.4	U								

Continuin	g Calibra	ation B	lank Res	ults								
Instrument	:	VAA	_				Units:		ug/L			
Chart Num	ber: <u>05</u>	18HGA.	PRN									
Standard S	ource:					<b>-</b>	Standa	ard ID	):			
			CCB: 5/18/0 10:55 A	00	CCB2 5/18/00 11:18 A	0						
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	Q	Found	Q	Found	Q
Mercury	253.7	0.2	0.0	T I	-0.1	В						

U Result is less than the MDL

B Result is between MDL and RL

Continuing Calibration Blank Results	
Instrument: ICP	Units: ug/L
Chart Number: J00519A.ARC	
Standard Source:	Standard ID:

			CCB1 5/19/00 9:33 AN		CCB2 5/19/00 9:46 AN	)						
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	Q	Found	Q	Found	Q
Arsenic	193.696	500	32.7	В	30.3	υ						
Barium	493.409	10000	0.8	В	0.6	В						
Cadmium	228.802	100	2.8	U	2.8	U						
Chromium	267.716	500	3.8	U	3.8	U						
Lead	220.353	500	24.6	U	24.6	U					1	
Silver	328.068	500	3.1	U	3.1	U						

### 98 657

# STL-Pittsburgh

# Metals Data Reporting Form

Continuing	g Calibra	tion B	lank Resi	ılts								
Instrument:	I	CP					Units:		ug/L			
Chart Numl	ber: <u>J0</u>	0519B.A	ARC									
Standard So	ource:						Standa	ard ID	):			
	Standard Source:			) M	CCB2 5/19/00 4:01 PM							
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	_Q_	Found	Q	Found	Q
Selenium	196.026	250	67.4	U	67.4	U						

Form 3 Equivalent B Result is between MDL and RL

## Metals Data Reporting Form

Preparation Blank Results

Lab Sample ID: DDA9GBT

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139097

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.000045	0.00020	0.000063	В	1	CVAA	5/18/00	11:01

## Metals Data Reporting Form

Preparation Blank Results

Lab Sample ID:

DDC2GBT

Matrix: Water

Units: mg/L

5/18/00 Prep Date:

Prep Batch: 0139097

Weight:

NA

Volume:

100

Percent Moisture:

Element	WL/ Mass	MDL	Report Limit	Сопс	Q	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.000045	0.00020	-0.000090	В	1	CVAA	5/18/00	10:57

### Metals Data Reporting Form

Preparation Blank Results

Lab Sample ID: DDA9GBT

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	193.696	0.030	0.50	0.18	В	1	ICP	5/19/00	9:08
Barium	493.409	0.00041	10.0	0.037	В	1	ICP	5/19/00	9:08
Cadmium	228.802	0.0028	0 10	0.0028	Ų	1	ICP	5/19/00	9:08
Chromium	267.716	0.0038	0 50	0.0038	U	1	ICP	5/19/00	9:08
Lead	220.353	0.025	0 50	0.025	U	1	ICP	5/19/00	9:08
Selenium	196.026	0.067	0.25	0.067	υ	1	ICP	5/19/00	15:23
Silver	328.068	0.0031	0 50	0.0031	U	1	ICP	5/19/00	9:08

### Metals Data Reporting Form

Preparation Blank Results

Lab Sample ID: DDC39BT

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	193.696	0.030	0.50	0.030	U	1	ICP	5/19/00	9:05
Barium	493.409	0.00041	10.0	0.00041	บ	1	ICP	5/19/00	9:05
Cadmium	228.802	0.0028	0.10	0.0028	U	1	ICP	5/19/00	9:05
Chromium	267.716	0.0038	0.50	0.0038	บ	1	ICP	5/19/00	9:05
Lead	220.353	0.025	0.50	0.025	ט	1	ICP	5/19/00	9:05
Selenium	196.026	0.067	0.25	0 067	บ	1	ICP	5/19/00	15:20
Silver	328.068	0.0031	0.50	0.0031	ט	1	ICP	5/19/00	9:05

#### Metals Data Reporting Form

Matrix Spike Sample Results

Spike Sample ID: DD6A4ST

Original Sample ID: DD6A4T Client ID: DF/S1/0133/SDC/001 AS

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139097

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	OS Conc	Q	MS Cone	Q	Spike Level	% Rec	OS DF	MS DF	Instr	OS Anal Date	OS Anal Time	MS Anal Date	MS Anal Time
Mercury	253.7	0.000045	υ	0.0053		0.005		1	1	CVAA	5/18/00	11:03	5/18/00	11:04

Comments: \_\_\_\_ Version 3.63.3

U Result is less than the MDL

Form 5A Equivalent

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

### Metals Data Reporting Form

Matrix Spike Duplicate Sample Results

Spike Sample ID: DD6A4DT

Original Sample ID: DD6A4T Client ID: DF/S1/0133/SDC/001 AD

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139097

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	OS Conc	Q	MSD Conc	Q	Spike Level	% Rec	OS DF	MSD DF	Instr	OS Anal Date	OS Anal Time	MSD Anal Date	MSD Anal Time
Mercury	253.7	0.000045	U	0.0052		0.005	104.8	1	I	CVAA	5/18/00	11:03	5/18/00	11:07

Comments:

Version 3.63.3

U Result is less than the MDL

Form 5A Equivalent

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

\* Duplicate analysis RPD was not within limits

## 657 105

#### STL-Pittsburgh

#### Metals Data Reporting Form

### Matrix Spike Sample Results

Spike Sample ID:

DD6A4ST

Original Sample ID:

DD6A4T

Client ID:

DF/S1/0133/SDC/001 AS

Matrix:

Water

Units:

mg/L

Prep Date:

5/18/00

Prep Batch:

0139114

Weight:

NA

Volume:

50

Percent Moisture:

NA

Element	WL/ Mass	OS Conc	Q	MS Conc	Q	Spike Level	% Rec	OS DF	MS DF	Instr	OS Anal Date	OS Anal Time	MS Anal Date	MS Anal Time
Arsenic	193.7	0.15	В	4.9		5	94.8	1	1	ICP	5/19/00	9:14	5/19/00	9:20
Barium	493.4	1.2	В	43.3		50	84.1	1	1	ICP	5/19/00	9:14	5/19/00	9:20
Cadmium	228.8	0.0028	U	0.94		1	94.4	1	1	ICP	5/19/00	9:14	5/19/00	9:20
Chromium	267.7	0.0038	ប	4.3		5	85 3	1	1	ICP	5/19/00	9:14	5/19/00	9:20
Lead	220.4	0.025	U	4.4		5	88.7	1	1	ICP	5/19/00	9:14	5/19/00	9:20
Selenium	196.0	0.067	U	1.1		1	104.8	1	1	ICP	5/19/00	15:29	5/19/00	15:36
Silver	328.1	0.0031	U	0.90		1	90.4	1	1	ICP	5/19/00	9:14	5/19/00	9:20

Comments:

Version 3.63.3

\* Duplicate analysis RPD was not within limits

U Result is less than the MDL

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

## Metals Data Reporting Form

Matrix Spike Duplicate Sample Results

Spike Sample ID: DD6A4DT

Original Sample ID: DD6A4T Client ID: DF/S1/0133/SDC/001 AD

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	OS Cone	Q	MSD Conc	Q	Spike Level	% Rec	OS DF	MSD DF	Instr	OS Anal Date	OS Anal Time	MSD Anal Date	MSD Anal Time
Arsenic	193.7	0.15	В	5 0		5	96.2	1	1	ICP	5/19/00	9:14	5/19/00	9:24
Barium	493.4	1.2	В	43.8		50	85.2	i	1	ICP	5/19/00	9:14	5/19/00	9:24
Cadmium	228.8	0.0028	U	0.96		1	95.5	1	1	ICP	5/19/00	9:14	5/19/00	9:24
Chromium	267.7	0.0038		4.3		5	86.3	1	1	ICP	5/19/00	9:14	5/19/00	9:24
Lead	220.4	0.025		4.5		5	89.8	1	1	ICP	5/19/00	9:14	5/19/00	9.24
Selenium	196.0	0.067	υ	1.1		1	105.2	1	1	ICP	5/19/00	15:29	5/19/00	15:39
Silver	328.1	0.0031	υ	0.91		1	91.2	1	1	ICP	5/19/00	9:14	5/19/00	9:24

Comments:

Version 3.63.3

Form 5A Equivalent

U Result is less than the MDL

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

Duplicate analysis RPD was not within limits

### Metals Data Reporting Form

Matrix Spike Duplicate RPD Report

Matrix Spike Duplicate Sample ID: DD6A4DT

Matrix Spike Sample ID: DD6A4ST Client ID: DF/S1/0133/SDC/001 AD

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139097

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	MS Conc	Q	MSD Conc	Q	RPD	MS DF	MSD DF	Instr	MS Anal Date	MS Anal Time	MSD Anal Date	MSD Anal Time
Mercury	253.7	0.0053		0.0052		1.7%	1	1	CVAA	5/18/00	11:04	5/18/00	11:07

Comments: \_\_\_\_ Version 3.63.3

U Result is less than the MDL

Form 6 Equivalent

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

\* Duplicate analysis RPD was not within limits

# STL-Pittsburgh

# Metals Data Reporting Form

Matrix Spike Duplicate RPD Report

Matrix Spike Duplicate Sample ID: DD6A4DT

Matrix Spike Sample ID: DD6A4ST Client ID: DF/S1/0133/SDC/001 AD

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MS Conc	Q	MSD Cone	Q	RPD	MS DF	MSD DF_	Instr	MS Anal Date	MS Anal Time	MSD Anal Date	MSD Anal Time
Arsenic	193.696	4.9		5.0		1.5 %	1	1	ICP	5/19/00	9:20	5/19/00	9:24
Barium	493.409	43.3		43.8	:	1.4 %	1	1	ICP	5/19/00	9:20	5/19/00	9:24
Cadmium	228.802	0.94		0.96		1.1 %	1	1	ICP	5/19/00	9:20	5/19/00	9:24
Chromium	267.716	4.3		4.3		1.2 %	1	1	ICP	5/19/00	9:20	5/19/00	9:24
Lead	220.353	4.4		4.5		1.2 %	1	1	ICP	5/19/00	9:20	5/19/00	9:24
Selenium	196.026	1.1		11		10.2 %	ì	1	ICP	5/19/00	15:36	5/19/00	15:39
Silver	328.068	0 90		0.91		1.0 %	1	1	ICP	5/19/00	9:20	5/19/00	9:24

Comments: \_\_\_\_\_
Version 3.63.3

U Result is less than the MDL

Form 6 Equivalent

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

\* Duplicate analysis RPD was not within limits

# STL-Pittsburgh

# Metals Data Reporting Form

Laboratory Control Sample Results

Lab Sample ID: DDC2GCT

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139097

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	Spike Level	Conc	Percent Recovery	Q	Range	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.0025	0.0028	110.0		80-120	1	CVAA	5/18/00	10:58

# STL-Pittsburgh

# Metals Data Reporting Form

Laboratory Control Sample Results

Lab Sample ID: DDC39CT

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	Spike Level	Conc	Percent Recovery	Q	Range	DF	Instr	Anal Date	Anal Time
Arsenic	193 696	2.0	2.2	110.1		80-120	1	ICP	5/19/00	9:11
Barium	493 409	2.0	1.9	93.7	В	80-120	1	ICP	5/19/00	9:11
Cadmium	228.802	0.050	0.046	91.3	В	80-120	1	ICP	5/19/00	9:11
Chromium	267.716	0.20	0.19	94.7	В	80-120	1	ICP	5/19/00	9:11
Lead	220.353	0.50	0.47	94.5	В	80-120	1	ICP	5/19/00	9:11
Selenium	196.026	2.0	2.2	107.8		80-120	1	ICP	5/19/00	15:26
Silver	328.068	0.050	0 047	94.3	В	80-120	1	ICP	5/19/00	9:11

GENERAL CHEMISTRY SUMMARY

## Client Sample ID: DF/S1/0133/SDC/001 A

## General Chemistry

Lot-Sample #...: COE130194-001 Work Order #...: DD6A4

Matrix..... SOLID

Date Sampled...: 05/12/00

Date Received..: 05/13/00 **% Moisture....:** 10

PARAMETER	RESULT	RL_	UNITS	METHO	··	PREPARATION- ANALYSIS DATE	PREP BATCH #
рH	7.8 Dilution	Factor: 1	No Units MS Run #		<b>9045C</b> 0136107	05/15/00	0136242
Ignitability	INO Dilution	 Factor: 1	No Units MS Run #		SECTION 7.1 0136105	05/15/00	0136237
Percent Solids	<b>89.6</b> Dilution	1.0 Factor: 1	<b>%</b> MS Run #		<b>160.3 MOD</b> 0140053	05/19-05/20/00	0140168
Reactive Cyanide	ND Dilution	200 Factor: 1	mg/kg Ms Run#		7.3.3 0138162	05/17/00	0138394
Reactive Sulfide	ND Dilution	200 Factor: 1	mg/kg Ms Run #		7.3.4 0138161	05/17/00	0138393

Matrix....: SOLID

## UXB INTERNATIONAL

# Client Sample ID: DF/S1/0133/SDC/001 B

# General Chemistry

Lot-Sample #...: COE130194-002 Work Order #...: DD6A5

Date Sampled...: 05/12/00

Date Received..: 05/13/00

\* Moisture....: 8.6

PARAMETER	RESULT	RL,	UNITS	METHO	D	PREPARATION- ANALYSIS DATE	PREP BATCH #
рн	8.1 Dilution F	actor: 1	No Units MS Run #		<b>9045C</b> 0136107	05/15/00	0136242
Ignitability	<b>INO</b> Dilution F		No Units MS Run #.		SECTION 7.1 0136105	05/15/00	0136237
Percent Solids	91.4 Dilution F		<b>%</b> MS Run #.		<b>160.3 MOD</b> 0140053	05/19-05/20/00	0140168
Reactive Cyanide	ND :	200	mg/kg	SM046	7 2 2	05/17/00	0120204
	Dilution F		MS Run #.			03/17/00	0138394

# Client Sample ID: DF/S1/0133/SDC/001 C

## General Chemistry

Lot-Sample #...: COE130194-003 Work Order #...: DD6A6

Date Sampled...: 05/12/00

Date Received..: 05/13/00

Matrix..... SOLID

**\* Moisture....:** 10

PARAMETER	RESULT	RL	UNITS	METHO	<u>D</u>	PREPARATION- ANALYSIS DATE	PREP BATCH #
рH	8.1 Dilution	Factor: 1	No Units MS Run #	SW846	<b>9045C</b> 0136107	05/15/00	0136242
Ignitability	INO Dilution	 Factor: 1	No Units MS Run #	SW846	SECTION 7.1 0136105	05/15/00	0136237
Percent Solids		1.0 Factor: 1	% MS Run #	MCAWW	160.3 MOD 0140053	05/19-05/20/00	0140168
Reactive Cyanide	ND Dilution	200 Factor: 1	mg/kg MS Run a	SW846	7.3.3 0138162	05/17/00	0138394
Reactive Sulfide	ND Dilution	200 Factor: 1	mg/kg MS Run a	SW846	7.3.4 0138161	05/17/00	0138393

657 115

# Client Sample ID: DF/S1/0133/SDC/001 D

## General Chemistry

Lot-Sample #...: C0E130194-004 Work Order #...: DD6A7 Matrix.....: SOLID

\* Moisture....: 11

PARAMETER	RESULT	RL	UNITS	METHOD	PREPARATION- ANALYSIS DATE	PREP BATCH #
рн	7.9 Dilution	Factor: 1	No Units MS Run #	SW846 9045C : 0136107	05/15/00	0136242
Ignitability	<b>NO</b> Dilution	Factor: 1	No Units MS Run #	SW846 SECTION 7.1	05/15/00	0136237
Percent Solids	89.2 Dilution	1.0 Factor: 1	∜ MS Run #	MCAWW 160.3 MOD : 0140053	05/19-05/20/00	0140168
Reactive Cyanide	ND Dilution	200 Factor: 1		SW846 7.3.3 : 0138162	05/17/00	0138394
Reactive Sulfide	NID Dilution	200 Factor: 1	mg/kg MS Run #	SW846 7.3.4	05/17/00	0138393

# General Chemistry

Client Lot #...: C0E130194

Matrix ..... SOLID

PARAMETER	RESULT	REPORTING LIMIT	UNITS	METHOD		PREPARATION- ANALYSIS DATE	PREP BATCH #
Reactive Cyanide	ND Di	Work Order 200 lution Factor: 1	#: DDARD101 mg/kg	MB Lot-Sample SW846 7.3.3	#:	A0E170000-394 05/17/00	0138394
Reactive Sulfide	ND Di	Work Order 200 lution Factor: 1	#: DDARC101 mg/kg	MB Lot-Sample SW846 7.3.4	#:	A0E170000-393 05/17/00	0138393

NOTE (S):

Calculations are performed before rounding to avoid round-off errors in calculated results.

## LABORATORY CONTROL SAMPLE EVALUATION REPORT

## General Chemistry

Client Lot #...: C0E130194

Matrix....: SOLID

PERCENT PARAMETER

RECOVERY LIMITS

PREPARATION-

PREP

RECOVERY

METHOD

SW846 9045C

ANALYSIS DATE

BATCH #

Нq

Work Order #: DD70J101 LCS Lot-Sample#: C0E150000-242

100

(85 - 115) Dilution Factor: 1

05/15/00

0136242

NOTE (S):

Calculations are performed before rounding to avoid round-off errors in calculated results.

STL Pittsburgh

117

657 118

General Chemistry

Client Lot #...: COE130194 Work Order #...: DDAJ6-SMP Matrix.....: SOLID

DDAJ6-DUP

Date Sampled...: 04/12/00 Date Received..: 05/17/00

\* Moisture....: 20

PREPARATION-PREP RPD DUPLICATE <u>LIMI</u>T ANALYSIS DATE BATCH # PARAM RESULT RESULT UNITS RPD METHOD SD Lot-Sample #: C0E170244-030 Percent Solids (0-20) MCAWW 160.3 MOD 05/19-05/20/00 0140168 79.9 81.7 욯 2.2

Dilution Factor: 1

Prep Date....: 0140053 Analysis Date..: Prep Batch #...:

#### SAMPLE DUPLICATE EVALUATION REPORT

#### General Chemistry

Client Lot #...: COE130194 Work Order #...: DD6DM-SMP Matrix....: SOLID

Prep Date....: 0138161 Analysis Date..:

DD6DM-DUP

Prep Batch #...:

Date Sampled...: 05/11/00 Date Received..: 05/13/00

\* Moisture....: 0.0 DUPLICATE RPD PREPARATION-PREP RESULT PARAM RESULT RPD UNITS LIMIT METHOD ANALYSIS DATE BATCH # Reactive Cyanide SD Lot-Sample #: G0E130196-003 ND ND mg/kg 0 SW846 7.3.3 05/17/00 (0-20)0138394 Dilution Factor: 1 Prep Date....: 0138162 Analysis Date..: Prep Batch #...: Reactive Sulfide SD Lot-Sample #: G0E130196-003 ND ND mg/kg 200 (0-20) SW846 7.3.4 05/17/00 0138393 Dilution Factor: 1

GC/MS VOLATILE DATA

# GC/MS VOLATILE QC SUMMARY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

QESSDG:

Lot #: C0E130194

-	CLIENT ID.	SRG01	SRG02	SRG03	SRG04	TOT OUT
	=======================================	a=====	======	======	======	======
	DF/S1/0133/SDC/001 A	110	103	93	103	00 j
	DF/S1/0133/SDC/001 B	108	99	89	104	.00
03	DF/S1/0133/SDC/001 C	106	106	93	102	00
04	DF/S1/0133/SDC/001 D	119	96	91	107	00
05	METHOD BLK. DD95L101	104	100	90	98	00
06	LCS DDA08101	103	104	103	102	00
07	DF/S1/0133/SDC/001 A D	103	103	102	102	00
08	DF/S1/0133/SDC/001 A S	102	104	102	103	00

SURROGE	<u>ates</u>	QC LIMITS
SRG01	= 1,2-Dichloroethane-d4	(77-120)
SRG02	= Toluene-d8	( 78-111)
SRG03	= 4-Bromofluorobenzene	( 80-114)
SRG04	= Dibromofluoromethane	( 78-110)

FORM II

<sup>#</sup> Column to be used to flag recovery values

<sup>\*</sup> Values outside of required QC Limits

D System monitoring Compound diluted out

657 123

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Lot #: C0E170000

WO #: DDA08101 BATCH: 0138262

	SPIKE	SAMPLE		QC	!
G01-7-4-7-	ADDED	CONCENT.	ક	LIMITS	1
COMPOUND	(mg/L )	(mg/L )	REC	REC	QUAL
	==	========	====		-=======
Benzene	0.500	0.546	109	79 - 116	ĺ
2-Butanone	0.500	0.503	101	35 - 156	
Carbon tetrachloride	0.500	0.567	113	72 - 133	
Chlorobenzene	0.500	0.543	109	81 - 115	
Chloroform	0.500	0.540	108	81- 122	
1,2-Dichloroethane	0.500	0.552	110	73 - 127	
1,1-Dichloroethene	0.500	0.510	102	65 - 119	
Tetrachloroethene	0.500	0.566	113	78 - 131	
Trichloroethene	0.500	0.545	109	80 - 122	j
Vinyl chloride	0.500	0.627	125	53 - 134	

NOTES (S):
* Values outside of QC limits
Spike Recovery:0 out of10 outside limits
COMMENTS:
FORM III

657 124 SW846 8260B MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A

Level: (low/med) LOW

Lot #: COE130194

WO #: DD6A410P BATCH: 0138262

	SPIKE	SAMPLE	MS	MS		
	ADDED	CONCENT.	CONCENT.	ક	LIMITS	
COMPOUND	(mg/L)	(mg/L )	(mg/L)	REC	REC	QUAL
	=======		=======	=====	=======	========
Benzene	0.500	ND_	0.544	109	73- 123	i
2-Butanone	0.500	0.027	0.532	101	10- 151	
Carbon tetrachloride	0.500	ND	0.563	113	61 - 143	·
Chlorobenzene	0.500	ND	0.546	109	70 - 122	! <del></del>
Chloroform	0.500	ND	0.536	107	65 - 131	
1,2-Dichloroethane	0.500	ND	0.552	110	67 - 132	
1,1-Dichloroethene	0.500	ND	0.518	104	57- 138	·
Tetrachloroethene	0.500	ND	0.555	111	70 - 130	
Trichloroethene	0.500	ND	0.545	109	58- 141	
Vinyl chloride	0.500	ND	0.635	127	51- 133	

NOTES (S):	

<sup>#</sup> Column to be used to flag recovery and RPD values with an asterisk \* Values outside of QC limits RPD: 0 out of 0 outside limits Spike Recovery: 0 out of 10 outside limits COMMENTS:

657 125

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A

Level: (low/med) LOW

Lot #: C0E130194

WO #: DD6A410Q BATCH: 0138262

	SPIKE	MSD	MSD				
!	ADDED	CONCENT.	용	ફ	QC 1	LIMITS	i
COMPOUND	(mg/L)	(mg/L)	REC	RPD	RPD	REC	QUAL
=======================================	========		=====	======	====		
Benzene	0.500	0.552	110	1.5	20	73 - 123	i i
2-Butanone	0.500	0.526	100	1.1	34	10 - 151	<del></del>
Carbon tetrachloride	0.500	0.571	114	1.5	20	61- 143	
Chlorobenzene	0.500	0.550	110	0.69	20	70~ 122	
Chloroform	0.500	0.541	108	0.79	20	65 - 131	
1,2-Dichloroethane	0.500	0.561	112	1.6	20	67- 132	
1,1-Dichloroethene	0.500	0.522	104	0.78	20		
Tetrachloroethene	0.500	0.567	113	2.2	20		 
Trichloroethene	0.500	0.551	110	1.1	20	58 - 141	 
Vinyl chloride	0.500	0.638	128	0.50	20	51 - 133	

# Column to be used to flag recovery and RPD values with an asterisk * Values outside of QC limits
values outside of QC limits
RPD: 0 out of 10 outside limits
Spike Recovery: 0 out of 10 outside limits
COMMENTS:

FORM III

NOTES (S):

657 126 SW846 8260B METHOD BLANK SUMMARY

Lab Name: Severn Trent Laboratories, Inc.

Lab Code: QESPIT

SDG Number:

Lab File ID: 4051703.d

Lot Number: COE130194

Date Analyzed: 05/17/00

Time Analyzed: 14:39

Matrix: SOLID

Date Extracted:05/17/00

GC Column: RTX-624 , ID: .18

Extraction Method: 1311/5030B

BLANK WORKORDER NO.

DD95L101

Instrument ID: HP4

Level: (low/med) LOW

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, LCS, LCSD, MS, MSD:

CLIENT ID. WORK ORDER # FILE ID ANALYZED		SAMPLE	LAB	DATE	DTM:	
		CLIENT ID		<del></del>		
01 CHECK SAMPLE   DDA08101   C   4051707.d   05/17/00   17:04   02 DF/S1/0133/SDC/001   DD6A410P   S   4051705.d   05/17/00   16:07   03 DF/S1/0133/SDC/001   DD6A410Q   D   4051706.d   05/17/00   16:35   04 DF/S1/0133/SDC/001   DD6A4101   4051708.d   05/17/00   17:39   05 DF/S1/0133/SDC/001   DD6A5101   4051709.d   05/17/00   18:08   06 DF/S1/0133/SDC/001   DD6A6101   4051712.d   05/17/00   19:29   07 DF/S1/0133/SDC/001   DD6A7101   4051711.d   05/17/00   19:00   08		,	1	I		
02   DF/S1/0133/SDC/001 A   DD6A410P S   4051705.d   05/17/00   16:07   03   DF/S1/0133/SDC/001 A   DD6A410Q D   4051706.d   05/17/00   16:35   04   DF/S1/0133/SDC/001 A   DD6A4101   4051708.d   05/17/00   17:39   05   DF/S1/0133/SDC/001 B   DD6A5101   4051709.d   05/17/00   18:08   06   DF/S1/0133/SDC/001 C   DD6A6101   4051712.d   05/17/00   19:29   07   DF/S1/0133/SDC/001 D   DD6A7101   4051711.d   05/17/00   19:00   08	01	CHECK SAMPLE		!		
03   DF/S1/0133/SDC/001 A   DD6A410Q D   4051706.d   05/17/00   16:35   04   DF/S1/0133/SDC/001 A   DD6A4101   4051708.d   05/17/00   17:39   05   DF/S1/0133/SDC/001 B   DD6A5101   4051709.d   05/17/00   18:08   06   DF/S1/0133/SDC/001 C   DD6A6101   4051712.d   05/17/00   19:29   07   DF/S1/0133/SDC/001 D   DD6A7101   4051711.d   05/17/00   19:00   08				!	· <del></del>	
04   DF/S1/0133/SDC/001 A   DD6A4101   4051708.d   05/17/00   17:39   05   DF/S1/0133/SDC/001 B   DD6A5101   4051709.d   05/17/00   18:08   06   DF/S1/0133/SDC/001 C   DD6A6101   4051712.d   05/17/00   19:29   07   DF/S1/0133/SDC/001 D   DD6A7101   4051711.d   05/17/00   19:00   08						
06 DF/S1/0133/SDC/001 C DD6A6101 4051712.d 05/17/00 19:29 07 DF/S1/0133/SDC/001 D DD6A7101 4051711.d 05/17/00 19:00 08	04	DF/S1/0133/SDC/001 A	DD6A4101			:
07 DF/S1/0133/SDC/001 D       DD6A7101       4051711.d       05/17/00       19:00         09				4051709.d	05/17/00	18:08
08       ————————————————————————————————————				4051712.d	05/17/00	19:29
09       10       11       12       13       14       15       16       17       18       19       20       21       22       23       24       25		DF/S1/0133/SDC/001 D	DD6A7101	4051711.d	05/17/00	19:00
10	,				l	
11       12         13       3         14       4         15       4         16       4         17       4         18       4         19       4         20       4         21       4         22       4         23       4         25       4	•					
12						
13				<u> </u>		
14       15       16       17       18       19       20       21       22       23       24       25		<u> </u>				
15				<u> </u>		
16       17       18       19       20       21       22       23       24       25		j		<u></u> ;		
17       18       19       20       21       22       23       24       25		<del></del>		<u> </u>		
18       19       20       21       22       23       24       25	-			 		
19		<u> </u>	<u></u>	ļ	ļ	
20		<u> </u>	<u></u>		<u></u> 	<u>-</u>
21					 	
23				 	!	
23	22				1	l
24			' <del></del>			
· · · · · · · · · · · · · · · · · · ·			<del></del>			
· · · · · · · · · · · · · · · · · · ·					i ———	
49	26					·
27	27			<u> </u>	¦	<del></del>
28	28			<u> </u>	<u> </u>	<del> </del>
29	29				ļ	<del></del>
30	30					

COMMENTS:				
	 	 	··-	
<del></del>	 <u></u>	 FORM		

#### 5A VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK BROMOFLUOROBENZENE (BFB)

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: C0E130194

Lab File ID: BF40517

BFB Injection Date: 05/17/00

Instrument ID: HP4

BFB Injection Time: 1041

GC Column: DB624 20M ID: 0.18 (mm) Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
a====		~
50	15.0 - 40.0% of mass 95	25.6
75	30.0 - 60.0% of mass 95	43.0
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	6.9
173	Less than 2.0% of mass 174	$0.0 \ \overline{(0.0)1}$
174	50.0 - 100.0% of mass 95	56.6
175	5.0 - 9.0% of mass 174	4.2 (7.3)1
176	95.0 - 101.0% of mass 174	54.9 ( 97.1)1
177	5.0 - 9.0% of mass 176	3.6 ( 6.5)2
-''		
·	1-Value is % mass 174 2-Value is % mass	176 .

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

I	EPA	LAB	LAB	DATE	TIME (
,	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED	ANALYZED
		=======================================	==========	========	========
01	VSTD50	VSTD50	CC40517	05/17/00	1108
02	VSTD5	VSTD5	1A40517	05/17/00	1145
03	VSTD20	VSTD20	1B40517	05/17/00	1209
04	VSTD100	VSTD100	1D40517	05/17/00	1234
05	VSTD200	VSTD200	1E40517	05/17/00	1303
06	INTRA-LAB BL	DD95L101	4051703	05/17/00	1439
DF(51/0133/50001A MO7	<del>DF 001</del>	DD6A410P	4051705	05/17/00	1607
V Aµ5008	<del>DP-00</del> 1	DD6A410Q	4051706	05/17/00	1635
, 09	INTRA-LAB CH	DDA08101	4051707	05/17/00	1704
OF/sib133/socko1A10	<del>DF-00</del> 1	DD6A4101	4051708	05/17/00	1739
i 1311	<del>S1-001</del>	DD6A5101	4051709	05/17/00	1808
7012	<del>SDC/001</del>	DD6A7101	4051711	05/17/00	1900
<b>√</b> C13	0 <del>133/00</del> 1	DD6A6101	4051712	05/17/00	1929
14		1			
15 مرکب			<u> </u>		
A24/00 16				<u> </u>	
<b>74"</b> 17					
18					
19					
20					<u> </u>
21					
22					

page 1 of 1

FORM V VOA

OLMO3.0

Lab Name: STL-PITTSBURGH Contract:

SAS No.: 40325 SDG No.: C0E130194 Lab Code: STLPIT Case No.:

Date Analyzed: 05/17/00 Lab File ID (Standard): CC40517

Time Analyzed: 1108 Instrument ID: HP4

Heated Purge: (Y/N) N (mm) ID: 0.18 GC Column: DB 624

ļ		IS1 (CBZ)	RT #	IS2 (DCB) AREA #	RT #	IS3 AREA #	RT #
		AREA #	K1 #	AKEA T	=======	========	======
	12 HOUR STD UPPER LIMIT LOWER LIMIT	149779 299558 74890	10.12 10.32 9.92	227254 454508 113627	12.46 12.66 12.26	660677 1321354 330338	6.99 7.19 6.79
	EPA SAMPLE NO.		======			=======	======
01 DF   013   2002 03 04 DF   51   013   200   05 06 07 08 09 10 11 12 13 14 15 16 17	INTRA-LAB BL DF-001 A MA DF-001 A MA INTRA-LAB CH DF-001 B S1-001 B SDC/001 D 0133/001 C	161550	10.12 10.12 10.13 10.13 10.13 10.13 10.13	210752 229690 232577 229414 173160 180580 170776 196867	12.46 12.46 12.46 12.46 12.46 12.46 12.46	792190 736767 739353 743134 666337 677588 585666 736808	7.01 7.00 7.00 7.00 7.01 7.01 7.01 7.01
20 21 22							

(CBZ) = Chlorobenzene-d5 IS1

IS2 (DCB) = 1,4-Dichlorobenzene-d4

= Fluorobenzene IS3

AREA UPPER LIMIT = +100% of internal standard area AREA LOWER LIMIT = - 50% of internal standard area RT UPPER LIMIT = + 0.20 minutes of internal standard RT RT LOWER LIMIT = - 0.20 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

page 1 of 1

FORM VIII VOA

OLM03.0

# GC/MS VOLATILE SAMPLE DATA

657 130

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 001

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD6A4101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

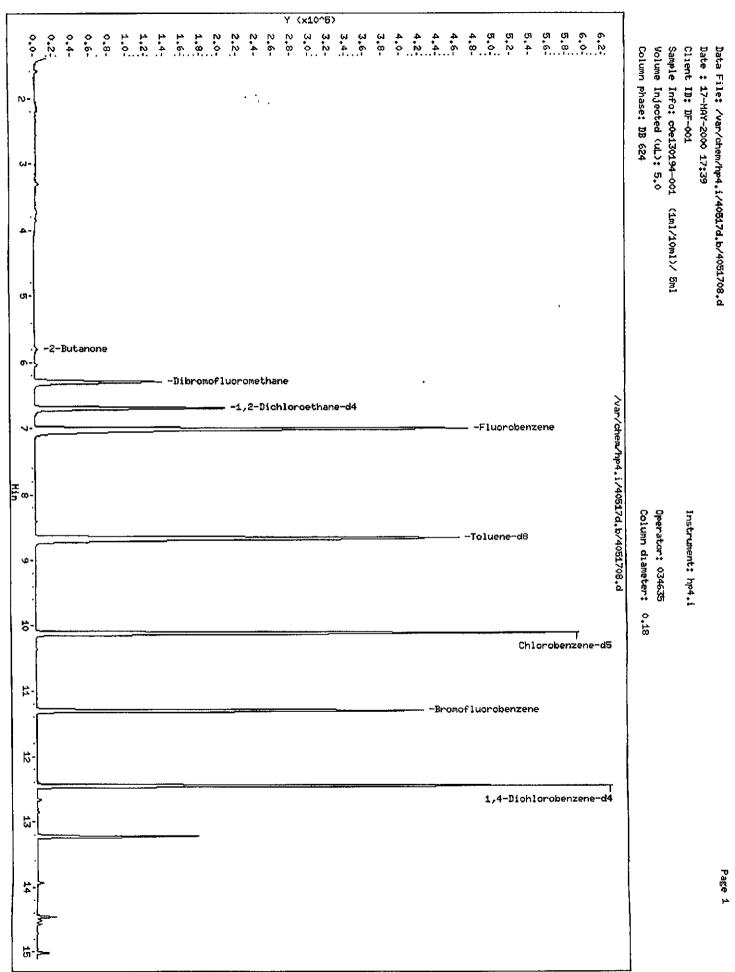
Moisture %:10

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 A

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L o	r ug/kg) mg/L	Q
71-43-2	Benzene	0.050	<u> </u>
78-93-3	2-Butanone	0.027	<u>J</u>
56-23-5	Carbon tetrachloride	0.050	<u>"</u> ן
108-90-7	Chlorobenzene	0.050	UU
67-66-3	Chloroform	0.050	<u> </u>
107-06-2	1,2-Dichloroethane	0.050	ַ
75-35-4	1,1-Dichloroethene	0.050	<u> </u>
127-18-4	Tetrachloroethene	0.050	ַ
79-01-6	Trichloroethene	0.050	<u>"</u>
75-01-4	Vinyl chloride	0.050	



Data File: /var/chem/hp4.i/40517d.b/4051708.d Page 1

Report Date: 17-May-2000 19:09

#### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file: /var/chem/hp4.i/40517d.b/4051708.d

Lab Smp Id: DD6A4101 Client Smp ID: DF-001

Inj Date : 17-MAY-2000 17:39

Operator : 034635 Inst ID: hp4.i

Smp Info : c0e130194-001 (1ml/10ml)/5mlMisc Info : dd6a4101,40517d.b,med1000.m

Comment

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m

Quant Type: ISTD Cal File: 1e40517.d Meth Date: 17-May-2000 14:06 dudeckk Cal Date : 17-MAY-2000 13:03 Als bottle: 14

Dil Factor: 1.00000 Integrator: HP RTE

Compound Sublist: tclp.sub

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

Name	Value	Description
DF Vf Vo Vi	5.000 5.000	Dilution Factor FinalVolume Sample Volume Purge Volume

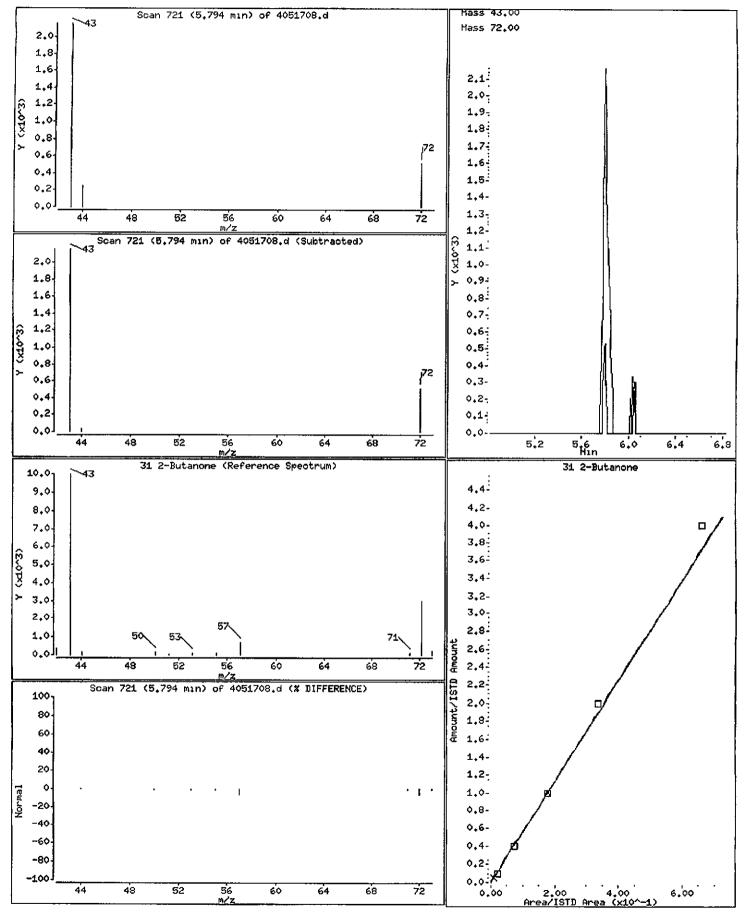
					CONCENTRA	ATIONS
	QUANT SIG				ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	{ UG/L}
	====	==				
* 46 Fluorobenzene	96	7.011	6.994 (1.000)	666337		
* 69 Chlorobenzene-d5	119	10.129	10.124 (1 000)	131368		
* 92 1,4-Dichlorobenzene-d4	152	12 458	12.460 (1 000)	173160		
\$ 43 1,2-Dichloroethane-d4	65	6 693	6.683 (0 955)	171017	273.909	54.78
\$ 39 Dibromofluoromethane	113	6.296	6.285 (0.898)	117491	257.714	51.54
\$ 59 Toluene-d8	98	8 662	8 645 (0.855)	568613	257.668	51.53
\$ 80 Bromofluorobenzene	95	11 309	11 310 (1 116)	198266	232.569	46.51
3 Vinyl Chloride	62.00	Com	pound Not Detecte	ed		
12 1,1-Dichloroethene	96.00	Соп	pound Not Detect	eđ.		
31 2-Butanone	43	5.794	5 821 (0 826)	6420	13.5520	2.710
37 Chloroform	83.00	Соп	pound Not Detecte	⊵đ.		
41 Carbon Tetrachloride	117 00	Con	pound Not Detecte	ed.		
42 Benzene	78 00	Com	pound Not Detecte	ed.		
45 1,2-Dichloroethane	62 00	Соп	pound Not Detecte	eđ.		,
47 Trichloroethene	130.00	Соп	pound Not Detect	ed.		

Page 2

Data File: /var/chem/hp4.i/40517d.b/4051708.d Report Date: 17-May-2000 19:09

					CONCENTRATIONS		
	QUANT SIG				ON-COLUMN	FINAL	
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( UG/L)	
		= =					
65 Tetrachloroethene	164.00	Compound Not Detected					
70 Chlorobenzene	112 00	Compound Not Detected					





Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 002

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD6A5101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

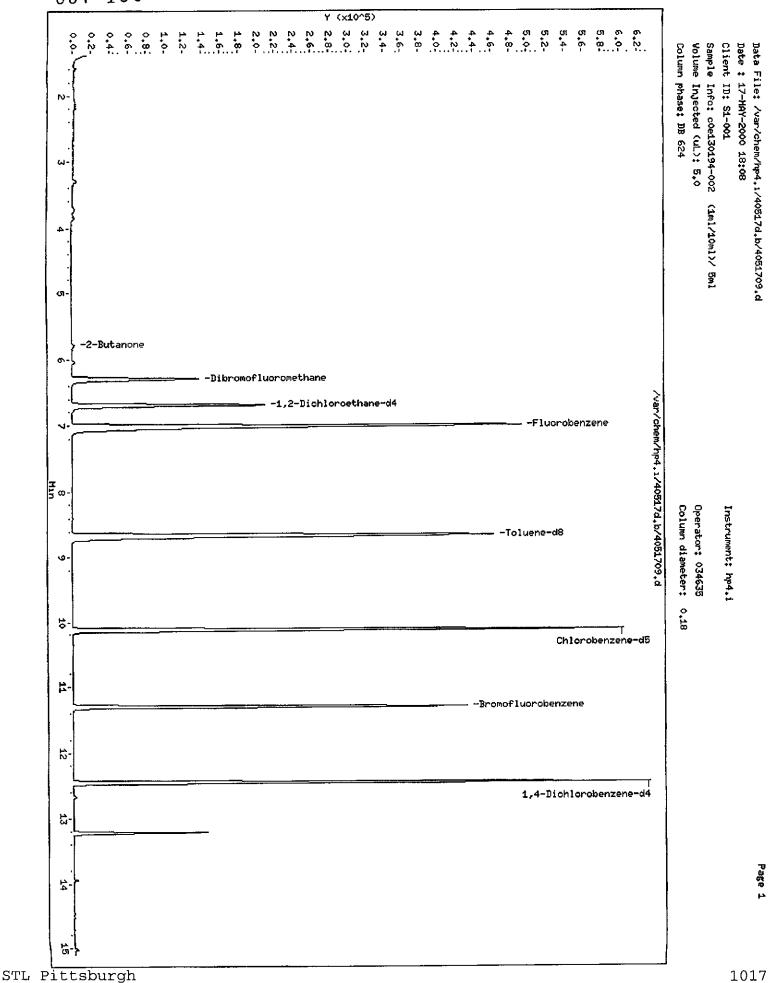
Moisture %:8.6

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 B

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L	<u>Q</u>
71-43-2	Benzene	0.050	<u></u>
78-93-3	2-Butanone	0.027	J
56-23-5	Carbon tetrachloride	0.050	<u>" </u>
108-90-7	Chlorobenzene	0.050	<u> </u>
67-66-3	Chloroform	0.050	<u> </u>
107-06-2	1,2-Dichloroethane	0.050	<u></u>
75-35-4	1,1-Dichloroethene	0.050	<u> </u>
127-18-4	Tetrachloroethene	0.050	<u> </u>
79-01-6	Trichloroethene	0.050	<u> </u>
75-01-4	Vinyl chloride	0.050	\



Data File: /var/chem/hp4.i/40517d.b/4051709.d Page 1

Report Date: 17-May-2000 19:09

#### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file: /var/chem/hp4.i/40517d.b/4051709.d Lab Smp Id: DD6A5101 Client Client Smp ID: S1-001

Inj Date : 17-MAY-2000 18:08

Operator: 034635 Inst ID: hp4.i

Smp Info : c0e130194-002 (1ml/10ml)/5ml Misc Info : dd6a5101,40517d.b,med1000.m

Comment

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m

Meth Date: 17-May-2000 14:06 dudeckk Quant Type: ISTD Cal File: 1e40517.d Cal Date : 17-MAY-2000 13:03

Als bottle: 15

Dil Factor: 1.00000 Integrator: HP RTE

Compound Sublist: tclp.sub

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

			• • • • • • • • • • • • • • • • • • • •	. 1
	Name	Value	Description	11/1/1
-	DF Vf		Dilution Factor FinalVolume	Diloi
	Vo	5.000	Sample Volume	<b>T</b>
	Vi	5.000	Purge Volume	•

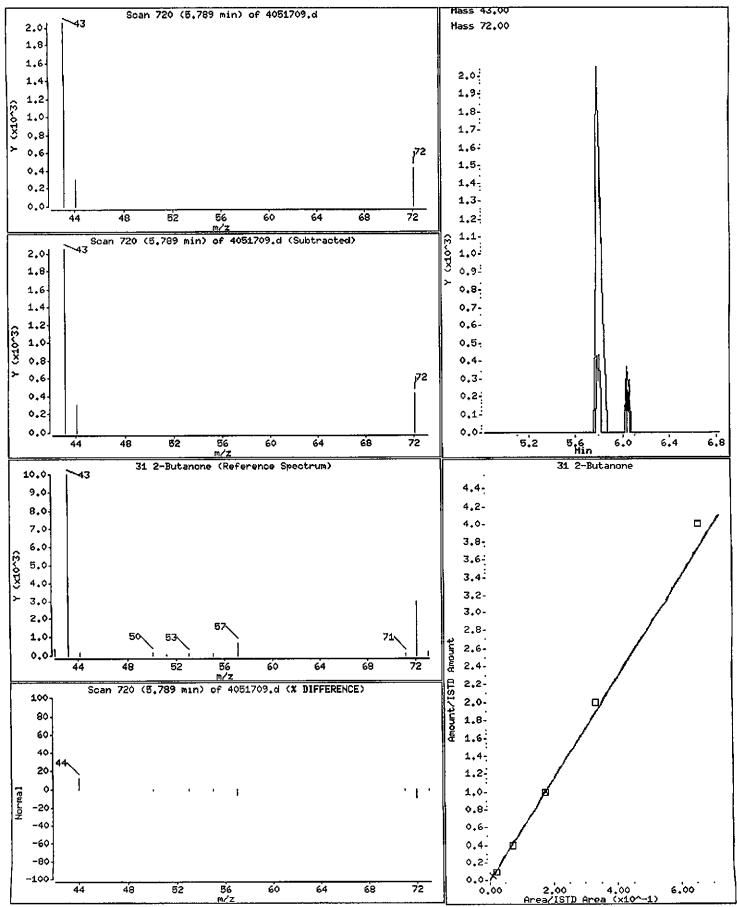
						CONCENTRA	TIONS
		QUANT SIG				ON-COLUMN	FINAL
Compounds		MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( UG/L)
	****			医医医内耳 经过户报过单			======
* 46 Fluorobenzene		96	7.012	6 994 (1.000)	677588		
* 69 Chlorobenzene-	d5	119	10.130	10.124 (1 000)	135805		
* 92 1,4-Dichlorobe	enzene-d4	152	12.459	12 460 (1.000)	180580		
\$ 43 1,2-Dichloroet	:hane-d4	65	6 694	6.683 (0 955)	170681	268 832	53.77
\$ 39 Dibromofluorom	nethane	113	6 297	6.285 (0.898)	120533	259.997	52 00
\$ 59 Toluene-d8		98	8 663	8.645 (0 855)	563521	247.017	49.40
\$ 80 Bromofluorober	ızene	95	11.310	11.310 (1.116)	196647	223 133	44.63
3 Vinyl Chloride	<b>:</b>	62.00	Com	pound Not Detect	eđ.		
12 1,1-Dichloroet	thene	96 00	Con	pound Not Detect	ed		
31 2-Butanone		43	5.789	5.821 (0.826)	6520	13.5346	2 707
37 Chloroform		83.00	Соп	pound Not Detect	ed		
41 Carbon Tetrac	nloride	117.00	Com	pound Not Detect	ed		
42 Benzene		78.00	Con	pound Not Detect	ed.		
45 1,2-Dichloroe	thane	62 00	Con	pound Not Detect	ed		
47 Trichloroether	ne	130 00	Con	pound Not Detect	ed.		

Data File: /var/chem/hp4.i/40517d.b/4051709.d Report Date: 17-May-2000 19:09

Page 2

					CONCENTRA	ATIONS
	QUANT SIG				ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( UG/L)
**********	<b>35.0</b> 0		=======================================	********	B=S====	
65 Tetrachloroethene	164.00	Co	mpound Not Detect	.ed		
70 Chlorobenzene	112.00	Co	mpound Not Detect	ed		





657 140

#### UXB INTERNATIONAL

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 003

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD6A6101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

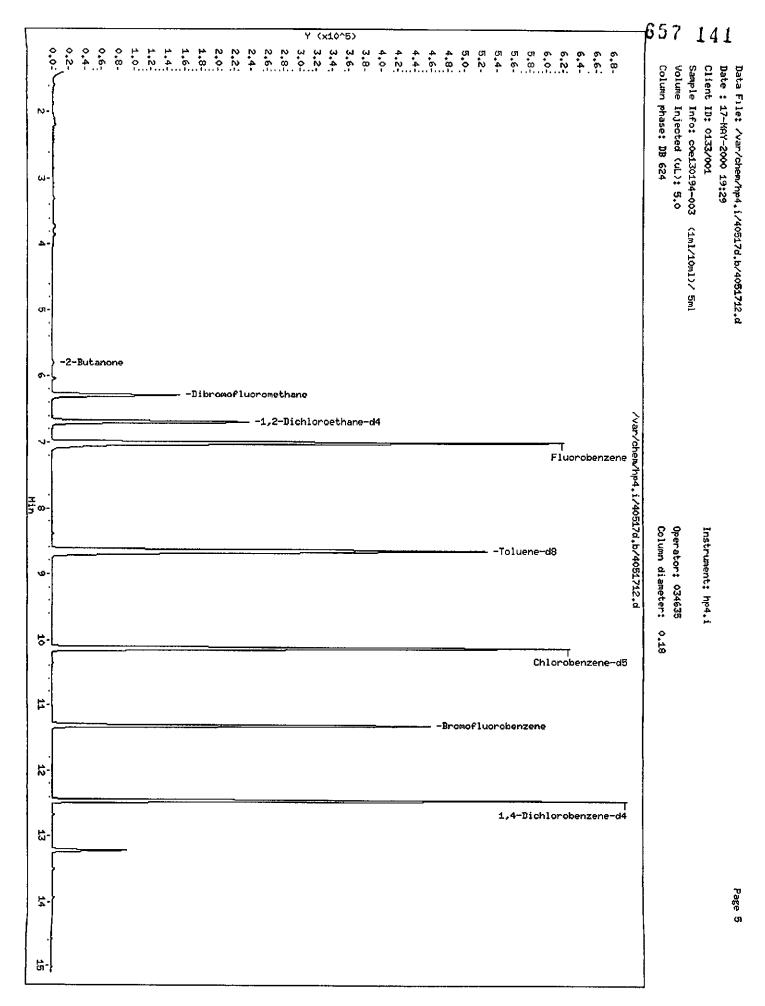
Moisture %:10

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 C

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L c	or ug/kg) mg/L	Q
71-43-2	Benzene	0.050	ט
78-93-3	2-Butanone	0.023	J
56-23-5	Carbon tetrachloride	0.050	וט
108-90-7	Chlorobenzene	0.050	<u>י</u>
67-66-3	Chloroform	0.050	וש ו
107-06-2	1,2-Dichloroethane	0.050	ט
75-35-4	1,1-Dichloroethene	0.050	ן ט
127-18-4	Tetrachloroethene	0.050	<u>"</u>
79-01-6	Trichloroethene	0.050	ט
75-01-4	Vinyl chloride	0.050	<u>"</u>



Data File: /var/chem/hp4.i/40517d.b/4051712.d

Report Date: 17-May-2000 19:47

Page 1

1 /2

#### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file : /var/chem/hp4.i/40517d.b/4051712.d

Client Smp ID: 0133/001 Lab Smp Id: DD6A6101

Inj Date : 17-MAY-2000 19:29

Inst ID: hp4.i Operator: 034635

Smp Info : c0e130194-003 (1ml/10ml)/5ml Misc Info : dd6a6101,40517d.b,med1000.m

Comment

: /var/chem/hp4.i/40517d.b/8260bh2o.m Method

Meth Date: 17-May-2000 14:06 dudeckk Quant Type: ISTD Cal File: 1e40517.d Cal Date : 17-MAY-2000 13:03

Als bottle: 18

Dil Factor: 1.00000 Integrator: HP RTE

Compound Sublist: tclp.sub

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

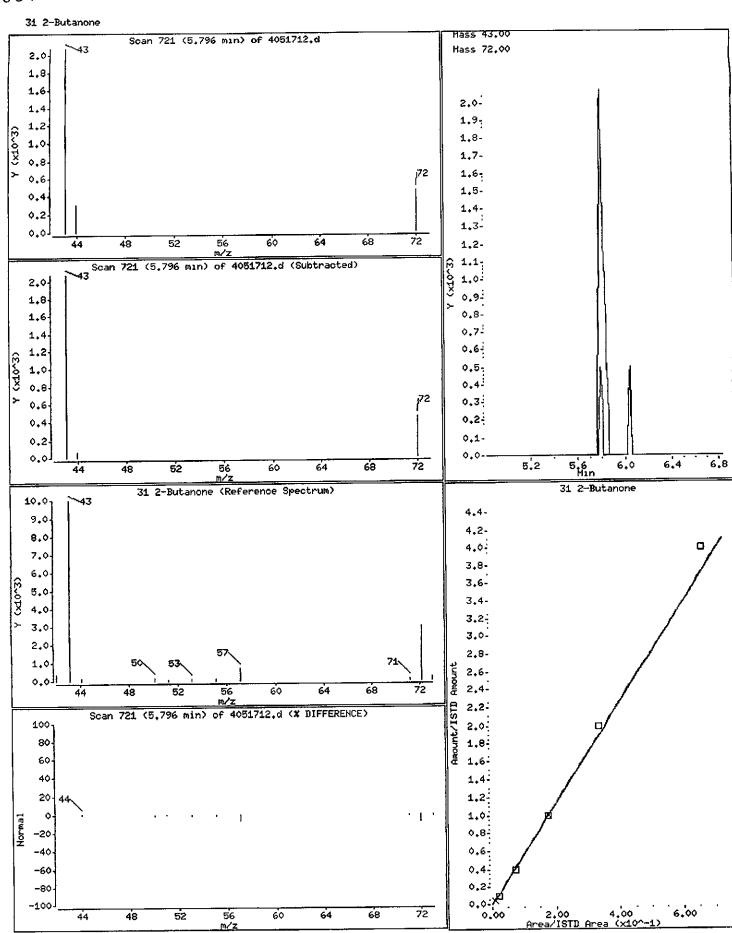
Name	Value	Description	VIIZINIA
DF Vf Vo Vi	5.000 5.000	Dilution Factor FinalVolume Sample Volume Purge Volume	ħ,

						CONCENTRA	RYTIONS
		QUANT SIG				ON-COLUMN	FINAL
Co	ompounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( UG/L)
==		2225	33		2002544	****	****
*	46 Fluorobenzene	96	7 007	6 994 (1 000)	736808		
*	69 Chlorobenzene-d5	119	10.131	10 124 (1.000)	142009		
*	92 1,4-Dichlorobenzene-d4	152	12.460	12 460 (1 000)	196867		
\$	43 1,2-Dichloroethane-d4	65	6.689	6 683 (0 955)	182429	264.241	52.85
\$	39 Dibromofluoromethane	113	6.292	6 285 (0.898)	128727	255 354	51 07
\$	59 Toluene-d8	98	8.658	8.645 (0.855)	633276	265.467	53.09
\$	80 Bromofluorobenzene	95	11.311	11 310 (1.116)	214408	232 658	46 53
	3 Vinyl Chloride	62.00	Con	pound Not Detect	ed.		
	12 1,1-Dichloroethene	96.00	Con	pound Not Detect	eđ		
	31 2-Butanone	43	5.796	5 821 (0.827)	6050	11 5495	2.310
	37 Chloroform	83.00	Сол	pound Not Detect	eđ		
	41 Carbon Tetrachloride	117.00	Con	pound Not Detect	ed		
	42 Benzene	78 00	Con	pound Not Detect	ed		
	45 1,2-Dichloroethane	62 00	Con	mpound Not Detect	ed.		
	47 Trichloroethene	130 00	Con	mpound Not Detect	ed		

Page 2

Data File: /var/chem/hp4.i/40517d.b/4051712.d Report Date: 17-May-2000 19:47

					CONCENTRATIONS	
	QUANT SIG				ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( UG/L)
	4555	==	*****		**==**	======
65 Tetrachloroethene	164 00	Compound Not Detected				
70 Chlorobenzene	112 00	Cor	mpound Not Detect	ed		



Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 004

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD6A7101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

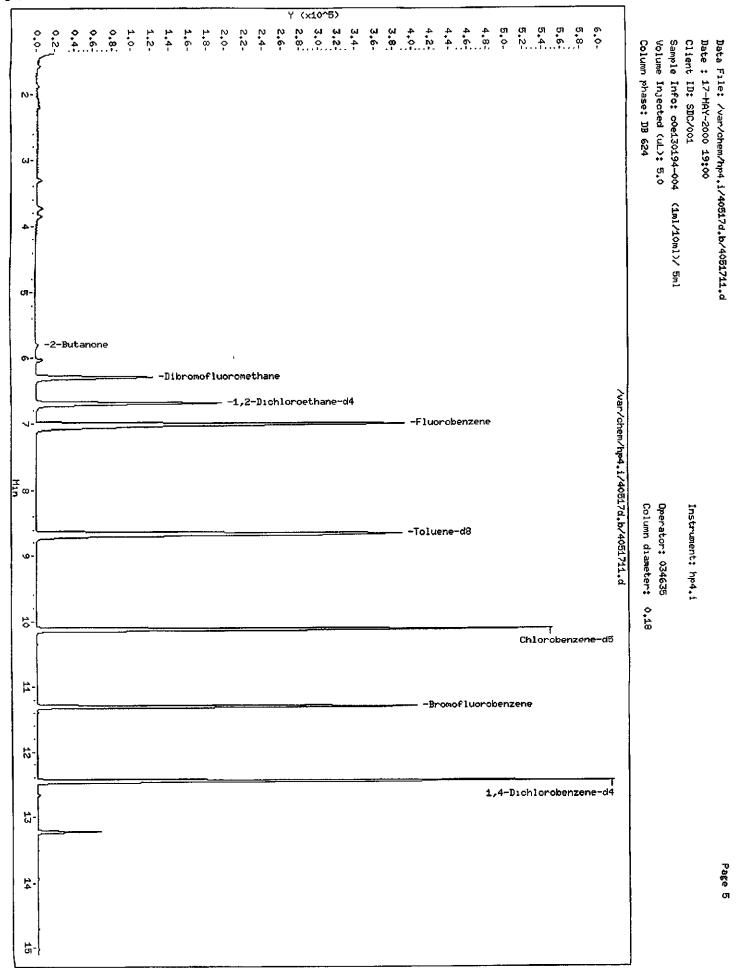
Moisture %:11

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 D

## CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L o	r ug/kg) mg/L	Q
71-43-2	Benzene	0.050	ן די
78-93-3	2-Butanone	0.039	jσ
56-23-5	Carbon tetrachloride	0.050	וס
108-90-7	Chlorobenzene	0.050	ט
67-66-3	Chloroform	0.050	יט
107-06-2	1,2-Dichloroethane	0.050	ו ט
75-35-4	1,1-Dichloroethene	0.050	i vi
127-18-4	Tetrachloroethene	0.050	Ū
79-01-6	Trichloroethene	0.050	ט ו
75-01-4	Vinyl chloride	0.050	יט



Data File: /var/chem/hp4.i/40517d.b/4051711.d

Report Date: 17-May-2000 19:19

#### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file : /var/chem/hp4.i/40517d.b/4051711.d

Lab Smp Id: DD6A7101 Client Smp ID: SDC/001

Inj Date : 17-MAY-2000 19:00

Operator : 034635 Inst ID: hp4.i

Smp Info : c0e130194-004 (1ml/10ml)/5ml Misc Info : dd6a7101,40517d.b,med1000.m

Comment

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m Meth Date : 17-May-2000 14:06 dudeckk Quant T Cal Date : 17-MAY-2000 13:03 Cal Fil Quant Type: ISTD Cal File: 1e40517.d

Als bottle: 17

Dil Factor: 1.00000 Integrator: HP RTE

Compound Sublist: tclp.sub

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

Name	Value	Description
DF Vf Vo Vi	5.000 5.000	Dilution Factor FinalVolume Sample Volume Purge Volume

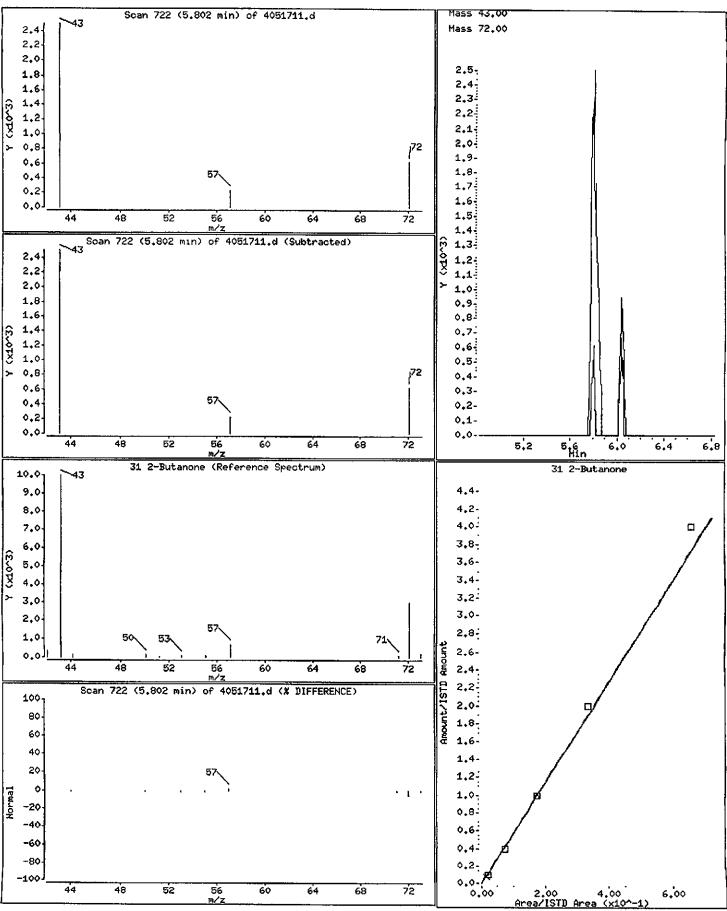
DI 2/12/00

FINAL ( UG/L)
( UG/L)
0022022
59.47
53 73
48 08
45 47
3 927

Data File: /var/chem/hp4.i/40517d.b/4051711.d Report Date: 17-May-2000 19:19

					CONCENTRATIONS			
	QUANT SIG				ON-COLUMN	FINAL		
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( UG/L)		
	====		***** ==== <b>*</b>			======		
65 Tetrachloroethene	164.00	Compound Not Detected						
70 Chlorobenzene	112 00	Co	mpound Not Detect	ed				

Page 2



# GC/MS VOLATILE CALIBRATION DATA

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.:

SAS No.: 40325 SDG No.: 40826D

Instrument ID: HP4

Calibration Date(s): 05/17/00 05/17/00

Heated Purge: (Y/N) N

Calibration Time(s): 1108

1303

GC Column: DB 624 ID: 0.18 (mm)

LAB FILE ID: RRF5 =1A40517 RRF20 =1B40517 RRF50 =CC40517 RRF100=1D40517 RRF200=1E40517							
COMPOUND	RRF5	RRF20	RRF50	RRF100		RRF	RSD
Dichlorodifluoromethane	0.114	0.118	0,127			0.122	5.1
Chloromethane	* 0.265						4.0
Vinyl Chloride	0.254	0.247					4.1
Bromomethane	0.036	0.036					6.7
Chloroethane .	0.041	0.044	0.049	0.046			
Trichlorofluoromethane	0.061	0.057					16.3
1,1-Dichloroethene	0.157						8.4
Methylene Chloride	0.208	0.183					10.1
trans-1,2-Dichloroethene	0.171	0.192	0.216		0.186		8.8
1,1-Dichloroethane	* 0.360	0.387					6.4
cis-1,2-dichloroethene	0.201	0.218	0.241			0.224	7.2
Chloroform	0.284	0.303	0.332		0.311	0.311	6.1
Bromochloromethane	0.076	0.088	0.094			0.088	8.2
1,1,1-Trichloroethane	0.192	0.224	0.246		0.240	0.230	9.9
Carbon Tetrachloride	0.111	0.133	0.159	0.162	0.158	0.145	15.3
1,2-Dichloroethane	0.226	0.247	0.274			0.143	7.6
Benzene	0.827	0.894	0.979		0.908	0.234	6.5
Trichloroethene	0.164	0.186			0.197	0.192	9.2
1,2-Dichloropropane	0.226	0.244	0.272	0.269			7.5
Bromodichloromethane	0.160	0.189	0.230	0.234	0.230	0.209	15.8
cis-1,3-Dichloropropene	0.253	0.299	0.356	0.358	0.348		14.1
Toluene	4.135	4.446		4.708	4.325	4.483	6.1
trans-1,3-Dichloropropene	0.985	1.213	1.380	1.418	1.353	1.270	13.9
1,1,2-Trichloroethane	0.759	0.873	0.920	0.913	0.859	0.865	7.5
Tetrachloroethene	0.524	0.571	0.631	0.636		0.593	7.9
Dibromochloromethane	0.406	0.512	0.633	0.674	0.667	0.578	20.1
Chlorobenzene_	* 2.534	2.729	2.956	2.920	2.728	2.773	6.1
Ethylbenzene	1.477	1.552	1.687		1.493	1.569	5.8
Styrene	2.631	3.130	3.468	3.459	3.030	3.144	11.0
Bromoform	* 0.185	0.240	0.335	0.377	0.376	0.303	28.5
1,1,2,2-Tetrachloroethane	* 0.710	0.788	0.865	0.857	0.807	0.805	7.7
1,3-Dichlorobenzene	1.201	1.257	1.393	1.376	1.191	1.284	7.4
1,4-Dichlorobenzene	1.230	1.291	1.377	1.364	1.282	1.309	4.7
1,2-Dichlorobenzene	1.112	1.167	1.278	1.267	1.215	1.208	5.8
Dibromomethane	0.094	0.102	0.118	0.113	0.106	0.107	8.7
1,2-Dibromoethane	0.650	0.746	0.811	0.816	0.752	0.755	8.9
1,1,1,2-Tetrachloroethane	0.543	0.641	0.753	0.762	0.704	0.681	13.3
Compounds with required min	! <u>_</u>			]			

\* Compounds with required minimum RRF and maximum %RSD values. All other compounds must meet a minimim RRF of 0.010.

page 1 of 2

FORM VI VOA

OLMO3.0

657 152

Lab Name: STL-PITTSBURGH Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: 40826D

Instrument ID: HP4 Calibration Date(s): 05/17/00 05/17/00

Heated Purge: (Y/N) N Calibration Time(s): 1108 1303

GC Column: DB 624 ID: 0.18 (mm)

LAB FILE ID: RRF5 =1A40517 RRF20 =1B40517 RRF50 =CC40517 RRF100=1D40517 RRF200=1E40517							
COMPOUND	RRF5	RRF20	RRF50	RRF100	RRF200	RRF	RSD
1,2,3-Trichloropropane	0.199	1	I	1	ł.	0.219	
1,2-Dibromo-3-chloropropane	0.056						
2,2-Dichloropropane	0.198						
1,1-Dichloropropene	0.225						8.0
1,3-Dichloropropane	1.359						6.6
n-Propylbenzene	0.801					0.878	6.0
Bromobenzene	0.571						
1,3,5-Trimethylbenzene	2.281				2.494		
2-Chlorotoluene	0.698			0.777			
4-Chlorotoluene	0.718	0.758		0.809	0.741		
tert-Butylbenzene	1.919	2.078		2.308	2.152		7.5
1,2,4-Trimethylbenzene	2.282	2.543			2.536		7.4
sec-Butylbenzene	3.229			3.725	3.426		
4-Isopropyltoluene	2.432	2.591		2.806	2.364		8.7
n-Butylbenzene	2.754	2.876	3.137		3.005		
1,2,4-Trichlorobenzene	0.606			0.568	0.598		
Hexachlorobutadiene	0.292	0.240	0.243	0.276	0.299		
Naphthalene	1.521	1.100	1.061	1.222	1.336		15.0
1,2,3-Trichlorobenzene	0.682	0.373	0.357	0.379			30.8
Acetone	0.114	0.102	0.098	0.088	0.114		10.8
Carbon Disulfide	0.450	0.486		0.534	0.453		9.2
2-Butanone	0.199	0.181		0.167	0.166		7.6
4-Methyl-2-Pentanone	1.561	1.490	1.562	1.542	1.490		2.4
2-Hexanone	1.230	1.063	1.089	1.104	1.043		6.6
Methyl tert-butyl ether	0.401	0.474	0.506	0.500	0.493		9.1
Isopropylbenzene	3.890	4.473	4.947	4.903	4.413		9.5
1,2-Dichloroethene (total)	0.186	0.205	0.228	0.220	0.204	0.209	7.8
Xylenes (total)	1.645	1.809		2.029	1.857		8.4
=======================================	======	======	======	======	=====		=====
Dibromofluoromethane	0.192	0.182	0.170	0.167	0.145	0.171	10.2
1,2-Dichloroethane-d4	0.265	0.245	0.235	0.227	0.199		10.2
Toluene-d8	4.746	4.524	4.201	4.068	3.458	4.199	11.7
Bromofluorobenzene	1.763	1.721	1.643	1.610	1.375	1.622	9.3
Compounds with required mini							

\* Compounds with required minimum RRF and maximim %RSD values. All other compounds must meet a minimim RRF of 0.010.

page 2 of 2

FORM VI VOA

OLM03.0

## INITIAL CALIBRATION REPORT

657 153

Instrument ID: hp4.i Lab File ID: 1a40517.d Analysis Type: WATER

Injection Date: 17-MAY-2000 11:45
 Lab Sample ID: vstd5
Method File: /var/chem/hp4.i/40517d.b/8260bh2

COMPOUND	TR
= = # - = = = = = = = = = = = = = = = =	
(Xylenes (total)	8
1,2-Dichloroethene (total)	
Dichlorodifluoromethane	1 9
Chloromethane	4
Vinyl Chloride	4
Bromomethane	6
Chloroethane	1 8
Trichlorofluoromethane	16
1,1-Dichloroethene	8
Carbon Disulfide	, 9
Acetone	10
Methylene Chloride	10
trans 1,2-Dichloroethene	1 8
Methyl tert-butyl ether	9
1,1-Dichloroethane	j e
2,2 Dichloropropane	11
cls-1,2-dichloroethene	1 7
2-Butanone	1 7
Bromochloromethane	8
Chloroform	6
1,1,1-Trichloroethane	, 9
Dibromofluoromethane	10
Carbon Tetrachloride	15
1,1-Dichloropropene	8
Benzene	6
1,2-Dichloroethane-d4	10
1,2-Dichloroethane	7
Trichloroethene	9
1,2-Dichloropropane	. 7
Dibromomethane	8
Bromodichloromethane	15
cis-1,3-Dichloropropene	14
4-Methyl-2-Pentanone	2
Toluene-d8	11
Toluene	6
trans-1,3-Dichloropropene	13
1,1,2-Trichloroethane	1 7
Tetrachloroethene	1 7
1,3-Dichloropropane	,

Report Date: 05/17/2000

# 657 154

## INITIAL CALIBRATION REPORT

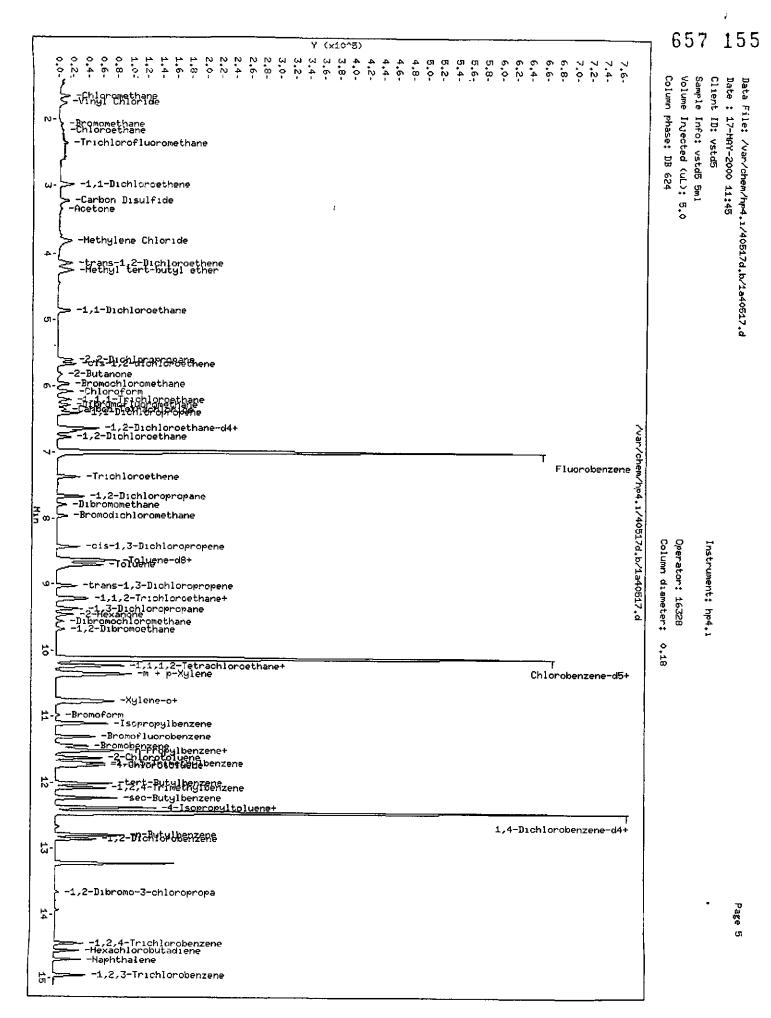
Instrument ID: hp4.i Lab File ID: 1a40517.d Analysis Type: WATER

Injection Date: 17-MAY-2000 11:45

Lab Sample ID: vstd5
Method File: /var/chem/hp4.i/40517d.b/8260bh2

COMPOUND		%RS	D
	≖=== = 		6
Dibromochloromethane	í	20	) 1
1,2-Dibromoethane	i	8	, 9
Chlorobenzene		6	: 1
Ethylbenzene	ĺ	5	В
1,1,1,2-Tetrachloroethane	Ĺ	13	3
m + p-Xylene	i	8	0
Xylene o		8	4
Styrene	ĺ	11	0
Bromoform	i	28	5
Isopropylbenzene	j	9	5
Bromofluorobenzene	1	9	3
Bromobenzene	Ť	8	0
1,1,2,2-Tetrachloroethane	İ	7	7
n-Propylbenzene	Ĺ	6	0
1,2,3-Trichloropropane	ļ	6	9
2-Chlorotoluene	1	5	2
1,3,5 Triwethylbenzene	1	6	6
4-Chlorotoluene		5	9
tert-Butylbenzene	1	7	5
1,2,4-Trimethylbenzene	1	7	4
sec-Butylbenzene	1	6	2
4 Isopropyltoluene	,	8	7
1,3-Dichlorobenzene	1	7	4
1,4-Dichlorobenzene	1	4	7
n-Butylbenzene	1	6	1
1,2 Dichlorobenzene		5	8
1,2-Dibromo-3-chloropropane	I	26	1
1,2,4-Trichloropenzene	l	8	6
Hexachlorobutadiene	ļ	10	2
Naphthalene		15	0
1,2,3-Trichlorobenzene		30	8

The average of all %RSD's in the initial calibration is 9 4



Data File: /var/chem/hp4.i/40517d.b/la40517.d Page 1

Report Date: 17-May-2000 13:43

#### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file : /var/chem/hp4.i/40517d.b/1a40517.d

Lab Smp Id: vstd5 Inj Date : 17-MAY-2000 11:45 Client Smp ID: vstd5

Operator : 16328 Inst ID: hp4.i

Smp Info : vstd5 5ml

Misc Info : ,40517d.b, med1000.m

K 5/17/0 Comment Method : /var/chem/hp4.i/40517d.b/8260bh2o.m Meth Date : 17-May-2000 13:43 dudeckk Quant

Quant Type: ISTD

Cal Date : 17-MAY-2000 13:03 Cal File: 1e40517.d

Als bottle: 5 Calibration Sample, Level: 1

Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: 2-padep.sub

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

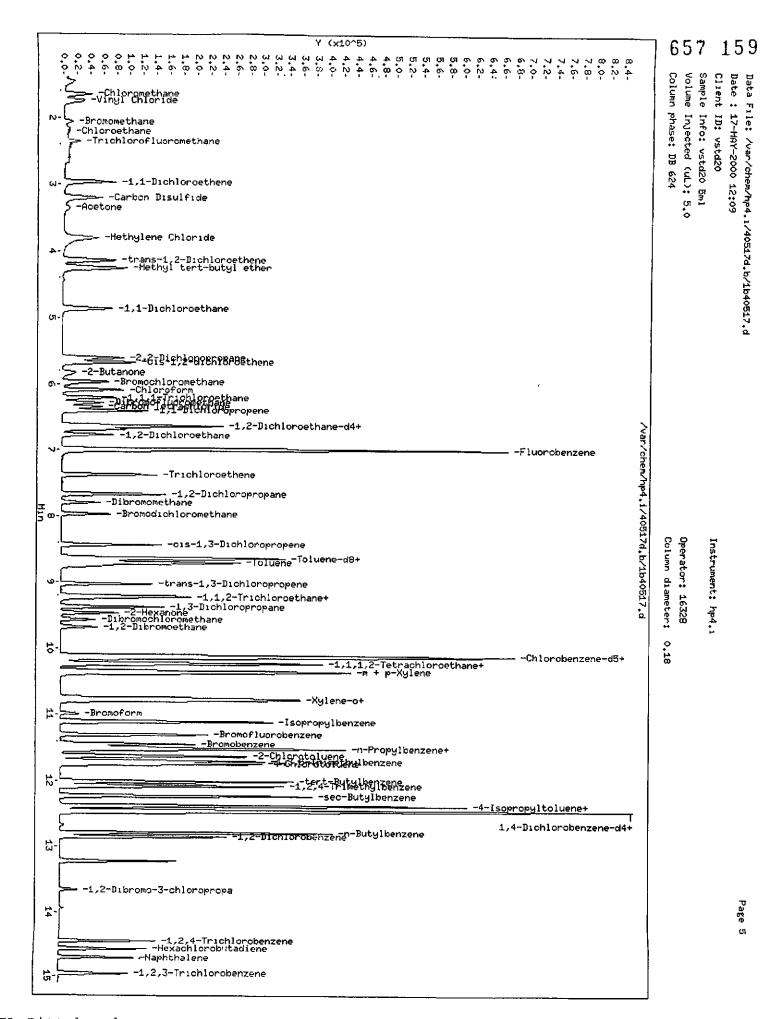
Name	Value	Description
DF Vf Vo Vı	5.000 5.000	Dilution Factor FinalVolume Sample Volume Purge Volume

						AMOUN	TS
		QUANT SIG				CAL-AMT	ON-COL
Com	pounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
===	=	====	==	=======================================	======================================	*****	======
*	46 Fluorobenzene	96	6 994	6 996 (1 000)	695137		
*	69 Chlorobenzene-d5	119	10 124	10 126 (1 000)	149448		
•	92 1,4 Dichlorobenzene-d4	152	12 459	12.462 (1 000)	215803		
\$	39 Dibromofluoromethane	113	6 285	6 287 (0 899)	13325	25 0000	28 02
\$	43 1,2-Dichloroethane-d4	65	6 682	6 684 (0 955)	18405	25 0000	28.26
\$	59 Toluene-d8	98	8 645	8 647 (0.854)	70926	25.0000	28.25
\$	80 Bromofluorobenzene	95	11 304	11 312 (1.117)	26349	25 0000	27 17
	1 Dichlorodifluoromethane	85	1 455	1 488 (0 208)	7900	25 0000	23 31
	2 Chloromethane	50	1.645	1 647 (0 235)	18434	25 0000	24 13
	3 Vinyl Chloride	62	1 743	1 769 (0.249)	17657	25 0000	24.41
	4 Bromomethane	94	2.055	2 069 (0.294)	2501	25 0000	23 85
	5 Chloroethane	64	2.158	2 203 (0 309)	2879	25 0000	23 58
	6 Trichlorofluoromethane	101	2 360	2.350 (0 337)	4237	25 0000	21.85
:	12 1,1-Dichloroethene	96	2 984	2 943 (0.427)	10943	25 0000	24 25
:	13 Acetone	43	3 351	3 377 (0 479)	7965	25 0000	27 72

					AMOŲI	VTS
0	QUANT SIG				CAL-AMT	ON-COL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	(ng)
	<b>基拉</b> 莱森	==		<b>t</b> ======	***====	*====a
15 Carbon Disulfide	76	3 222	3 182 (0 461)	31285	25 0000	22 76
18 Methylene Chloride	84	3 815	3 799 (0 545)	14444	25 0000	26 02
19 trans 1,2-Dichloroethe		4 151	4.129 (0 594)	11915	25 0000	22 08
20 Methyl tert butyl ethe		4 255	4 270 (0 608)	27866	25 0000	21 11
24 1,1-Dichloroethane	63	4 873	4 857 (0 697)	25017	25 0000	22 64
27 2,2-Dichloropropane	77	5.613	5 609 (0 802)	13751	25 0000	20 62
28 cis-1,2-dichloroetnene		5.674	5 670 (0 811)	13950	25 0000	22.44
M 29 1,2-Dichloroethene (to				25865	50 0000	44 54
30 Bromochloromethane	128	5 979	5 975 (0 855)	5304	25 0000	21 55
31 2-Butanone	43	5.826	5 847 (0.833)	13827	25 0000	27 98
37 Chloroform	83	6 089	6 091 (0 871)	19719	25 0000	22.82
38 1,1,1-Trichloroethane	97	6 212	6 202 (0 888)	13364	25 0000	20 91
40 1,1-Dichloropropene	75	6 407	6 403 (0 916)	15669	25 0000	22 04
41 Carbon Tetrachloride	117	6.352	6 342 (0 908)	7703	25 0000	19.19
42 Benzene	78	6 640	6 636 (0 949)	57472	25 0000	22 64
45 1,2-Dichloroethane	62	6 774	6 770 (0 969)	15678	25 0000	22 18
47 Trichloroethene	130	7 385	7 381 (1 056)	11390	25 0000	21.35
49 1,2-Dichloropropane	63	7 673	7 675 (1 097)	15727	25 0000	22 32
50 Dibromomethane	93	7 801	7 797 (1 115)	6559	25 0000	22 07
53 Bromodichloromethane	83	7 966	7 974 (1 139)	11111	25 0000	19 12
57 cis-1,3-Dichloropropend	75	8 437	8 439 (1 206)	17617	25 0000	19 62
58 4-Methyl-2-Pentanone	43	8 626	8 635 (0 852)	23331	25 0000	25 52
60 Toluene	91	8 712	B 714 (0 861)	61794	25 0000	23 06
61 trans-1,3-Dichloroprope	ene 75	9 030	9 032 (0 892)	14723	25 0000	19 40
63 1,3-Dichloropropane	76	9 378	9.380 (0 926)	20309	25 0000	22 53
64 1,1,2-Trichloroethane	97	9 207	9.215 (0 909)	11338	25 0000	21 93
65 Tetrachloroethene	164	9 232	9 234 (0.912)	7824	25.0000	22 07
66 2-Hexanone	43	9 458	9 466 (0 934)	18386	25 0000	27 81
67 Dibromochloromethane	129	9 568	9.570 (0 945)	6062	25 0000	17 53
68 1,2 Dibromoethane	107	9 684	9 686 (0 957)	9720	25 0000	21 53
70 Chlorobenzene	112	10 149	10 157 (1.002)	37868	25.0000	22 84
71 1,1,1,2-Tetrachloroetha	ine 131	10.240	10 249 (1.011)	8120	25.0000	19 95
72 Ethylbenzene	106	10 228	10 236 (1 010)	22068	25 0000	23 53
73 m + p-Xylene	106	10 350	10 359 (1 022)	50538	50 0000	45.09
74 Xylene-o	106	10 754	10 756 (1 062)	24591	25.0000	22 01
M 75 Xylenes (total)	106			75129	25.0000	67 24
76 Styrene	104	10 784	10 787 (1 065)	39318	25 0000	20 92
77 Bromoform	173	10 986	10 988 (1 085)	2766	25 0000	15.28
78 Isopropylbenzene	105	11 108	11 111 (1.097)	58142	25 0000	21 49
79 Bromobenzene	156	11.445	11 453 (0 919)	12324	25 0000	22 17
81 n-Propylbenzene	120	11 518	11 526 (0 924)	17289	25 0000	22 81
82 2-Chlorotoluene	126	11 628	11 636 (0 933)	15069	25 0000	23 42
83 1,1,2,2-Tetrachloroetha	ne 83	11 487	11 490 (0 922)	15320	25 0000	22 04
84 1,2,3-Trichloropropane	110	11 530	11 538 (0 925)	4305	25 0000	22 73
85 4-Chlorotoluene	126	11.750	11 752 (0 943)	15507	25.0000	23 30
86 1,3,5-Trimethylbenzene	105	11 707	11 710 (0 940)	49224	25.0000	22 65
87 tert-Butylbenzene	119	12 013	12 015 (0.964)	41419	25 0000	22 31

Data File: /var/chem/hp4.i/40517d.b/1a40517.d Report Date: 17-May-2000 13:43 Page 3

					AMOUN	TS
	QUANT SIG			•	CAL-AMT	ON-COL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	(ng)
	====	==	**========	======	=======	=====
88 1,2,4-Trimethylbenzene	105	12 080	12 083 (0.970)	49237	25 0000	22 20
89 sec Butylbenzene	105	12 239	12 242 (0 982)	69676	25 0000	23 01
90 4-Isopropyltoluene	119	12 386	12 388 (0 994)	52479	25 0000	23.25
91 1,3-Dichlorobenzene	146	12 386	12 382 (0 994)	25919	25 0000	23 39
93 1,4 Dichlorobenzene	146	12.484	12 486 (1.002)	26544	25 0000	23.49
94 n-Butylbenzene	91	12 796	12 798 (1 027)	59437	25 0000	22.99
95 1,2-Dichlorobenzene	146	12.857	12 859 (1 032)	23998	25 0000	23 01
96 1,2-Dibromo-3-chloropropane	157	13.664	13 660 (1.097)	1220	25 0000	17 55
97 1,2,4-Trichlorobenzene	180	14 459	14 461 (1 160)	13080	25 0000	27 17
98 Hexachlorobutadiene	225	14 581	14 583 (1 170)	6293	25.0000	27 02
99 Naphthalene	128	14 715	14 711 (1 181)	32822	25 0000	30 47
100 1,2,3 Trichlorobenzene	180	14 966	14.962 (1 201)	14724	25 0000	38 61



Data File: /var/chem/hp4.1/40517d.b/1b40517.d

Report Date: 17-May-2000 13:44

#### STL-PITTSBURGH

Page 1

VOLATILE REPORT SW-846 Method

Data file : /var/chem/hp4.i/40517d.b/1b40517.d

Lab Smp Id: vstd20 Client Smp ID: vstd20

Inj Date : 17-MAY-2000 12:09

Operator : 16328 Smp Info : vstd20 5ml Inst ID: hp4.i

Misc Info: ,40517d.b, med1000.m

Kslalo Comment

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m Meth Date : 17-May-2000 13:44 dudeckk Quant 7 Cal Date : 17-MAY-2000 13:03 Cal Fil Quant Type: ISTD Cal File: 1e40517.d

Als bottle: 6 Calibration Sample, Level: 2

Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: 2-padep.sub

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/V1

Name	Value	Description
DF Vf Vo Vi		Dilution Factor FinalVolume Sample Volume Purge Volume

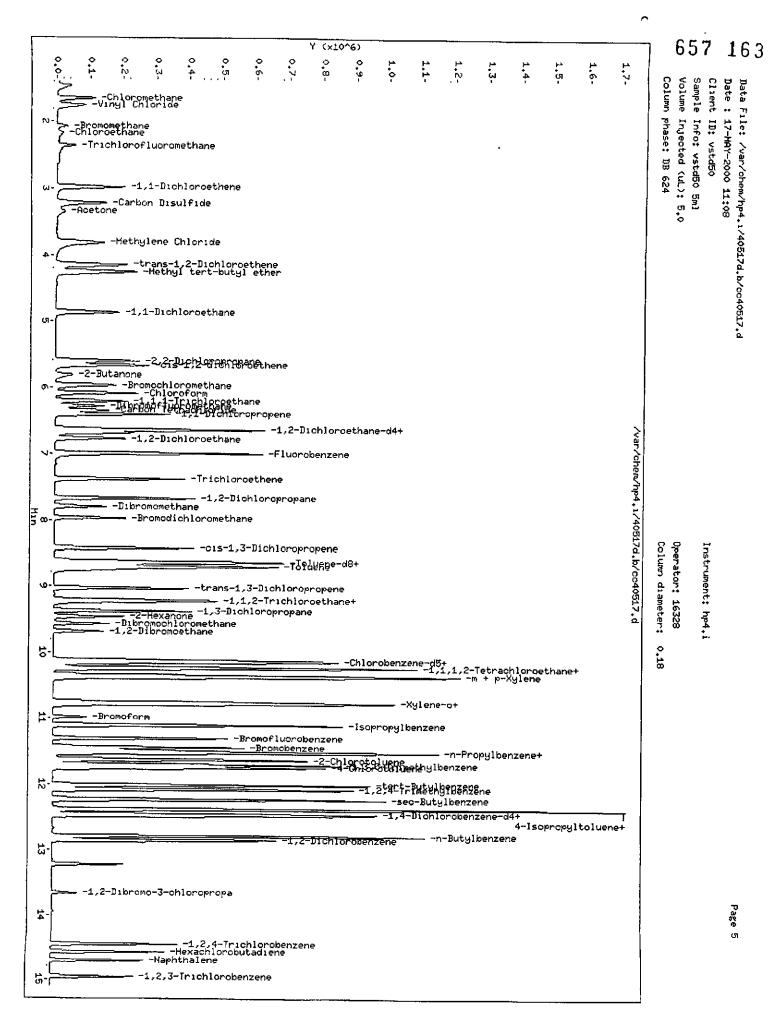
						MUOMA	TS
		QUANT SIG				CAL AMT	ON-COL
Co	abrucqmc	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
==	**************************************		==	======	두프를로로로로	=======	======
*	46 Fluorobenzene	96	6 993	6 996 (1.000)	693290		
*	69 Chlorobenzene-d5	119	10.123	10 126 (1 000)	149250		
*	92 1,4-Dichlorobenzene-d4	152	12 459	12 462 (1 000)	227595		
\$	39 Dibromofluoromethane	113	6 284	6 287 (0.899)	50380	100 000	106.2
\$	43 1,2-Dichloroethane-d4	65	6.688	6 684 (0.956)	67909	100.000	104.5
\$	59 Toluene d8	98	8.644	8 647 (0 854)	270089	100 000	107 7
\$	80 Bromofluorobenzene	95	11 309	11 312 (1.117)	102751	100 000	106 1
	1 Dichlorodifluoromethane	85	1 455	1 488 (0.208)	32883	100 000	97 29
	2 Chloromethane	50	1.644	1 647 (0 235)	74057	100.000	97 19
	3 Vinyl Chloride	62	1.742	1.769 (0 249)	68569	100 000	95 06
	4 Bromomethane	94	2 066	2 069 (0 295)	10004	100 000	95.66
	5 Chloroethane	64	2 152	2 203 (0.308)	12103	100.000	99 37
	6 Trichlorofluoromethane	101	2.353	2 350 (0 337)	15824	100 000	81 81
	12 1,1-Dichloroethene	96	2 965	2 943 (0 424)	44309	100 000	98.45
	13 Acetone	43	3 344	3 377 (0.478)	28419	100 000	99 17

Page 2

						AMOUN	me
		QUANT SIG				CAL-AMT	
Co	ompounds	MASS	RT	EXP RT REL RT	RESPONSE	· · · · · · ·	ON-COL
	异苯苯磺磺基 计连续 化氯化 计自己 化自己 化二甲基苯甲甲	E 2 2 3				( ng)	( ng)
	15 Carbon Disulfide	76	3 203	3 182 (0 458)	134764	100 000	00.70
	18 Methylene Chloride	84	3 808	3.799 (0 545)	50878	100 000	98 30
	19 trans-1,2-Dichloroethene	96	4 145	4.129 (0 593)	53357	100 000	91 91 99 13
	20 Methyl tert-butyl ether	73	4 261	4 270 (0.609)	131582	100 000	99 93
	24 1,1-Dichloroethane	63	4 866	4 857 (0 696)	107347	100 000	97 41
	27 2,2-Dichloropropane	77	5 612	5,609 (0 802)	63081	100 000	94 84
	28 cis-1,2-dichloroethene	96	5 673	5 670 (0 811)	60464	100 000	97.51
М	29 1,2-Dichloroethene (total)	100		0 0,0 (0 011,	113821	200 000	196 5
	30 Bromochloromethane	128	5 979	5.975 (0 855)	24435	100 000	99 53
	31 2-Butanone	43	5 826	5 847 (0 833)	50305	100 000	
	37 Chloroform	83	6 089	6 091 (0 871)	84073	100 000	102 1 97 56
	38 1,1,1-Trichloroethane	97	6 205	6.202 (0 B87)	62265	100.000	97.67
	40 1,1-Dichloropropene	75	6 407	6-403 (0 916)	68985	100.000	97.67
	41 Carpon Tetrachloride	117	6 345	6 342 (0 907)	36778	100 000	91 86
	42 Benzene	78	6.639	6 636 (0.949)	247867		97 88
	45 1,2-Dichloroethane	62	6 773	6 770 (0 969)	68487	100 000	97 17
	47 Trichloroethene	130	7 379	7.381 (1 055)	51621	100 000	97 01
	49 1,2 Dichloropropane	63	7 672	7 675 (1.097)	67722	100.000	96.38
	50 Dibromomethane	93	7 800	7 797 (1 115)	28403	100.000	95 82
	53 Bromodichloromethane	83	7 972	7.974 (1 140)	52518	100 000	90 64
	57 cis-1,3 Dichloropropene	75	8 436	8 439 (1 206)	82958	100 000	92 65
	58 4-Methyl 2-Pentanone	43	8 626	8 635 (0.852)	88960	100 000	97 43
	60 Toluene	91	8 711	8 714 (0 861)	265444	100 000	99 18
	61 trans-1,3-Dichloropropene	75	9 029	9.032 (0 892)	72424	100 000	95 54
	63 1,3-Dichloropropane	76	9 378	9 380 (0 926)	89640	100 000	99 59
	64 1,1,2-Trichloroethane	97	9 213	9 215 (0.910)	52125	100.000	101.0
	65 Tetrachloroethene	164	9 231	9 234 (0 912)	34098	100 000	96.32
	66 2-Hexanone	43	9 463	9-466 (0 935)	63481	100 000	96 16
	67 Dibromochloromethane	129	9 567	9 570 (0.945)	30569	100.000	88 52
	68 1,2-Dibromoethane	107	9 683	9 686 (0 957)	44571	100 000	98 86
	70 Chlorobenzene	112	10 148	10.157 (1 002)	162934	100 000	98 41
	71 1,1,1,2-Tetrachloroethane	131	10 240	10 249 (1.011)	38253	100 000	94 12
	72 Ethylbenzene	106	10 233	10 236 (1 011)	92680	100.000	98 95
	73 m + p-Xylene	106	10 356	10 359 (1 023)	226334	200 000	202 2
	74 Xylene-o	106	10 753	10.756 (1 062)	108012	100 000	96 80
М		106			334346	100.000	299 6
	76 Styrene	104	10 784	10 787 (1.065)	186878	100.000	99 58
	77 Bromoform	173	10 985	10 988 (1 085)	14349	100 000	79 36
	78 Isopropylbenzene	105	11 108	11.111 (1 097)	267060	100 000	98 85
	79 Bromobenzene	156	11 444	11 453 (0.919)	55402	100.000	94 51
	81 n-Propylbenzene	120	11.523	11.526 (0.925)	78527	100.000	98 22
	82 2 Chlorotcluene	126	11 627	11 636 (0 933)	65547	100 000	96 58
	83 1,1,2,2-Tetrachloroethane	83	11 487	11.490 (0 922)	71746	100 000	97.B6
	84 1,2,3-Trichloropropane	110	11 536	11 538 (0.926)	19847	100 000	99 36
	85 4-Chlorotoluene	126	11 750	11 752 (0 943)	69019	100 000	98 39
	86 1,3,5 Trimethylbenzene	105	11 707	11 710 (0 940)	223987	100 000	97.74
	87 tert-Butylbenzene	119	12 019	12 015 (0 965)	189222	100 000	96.65

Data File: /var/chem/hp4.1/40517d.b/1b40517.d Report Date: 17-May-2000 13:44

					AMOUN	TS
	QUANT SIG				CAL-AMT	ON COL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	(ng)
	====	= 4		========	=====	35=====
88 1,2,4 Trimethylbenzene	105	12 080	12 083 (0 970)	231554	100 000	98 99
89 sec-Butylbenzene	105	12 239	12 242 (0 982)	311341	100 000	97 51
90 4-Isopropyltoluene	119	12 385	12 388 (0 994)	235892	100 000	99.10
91 1,3-Dichlorobenzene	146	12 385	12.382 (0 994)	114443	100 000	97 93
93 1,4-Dichlorobenzene	146	12 483	12 486 (1 002)	117558	100 000	98 66
94 n-Butylbenzene	91	12 795	12 798 (1 027)	261868	100 000	96 05
95 1,2-Dichlorobenzene	146	12 856	12 859 (1 032)	106285	100 000	96 65
96 1,2 Dibromo-3-chloropropane	157	13 663	13 660 (1 097)	5822	100 000	79 40
97 1,2,4 Trichlorobenzene	180	14 458	14 461 (1 160)	45531	100 000	89 68
98 Hexachlorobutadiene	225	14 580	14 583 (1 170)	21823	100 000	88 84
99 Naphthalene	128	14 715	14 711 (1 181)	100135	100 000	88 14
100 1,2,3 Trichlorobenzene	180	14 959	14 962 (1 201)	34001	100 000	84 54



Data File: /var/chem/hp4.i/40517d.b/cc40517.d

Report Date: 17-May-2000 13:55

Page 1

Kistalo

#### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file : /var/chem/hp4.i/40517d.b/cc40517.d

Lab Smp Id: vstd50
Inj Date : 17-MAY-2000 11:08 Client Smp ID: vstd50

Operator: 16328 Inst ID: hp4.i

Smp Info : vstd50 5ml

Misc Info: ,40517d.b, med1000.m

Comment

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m

Meth Date : 17-May-2000 13:55 h Quant Type: ISTD Cal Date : 17-MAY-2000 13:03 Cal File: 1e40517.d

Als bottle: 4 Calibration Sample, Level: 3

Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: 2-padep.sub

Target Version: 3.40 Processing Host: hpuxcs21

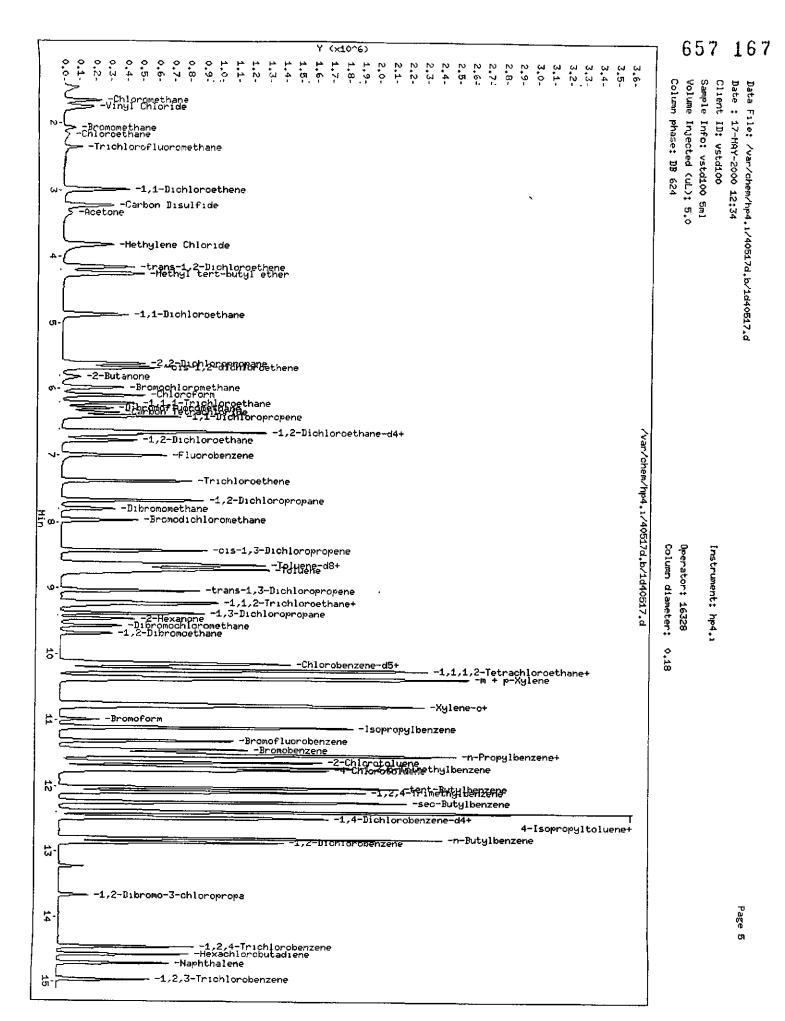
Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

Name	Value	Description
DF Vf Vo Vi	5.000 5.000	Dilution Factor FinalVolume Sample Volume Purge Volume

					AMOUN	TS
	QUANT SIG				CAL-AMT	ON-COL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
주프주주 3 11 대표 교교회 등 등 등 등 등 등 등 등 등 등 등 등 등 등 등 등 등 등	3022	==		======	<b>严强是</b> 法司告金	====###
* 46 Fluorobenzene	96	6.994	6.994 (1 000)	660677		
* 69 Chlorobenzene-d5	119	10 124	10 124 (1 000)	149779		
* 92 1,4-Dichlorobenzene-d4	152	12 460	12 460 (1 000)	227254		
\$ 39 Dibromofluoromethane	113	6.285	6 285 (0 899)	112189	250 000	248 2
\$ 43 1.2 Dichloroethane-d4	65	6 683	6 683 (0 955)	155579	250 000	251 3
\$ 59 Toluene-d8	98	8 645	8 645 (0 854)	629253	250 000	250 1
\$ 80 Bromofluorobenzene	95	11.310	11 310 (1 117)	246086	250 000	253 2
1 Dichlorodifluoromethane	85	1 450	1.450 (0 207)	84060	250 000	261 0
2 Chloromethane	50	1 645	1 645 (0.235)	189504	250 000	261 0
3 Vinyl Chloride	62	1.743	1 743 (0 249)	173382	250 000	252 2
4 Bromomethane	94	2.061	2 061 (0 295)	26679	250 000	267 7
5 Chloroethane	64	2 165	2 165 (0.309)	32667	250 000	281.4
6 Trichlorofluoromethane	101	2 354	2 354 (0.337)	54753	250 000	297 0
12 1,1-Dichloroethene	96	2 978	2 978 (0 426)	118932	250 000	277 3
13 Acetone	43	3 351	3 351 (0 479)	64935	250 000	237.8

					MOUN	ITS
0-	QUANT SIG				CAL-AMT	ON-COL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
15. On when the way 25. 1	====	==	*****	======	*****	
15 Carbon Disulfide	76	3 216	3 216 (0 460)	362443	250 000	277 4
18 Methylene Chloride	84	3 815	3 815 (0 545)	147271	250 000	279 2
19 trans-1,2-Dichloroetnene	96	4.152	4 152 (0 594)	142814	250 000	278 4
20 Methyl tert-butyl ether	73	4 249	4 249 (0 608)	334257	250 000	266 4
24 1,1-Dichloroethare	63	4 867	4 867 (0.696)	280382	250 000	267 0
27 2,2-Dichloropropane	77	5.613	5 613 (0 802)	167972	250 000	265 0
28 cis-1,2-dichloroethene	96	5.674	5 674 (0 811)	158989	250 000	269.1
M 29 1.2-Dichloroethene (total)	100			301803	500 000	546 8
30 Bromochloromethane	128	5.973	5 973 (0 854)	62236	250 000	266 0
31 2 Butanone	43	5 821	5.821 (0 832)	116124	250 000	247 2
37 Chloroform	83	6 090	6 090 (0 871)	219279	250 000	267 0
38 1.1,1-Trichloroethane	97	6 212	6 212 (0 888)	162494	250.000	267 5
40 1,1-Dichloropropene	75	6 407	6.407 (0 916)	181812	250.000	269 0
41 Carbon Tetrachloride	117	6 352	6 352 (0 908)	105078	250 000	275 4
42 Benzene	78	6 640	6 640 (0 949)	646947	250 000	268 1
45 1,2-Dichloroethane	62	6 774	6 774 (0.969)	181162	250 000	269 7
47 Trichloroethene	130	7 379	7 379 (1 055)	136342	250 000	268 9
49 1,2-Dichloropropane	63	7 673	7 673 (1 097)	179887	250 000	268 6
50 Dibromomethane	93	7 795	7 795 (1 114)	78170	250 000	276 7
53 Bromodichloromethane	83	7.972	7 972 (1 140)	152326	250 000	275 9
57 cis-1,3-Dichloropropene	75	B 437	8 437 (1 206)	234977	250 000	275 4
58 4 Methyl-2-Pentanone	43	8 627	8 627 (0 852)	234014	250.000	255 4
60 Toluene	91	8.712	8 712 (0 861)	718952	250 000	267 7
61 trans 1,3-Dichloropropene	75	9 030	9 030 (0 892)	206628	250 000	271 6
63 1,3-Dicaloropropane	76	9 379	9 379 (0 926)	240837	250 000	266 6
64 1,1,2-Trichloroethane	97	9 207	9 207 (0 909)	137716	250 000	265.8
65 Tetrachloroethene	164	9.232	9.232 (0 912)	94563	250 000	266 2
66 2-Hexanone	43	9 458	9 458 (0 934)	163100	250 000	246 2
67 Dibromochloromethane	129	9 568	9 568 (0 945)	94873	250 000	273 8
68 1,2-Dibromoethane	107	9 678	9 678 (0.956)	121424	250 000	268 4
70 Chlorobenzene	112	10 149	10 149 (1 002)	442746	250 000	266 5
71 1,1,1,2-Tetrachloroethane	131	10 247	10 247 (1 012)	112857	250 000	276.7
72 Ethylbenzene	106	10 234	10 234 (1.011)	252651	250 000	268 8
73 m + p-Xylene	106	10 363	10 363 (1 024)	611718	500 000	544 6
74 Xylene-o	106	10 754	10 754 (1 062)	300199	250 000	268 1
M 75 Xylenes (total)	106		, , , ,	911917	250 000	814 4
76 Styrene	104	10 785	10 785 (1 065)	519441	250.000	275 8
77 Bromoform	173	10 986	10 986 (1 085)	50250	250.000	
78 Isopropylbenzene	105	11 109	11.109 (1 097)	740920	250 000	277 0
79 Bromobenzene	156	11 445	11 445 (0 919)	153219		273 3
81 n-Propylbenzene	120	11 524	11 524 (0 925)	211369	250.000	261 8
82 2-Chlorotoluene	126	11 634	11.634 (0 934)	179841	250.000	264.8
83 1,1,2,2-Tetrachloroethane	83	11 488	11 488 (0 922)	196523	250 000	265 4
84 1,2,3-Trichloropropane	110	11 537	11 537 (0.926)	53924	250 000	268 4
85 4-Chlorotoluene	126		11 750 (0 943)		250 000	270.4
86 1,3,5-Trimetnylbenzene	105		11.708 (0.940)	188006	250 000	268 3
87 tert-Butylbenzene	119	12 013	12 013 (0 964)	606790	250 000	265 2
4	-17	10 013	12 013 (0 964)	521697	250 000	266.9

						AMOUN	TS
		QUANT SIG				CAL-AMT	ON - COL
Сотро	unds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
=====		====	==	****** ======	<b>5</b> #=====	======	
88	1,2,4-Trimethylbenzene	105	12 081	12 081 (0 970)	623203	250 000	266 8
89	sec-Butylbenzene	105	12 233	12 233 (0.982)	849111	250 000	266 3
90	4-Isopropyltoluene	119	12 386	12 386 (0 994)	654793	250 000	275 5
91	1,3 Dichlorobenzene	145	12 386	12.386 (0 994)	316567	250.000	271 3
93	1,4-Dichlorobenzene	146	12.484	12.484 (1 002)	312890	250 000	263 0
94	n Butylbenzere	91	12 796	12 796 (1 027)	713004	250 000	261 9
95	1,2-Dichlorobenzene	146	12 857	12 857 (1 032)	290497	250.000	264 6
96	1,2-Dibromo 3-chloropropane	157	13 664	13.664 (1 097)	18343	250 000	250 5
97	1,2,4-Trichlorobenzene	180	14 459	14.459 (1 160)	117260	250 000	231 3
98	Hexachlorobutadiene	225	14 581	14 581 (1.170)	55149	250.000	224 8
99	Naphthalene	128	14.709	14 709 (1 181)	241145	250.000	212.6
100	1.2.3-Trichlorobenzene	180	14 960	14 960 (1 201)	81058	250 000	201 8



Page 1

Data File: /var/chem/hp4.1/40517d.b/1d40517.d

Report Date: 17-May-2000 13:44

#### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file : /var/chem/hp4.i/40517d.b/1d40517.d

Lab Smp Id: vstd100 Client Smp ID: vstd100

Inj Date : 17-MAY-2000 12:34

Operator : 16328 Inst ID: hp4.i

Smp Info : vstd100 5ml

Misc Info: ,40517d.b, med1000.m

Kislinla Comment

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m

Meth Date: 17-May-2000 13:44 dudeckk Quant Type: ISTD Cal File: 1e40517.d Cal Date : 17-MAY-2000 13:03

Als bottle: 7 Calibration Sample, Level: 4

Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: 2-padep.sub

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

Name	Value	Description
DF Vf Vo Vi	5.000 5.000	Dilution Factor FinalVolume Sample Volume Purge Volume

						AMOUN	TS
		QUANT SIG				CAL-AMT	ON-COL
C	ompounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
=:	===##==================================		21			======	
*	46 Fluorobenzene	96	6.997	6.996 (1 000)	723860		
*	69 Chlorobenzene-d5	119	10 127	10 126 (1 000)	161827		
*	92 1,4-Dichlorobenzene-d4	152	12 456	12 462 (1 000)	244057		
\$	39 Dibromofluoromethane	113	6 288	6 287 (0.899)	241358	500.000	487 3
\$	43 1,2-Dichloroethane-d4	65	6 685	6.684 (0 955)	328172	500 000	483.8
\$	59 Toluene-d8	98	8 648	8.647 (0 854)	1316738	500 000	484 4
\$	80 Bromofluorobenzene	95	11 307	11 312 (1 117)	520956	500 000	496 1
	1 Dichlorodifluoromethane	85	1.452	1 488 (0.208)	186389	500 000	528.2
	2 Chloromethane	50	1.648	1 647 (0 235)	414921	500 000	521 5
	3 Vinyl Chloride	62	1 746	1 769 (0 249)	399176	500 000	530 0
	4 Bromomethane	94	2 064	2 069 (0.295)	58716	500 000	537.8
	5 Chloroethane	64	2 167	2 203 (0 310)	65977	500 000	518 8
	6 Trichlorofluoromethane	101	2 357	2 350 (0 337)	115846	500 000	573 6
	12 1 1 Dichloroethene	96	2 987	2 943 (0 427)	246289	500 000	524.1
	13 Acetone	43	3 341	3 377 (0 478)	126974	500 000	424.4



					AMOUN	<b>Y</b> TS
	QUANT SIG				CAL-AMT	ON-COL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	(ng)	(ng)
	23 F R	24 74	****	==n====	======	3555855
15 Carbon Disulfide	76	3 225	3 182 (0.461)	773680	500 000	540 5
18 Methylene Chloride	84	3 818	3 799 (0 546)	303463	500.000	525 0
19 trans-1,2-Dichloroethene	96	4 154	4 129 (0 594)	295490	500 000	525 8
20 Methyl tert butyl ether	73	4 252	4 270 (0 608)	724089	500 000	526 7
24 1,1-Dichloroethane	63	4 870	4 857 (0 696)	602338	500 000	523 5
27 2,2-Dichloropropane	77	5 615	5 609 (0 803)	378464	500.000	545 0
28 cis 1,2-dichloroethene	96	5 677	5 670 (0 811)	342945	500 000	529 7
M 29 1,2-Dichloroethene (total)	100			638435	1000 00	1056
30 Bromochloromethane	128	5 976	5 975 (0 854)	135466	500.000	528 5
31 2-Butanone	43	5 817	5.847 (0 831)	241800	500 000	469 8
37 Chloroform	83	6 086	6 091 (0 870)	469507	500 000	521 8
38 1,1,1-Trichloroethane	97	6 208	6 202 (0 887)	357218	500 000	536 7
40 1.1-Dichloropropene	75	6.410	6 403 (0 916)	396743	500 000	535 8
41 Carbon Tetrachloride	117	6 349	6 342 (0 907)	234217	500 000	560 3
42 Benzene	78	6 642	6 636 (0 949)	1386243	500 000	524 3
45 1.2-Dichloroethane	62	6 771	6 770 (0 968)	388818	500 000	528.4
47 Trichloroethene	130	7.382	7 381 (1 055)	298384	500 000	537 1
49 1,2-Dichloropropane	63	7 669	7 675 (1 096)	389801	500 000	531 3
50 Dibromomethane	93	7 798	7 797 (1.114)	163499	500.000	528 3
53 Bromodichloromethane	83	7 969	7 974 (1 139)	339420	500.000	561 0
57 cis-1,3-Dichloropropene	75	8 434	8 439 (1 205)	518893	500 000	555 0
58 4-Methyl-2-Pentanone	43	8 623	8 635 (0 851)	499267	500 000	504 3
60 Toluene	91	8.715	8 714 (0 861)	1523920	500 000	525 2
61 trans-1,3-Dichloropropene	75	9.033	9 032 (0 892)	458964	500 000	558 4
63 1,3-Dichloropropane	76	9 375	9 380 (0 926)	514213	500 000	526 9
64 1,1,2-Trichloroethane	97	9 210	9 215 (0 909)	295592	500 000	528.0
65 Tetrachloroethene	164	9 234	9 234 (0 912)	205773	500 000	536.1
66 2 Hexanone	43	9 461	9.466 (0 934)	357162	500 000	499 0
67 Dibromochloromethane	129	9 565	9 570 (0.944)	218270	500 000	582 9
68 1,2-Dibromoethane	107	9 681	9 686 (0 956)	264181	500 000	540.4
70 Chlorobenzene	112	10 151	10.157 (1 002)	944941	500 000	526 4
71 1,1,1,2-Tetrachloroethane	131	10 243	10 249 (1 011)	246692	500 000	559 8
72 Ethylbenzene 73 m + p Xylene	106	10 237	10 236 (1.011)	529388	500 000	521.3
• •	106	10 365	10 359 (1 024)	1288852	1000 00	1062
74 Xylene-o M 75 Xylenes (total)	106	10 757	10 756 (1 062)	656796	500 000	542 9
76 Styrene	106		(;)	1945648	500 000	1608
77 Bromoform	104	10 787	10 787 (1 065)	1119514	500 000	550 2
78 Isopropylbenzene	173	10 983	10 988 (1 085)	122056	500 000	622 6
79 Bromobenzene	105	11.111	11 111 (1 097)	1586989	500 000	541 8
81 n-Propylbenzene	156 120	11.448	11 453 (0 919)	333534	500 000	530 6
82 2-Chlorotoluene	126	11 521	11 526 (0.925)	451688	500.000	526 9
83 1,1,2,2-Tetrachloroethane	83	11 631	11 636 (0.934)	379255	500 000	521 1
84 1,2,3-Trichloropropane	110	11.490 11 533	11.490 (0 922)	418120	500 000	531 8
85 4-Chlorotoluene	126	11 533	11 538 (0 926)	112580	500 000	525.6
86 1,3,5-Trimethylbenzene	105	11.710	11 752 (0 944) 11 710 (0 940)	394786	500 000	524.6
87 tert-Butylbenzene	119	12 016	12 015 (0 965)	1308416	500 000	532 4
	***	15 010	TO (C05)	1126398	500 000	536 5

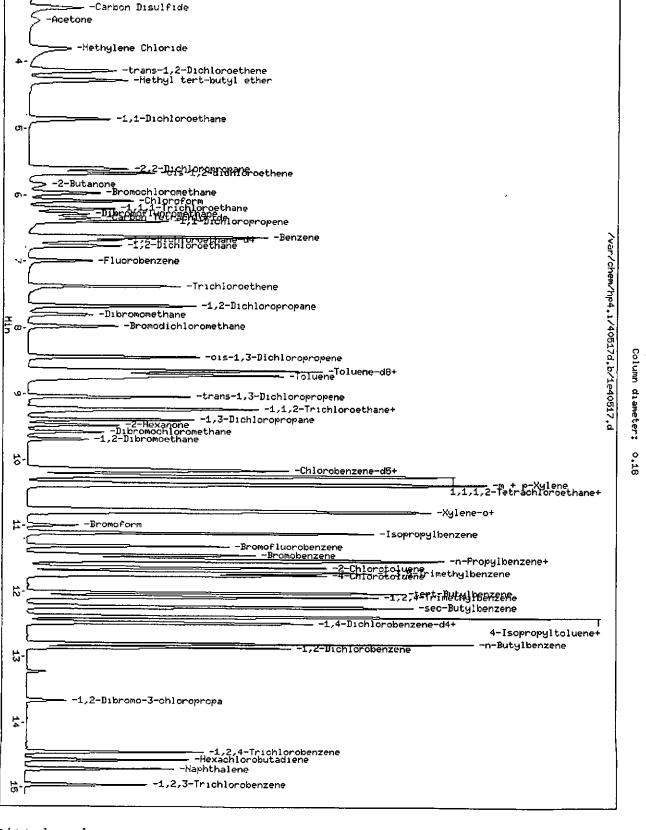
# 657 170

						AMOUN	TS
		QUANT SIG				CAL-AMT	ON-COL
Compo	unds	REAM	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
	=====###	====	==	###### ######	3======	======	
88	1,2,4 Trimethylbenzene	105	12 083	12 083 (0 970,	1339260	500 000	533 9
89	sec-Butylbenzene	105	12 236	12 242 (0 982)	1818067	500 000	531 0
90	4-Isopropyltoluene	119	12 383	12 388 (0.994)	1369572	500 000	536 5
91	1,3-Dichlorobenzene	146	12.383	12 382 (0 994)	671612	500 000	535 9
93	1,4-Dichlorobenzene	146	12.481	12.486 (1 002)	665773	500 000	521 0
94	n Butylbenzene	91	12 792	12 798 (1 027)	1562308	500 000	534 4
95	1,2-Dichlorobenzene	146	12 854	12 859 (1 032)	618380	500 000	524 4
96	1,2-Dibromo-3-chloropropane	157	13 661	13 660 (1 097)	45746	500.000	581 8
97	1,2,4-Trichlorobenzene	180	14 461	14 461 (1 161)	2 <b>7</b> 7052	500 000	508 9
98	Hexachlorobutadiene	225	14 578	14 583 (1 170)	134819	500 000	511.8
99	Naphthalene	129	14 712	14 721 (1 181)	596388	500.000	489.5
100	1,2,3-Trichlorobenzene	180	14.957	14.962 (1 201)	185090	500 000	429 2

Data File: /var/chem/hp4.i/40517d.b/le40517.d

5,6-





(x10^6) 2.8-3.0-

٠<u>,</u>8

== -Chloromethane -Vinyl Chloride

- −Bromomethane -Chloroethane - −Trichlorofluoromethane

-1,1-Dichloroethene

Data File: /var/chem/hp4.i/40517d.b/le40517.d

Report Date: 17-May-2000 13:44

#### STL-PITTSBURGH

Page 1

VOLATILE REPORT SW-846 Method

Data file : /var/chem/hp4.1/40517d.b/1e40517.d

Lab Smp Id: vstd200 Client Smp ID: vstd200

Inj Date : 17-MAY-2000 13:03
Operator : 16328 Inst ID: hp4.i

Smp Info : vstd200 5ml

Misc Info : ,40517d.b, med1000.m Ks/17/-Comment

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m

Meth Date: 17-May-2000 13:44 dudeckk Cal Date: 17-MAY-2000 13:03 Quant Type: ISTD Cal File: 1e40517.d

Als bottle: 8 Calibration Sample, Level: 5

Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: 2-padep.sub

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

Name	Value	Description
DF Vf Vo Vi		Dilution Factor FinalVolume Sample Volume Purge Volume

						AMOUN	rs
		QUANT SIG				CAL AMT	ON-COL
(	Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
•	본·소요요요요요요요요요요요요요요요요요요	====	==	<b>7388</b> 03 00000	BB=======	=====	========
4	* 46 Fluorobenzene	96	6 996	6 996 (1 000)	771765		
1	* 69 Chlorobenzene-d5	119	10 126	10 126 (1 000)	174678		
•	* 92 1,4-Dichlorobenzene-d4	152	12.462	12.462 (1 000)	245636		
Ş	\$ 39_Dibromofluoromethane	113	6 287	6 287 (0 899)	448691	1000 00	849 7
\$	\$ 43 1,2 Dichloroethane d4	65	6 684	6 684 (0.955)	615666	1000 00	851 4
\$	\$ 59 Toluene-dB	98	8.647	8 647 (0.854)	2416455	1000.00	823 5
\$	\$ 80 Bromofluorobenzene	95	11.312	11 312 (1 117)	960718	1000 00	847 5
	1 Dichlorodifluoromethane	85	1 488	1.488 (0 213)	374065	1000 00	994 2
	2 Chloromethane	50	1 647	1.647 (0.235)	627841	1000 00	976 0
	3 Vinyl Chloride	62	1 769	1 769 (0.253)	806063	1000 00	1004
	4 Bromomethane	94	2 069	2.069 (0 296)	109766	1000 00	942 9
	5 Chloroethane	64	2 203	2 203 (0 315)	121983	1000 00	899 7
	6 Trichlorofluoromethane	101	2 350	2 350 (0 336)	209443	1000 00	972 7
	12 1,1 Dichloroethene	96	2 943	2 943 (0 421)	444848	1000 00	887 9
	13 Acetone	43	3 377	3 377 (0 483)	350752	1000 00	1100

CAL-MR   MASS						AMOUN	ITS
15 Carbon Disulfide 76 3 182 3 182 (0 455) 1397397 1000 00 973 2 18 Methylene Chlorade 84 1399 3.799 5.799 (5 431) 538099 1000 00 973 2 19 trans-1.2 highloroethene 96 4 129 4.129 (0 500) 575254 1000 00 960 12 12 0 Methyl tetr-butyl éther 73 4 479 4 270 (0 610) 1320299 1000 00 1005 24 1.1-Dichloroethane 63 4 887 4 857 (0 684) 1233229 1000 00 1005 24 1.1-Dichloroethane 63 4 887 4 857 (0 684) 1233229 1000 00 1005 27 2.2-Dichloroethane 66 5 650 5 650 (0 802) 797232 1000 00 1005 27 2.2-Dichloroethane 76 5 659 5 650 (0 802) 797232 1000 00 1005 27 2.2-Dichloroethane 128 5 975 5.975 (0 884) 1233229 1000 00 1005 27 2.2-Dichloroethane 128 5 975 5.975 (0 884) 1233229 1000 00 1002 28 1.2-Dichloroethane 128 5 975 5.975 (0 884) 1233229 1000 00 1002 28 1.2-Dichloroethane 128 5 975 5.975 (0 884) 1233229 1000 00 1002 28 12 2-Dichloroethane 128 5 975 5.975 (0 884) 1239797 1000 00 1002 28 12 2-Dichloroethane 128 5 975 5.975 (0 884) 1239797 1000 00 1002 28 12 2-Dichloroethane 128 5 975 5.975 (0 884) 1239797 1000 00 1004 40 1.1 Dichloropropene 75 6 403 6 403 (0 915) 788055 1000 00 999.6 41 Carbon Tetrachloride 117 6 342 5 342 (0 906) 465056 1000 00 999.4 42 Bennene 78 6 6536 6.636 (0 948) 2803750 1000 00 1002 42 Bennene 78 6 6536 6.636 (0 948) 2803750 1000 00 1005 47 Tetrachloroethane 130 7 381 7 381 (1.055) 507270 1000 00 1005 47 Tetrachloroethane 130 7 381 7 381 (1.055) 507270 1000 00 1005 47 Tetrachloroethane 130 7 381 7 381 (1.055) 507270 1000 00 1005 47 Tetrachloroethane 130 7 394 7 794 (1 140) 711520 1000 00 1005 47 Tetrachloroethane 130 7 394 7 394 (1 140) 711520 1000 00 1005 47 Tetrachloroethane 131 7 394 7 394 (1 140) 711520 1000 00 1005 50 Dibromomethane 130 7 394 7 394 (1 140) 711520 1000 00 1005 50 Dibromomethane 131 7 394 7 394 (1 140) 711520 1000 00 1007 58 44 Methyl-2-Pentanoethane 131 7 394 7 394 (1 140) 711520 1000 00 1007 58 44 Methyl-2-Pentanoethane 131 7 394 7 394 (1 140) 711520 1000 00 1007 58 44 Methyl-2-Pentanoethane 132 9 350 1005 1009 1009 1000 1007 58 44 Methyl-2-Pentanoethane 140 9 686 9 686 (0 957) 52539						CAL-AMT	ON-COL
18 Carbon Dusulfide	-	MASS	ŔT	EXP RT REL RT	RESPONSE	(ng)	( ng)
18 Methylene Chiorade				#32222 HERRE	****	#4====#	======
19   Crans-1.2 Dichloroethene				3 182 (0 455)	1397397	1000 00	915 7
20 Methyl terr-butyl éther	· •		3 799	3.799 (0 543)	538099	1000 00	873 2
24 1,1-Dachloroethane 63 4 857 4 857 (0 694) 1233229 1000 00 1005 27 2,2-Dachloropropane 77 5 609 5 609 (0 802) 279232 1000 00 1007 28 cis-1,2-dachloroethene 96 5 670 6 509 609 (0 802) 279232 1000 00 1007 30 Bromochloromethane 123 5 975 5.975 (0 854) 279197 1000 00 1052 31 2-Datanone 43 5 847 6 840 5 511183 1000 00 991 6 37 Chloroform 83 6 691 6 091 (0 871) 295258 1000 00 1914 37 Chloroform 83 6 691 6 091 (0 871) 295258 1000 00 1004 40 1,1 Dachloropropene 75 6 403 6 403 (0 915) 788065 1000 00 1004 40 1,1 Dachloropropene 75 6 403 6 403 (0 915) 788065 1000 00 1005 41 Carbon Tetrachloride 117 6 342 6 342 (0 906) 486506 1000 00 1005 42 Bensene 78 6 6536 6 6636 (0 948) 2803750 1000 00 1005 47 Trichloroethane 62 6 670 6 770 (0 568) 788325 1000 00 1005 47 Trichloroethane 63 7 675 7 675 (1.097) 787044 1000 00 1006 50 Obbromomethane 93 7 797 7 797 (1 114) 328434 1000 00 1006 50 Obbromomethane 83 7 797 7 797 (1 114) 328434 1000 00 1006 50 Obbromomethane 83 7 797 7 797 (1 114) 328434 1000 00 1006 50 Obbromomethane 83 7 797 7 797 (1 114) 328434 1000 00 1006 50 Obbromomethane 83 7 797 7 797 (1 114) 328434 1000 00 1006 50 Obbromomethane 83 7 797 7 797 (1 114) 328434 1000 00 1006 50 Obbromomethane 83 7 797 8 8 794 (1 1400 711520 1000 00 1006 50 Obbromomethane 83 7 797 8 8 794 (1 1400 711520 1000 00 1006 50 Obbromomethane 84 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8 8			4 129	4.129 (0 590)	575254	1000 00	960 1
27 2,2-Dichloropropane	· · · · · · · · · · · · · · · · · · ·		4 270	4 270 (0 610)	1520929	1000 00	1038
8 cis-1,2-dichloroethene 96 5 670 (0 810) 684419 1000 00 991 6  8 29 1,2 Dichloroethene (tocal) 100 1025 31 2. Euchanone 128 5 975 (0 854) 279197 1000 00 1934 6 30 Bromochloromethane 128 5 975 5 975 (0 854) 279197 1000 00 1022 31 2. Euchanone 43 5 847 5 847 (0 836) 511183 1000 00 991 6 37 Chloroform 83 6 091 6 091 (0 871) 395258 1000 00 999 3 31 1.1.1-Trichloroethane 97 6 202 6 202 (0 886) 740791 1000 00 1044 60 1.1 Dichloropropene 75 6 401 6 403 (0 915) 788065 1000 00 998.2 61 1.2 Euchanone 75 6 636 (6 666) 60 980 998.2 61 1.2 Euchanone 78 6 636 (6 666) 60 980 998.2 61 1.2 Euchanone 78 6 636 (6 666) 60 980 998.2 61 1.2 Euchanone 78 6 636 (6 666) 60 980 998.2 61 1.2 Euchanone 78 6 636 (6 666) 60 980 00 1005 60 Dibromomethane 130 7 381 7 381 (1.055) 607870 1000 00 1026 60 Dibromomethane 93 7 797 7 797 (1 114) 328434 1000.00 1007 60 Dibromomethane 93 7 797 7 797 (1 114) 328434 1000.00 995 4 6 53 Bromodichloromethane 83 7 974 7 974 (1 140) 711520 1000 00 1007 60 Dibromomethane 83 7 974 7 974 (1 140) 711520 1000 00 1007 60 Dibromomethane 83 7 974 7 974 (1 140) 711520 1000 00 1007 60 Dibromomethane 83 7 974 7 974 (1 140) 711520 1000 00 1007 60 Dibromomethane 83 7 974 7 974 (1 140) 711520 1000 00 1007 60 Dibromomethane 83 7 974 7 974 (1 140) 711520 1000 00 1007 60 00 1007 60 00 00 00 00 00 00 00 00 00 00 00 00		_		4 857 (0 694)	1233229	1000 00	1005
W 29 1.2 Dichloroethene (total)   100   128   5 975   5.975 (0 854)   279197   1000 00   1954   1000 1000   1012   1124   1124   1128	• •		5 609	5 609 (0 802)	797232	1000 00	1077
1996   1997   1000 00   1022   12 - 1997   1000 00   1022   12 - 1997   1000 00   1022   13   2 - 1997   1000 00   1022   13   2 - 1997   1000 00   1023   13   13   15   15   15   15   15   1			5 670	5 670 (0 810)	684419	1000 00	991 6
31 2-Butanone	, , , , , , , , , , , , , , , , , , , ,				1259673	2000 00	1954
37 Chloroform 83 6 091 6 091 (0 871; 959258 1000 00 999 93 81 1.1.1-Trichloroethane 97 6 202 6 202 (0 886) 740791 1000 00 1044 40 1.1 Dichloropropene 75 6 403 6 403 (0 915) 788055 1000 00 998.2 41 Carbon Tetrachloride 117 6 342 6 342 (0 906) 446506 1000 00 998.2 42 Benzene 78 6 636 6 636 (0 948) 2803750 1000 00 1005 42 Benzene 78 6 636 6 636 (0 948) 2803750 1000 00 1005 42 Benzene 130 7381 7381 (1.055) 607870 1000 00 1005 49 1.2 Dichloroethane 62 6 770 6 770 (0 968) 788825 1000 00 1005 49 1.2 Dichloropropane 63 7 675 7 675 (1.097) 787044 1000 00 1006 50 Dibromomethane 93 7 797 7 797 (1 114) 328434 1000.00 995 4 53 Bromodichloromethane 83 7 974 7 974 (1 140) 711520 1000 00 1107 55 618-1,3-Dichloropropane 75 8 439 8 439 (1.206) 1073632 1000 00 1077 58 4 Methyl-2-Pentanone 43 8.635 8 635 (0 853) 1041651 1000.00 974 6 66 Toluen 99 8 71 8 714 8 714 (0 861) 302024 1000 00 984 6 61 trans-1,3-Dichloropropane 75 9 032 9 032 (0 892) 945299 1000 00 1065 63 1.3 Dichloropropane 76 9.280 9 380 (0 926) 1034970 1000 00 982 5 64 1,12-Trichloroethane 91 9.570 9 9.215 9 215 (0 910) 60359 1000 00 983 6 65 Tetrachloroethane 104 9 234 9 234 (0 912) 421277 1000 00 983 6 65 Tetrachloroethane 107 9 686 9 686 (0 935) 728700 1000 00 983 6 65 Tetrachloroethane 107 9 686 9 686 (0 935) 728700 1000 00 983 6 65 Tetrachloroethane 107 9 686 9 686 (0 935) 728700 1000 00 983 7 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 983 5 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 983 5 70 Chlorobenzene 106 10 236 10 236 (1 011) 1042947 1000 00 993 6 77 78 77 97 11 1.1.2-Tetrachloroethane 107 9 686 9 686 (0 935) 129397 1000 00 993 6 70 Chlorobenzene 106 10 236 10 236 (1 011) 1042947 1000 00 993 6 77 78 11 1.1.1.2-Tetrachloroethane 107 9 686 9 686 (0 935) 129397 1000 00 993 6 70 Chlorobenzene 106 10 236 10 236 (1 011) 1042947 1000 00 993 6 70 Chlorobenzene 106 10 236 10 236 (1 011) 1042947 1000 00 993 6 70 Chlorobenzene 107 10 10 10 10 10 10 10 10 10 10 10 10 10		128	5 975	5.975 (0 854)	279197	1000 00	1022
38   1.1.1-Trichloroethane		43	5 847	5 847 (0 836)	511183	1000 00	931 6
40 1,1 Dichloropropene 75 6 403 6 403 (0 915) 788065 1000 00 998.2 41 Carbon Tetrachloride 117 6 342 6 342 (0 906) 486506 1000 00 1092 42 Benrene 78 6 636 6 636 (0 948) 2803750 1000 00 1092 42 Benrene 62 6 770 6 770 (0 966) 788825 1000 00 1005 47 Trichloroethane 62 6 770 6 770 (0 966) 788825 1000 00 1005 47 Trichloroethane 130 7 381 7 381 (1.055) 607870 1000 00 1026 49 1.2 Dichloropropane 63 7 675 7 675 (1.097) 787044 1000 00 1006 50 Dibromomethane 93 7 797 7 797 (1 114) 328434 1000.00 995 4 53 Bromodichloromethane 83 7 797 7 797 (1 114) 328434 1000.00 995 4 53 Bromodichloromethane 83 7 974 7 974 (1 140) 711520 1000 00 1077 58 4 Mcthyl-2-Pentanone 43 8.635 8 635 (8 853) 1041451 1000.00 974 6 60 Toluene 91 8 714 8 714 (0 861) 3022024 1000 00 964 8 61 trans-1,3-Dichloropropane 75 9.032 9032 (0 892) 945299 1000 00 1065 63 1,3 Dichloropropane 76 9.380 9 380 (0 926) 1034970 1000 00 982 6 63 1,3 Dichloropropane 76 9.380 9 380 (0 926) 1034970 1000 00 982 6 65 Tetrachloroethane 97 9.215 9 215 (0 910) 600359 1000 00 993 6 65 Tetrachloroethane 129 9.570 9 570 (0 945) 465815 1000 00 1157 66 1,2-Dibromochlane 129 9.570 9 570 (0 945) 465815 1000 00 1156 62 -Rexanne 107 9 686 9 686 (0 957) 525391 1000 00 1157 66 1,2-Dibromochane 129 9.570 9 570 (0 945) 445815 1000 00 1157 66 1,2-Dibromochane 129 9.570 9 570 (0 945) 445814 1000 00 983 5 70 70 (0 945) 445814 1000 00 9		83	6 091	6 091 (0 871)	959258	1000 00	999 9
41 Carbon Tetrachloride 117 6 342 6 342 (0 906) 486506 1000 00 1092 42 Benxene 78 6 636 6.636 (0 948) 2803750 1000 00 994 6 45 1,2-Dichloroethane 62 6 770 (0 968) 788825 1000 00 1005 47 Trichloroethane 130 7 381 7 381 (1.055) 607870 1000 00 1026 49 1,2 Dichloropropane 63 7 675 7 675 (1.097) 787044 1000 00 1006 50 Dibromomethane 93 7 797 7 797 (1 114) 328434 1000.00 995 4 53 Bromodichloromethane 83 7 974 7 974 (1 140) 711520 1000 00 1077 58 4 Mcthyl-2-Pentanone 43 8.635 8 635 (0 853) 1041451 1000.00 974 6 60 Toluene 91 8 714 (0 861) 3022024 1000 00 964 8 61 trans-1,3-Dichloropropene 75 9 012 9012 (0 892) 945299 1000 00 964 8 61 trans-1,3-Dichloropropene 75 9 012 9012 (0 892) 945299 1000 00 992 5 64 1,1,2-Trichloropropene 75 9 012 9012 (0 892) 945299 1000 00 993 6 6 70 toluene 164 9 214 9 234 9 234 (0 912) 421277 1000 00 1017 66 2 -Mexanone 164 9 2.18 9 215 (0 910) 600359 1000 00 993 6 6 70 toluene 164 9 214 9 234 9 234 (0 912) 421277 1000 00 1017 66 2 -Mexanone 164 9 2.18 9 2.18 (0 912) 421277 1000 00 1017 66 2 -Mexanone 164 9 2.14 9 234 9 234 (0 912) 421277 1000 00 1017 66 2 -Mexanone 179 9 686 9 686 (0 935) 728700 1000 00 993 6 6 1,2-Dibromochloromethane 129 9 .570 9 570 (0 945) 455815 1000 00 1152 68 1,2-Dibromochloromethane 107 9 686 9 686 (0 957) 525331 1000 00 993 70 Chlorobenaene 112 10 157 10 157 (1 003) 1905796 1000 00 993 70 Chlorobenaene 112 10 157 10 157 (1 003) 1905796 1000 00 993 70 Chlorobenaene 112 10 157 10 157 (1 003) 1905796 1000 00 993 6 11 1,1,1,1,2-Tetrachloroethane 106 10 236 10.236 (1 011) 1042947 1000 00 993 6 17 1,1,1,1,2-Tetrachloroethane 107 10 988 10 988 (1 081) 2451464 2000 00 1034 124 124 124 124 124 124 124 124 124 12		97	6 202	6 202 (0 886)	740791	1000 00	1044
42 Benzene 78 6 636 6 636 6 0948) 2803750 1000 00 994 6 45 1,2-Dichloroethane 62 6 770 6 770 (0 968) 788825 1000 00 1005 47 Trichloroethene 130 7 381 7 381 (1.055) 607870 1000 00 1026 49 1,2 Dichloropropane 63 7 675 7 675 (1.097) 787044 1000 00 1006 50 Dibromomethane 93 7 797 7 797 (1 114) 328434 1000 00 995 4 53 Bromodichloromethane 63 7 974 7 974 (1 140) 711520 1000 00 1103 57 Cls-1,3-Dichloropropene 75 8 439 8 439 (1.206) 1073632 1000 00 1077 58 4 Mcthyl-2-Pentanone 43 8.635 8 635 (0 853) 1041451 1000 00 994 8 61 trans-1,3-Dichloropropene 75 9 032 9 032 (0 892) 945299 1000 00 1065 63 1,3 Dichloropropane 76 9.380 9 380 (0 926) 1034971 1000 00 982 8 64 1,1,2-Trichloroethane 97 9.215 9 215 (0 910) 600359 1000 00 993 8 65 Tetrachloroethene 164 9 234 9 234 (0 912) 421277 1000 00 1017 66 2-Bexanone 43 9.466 9 466 (0 935) 728700 1000 00 995 7 67 Dibromomethane 129 9.570 9 570 (0 945) 465815 1000 00 995 7 68 1,2-Dibromochloromethane 129 9.570 9 570 (0 945) 465815 1000 00 995 7 68 1,1,1,2-Tetrachloroethane 129 9.570 9 570 (0 945) 465815 1000 00 995 7 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000 00 995 7 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000 00 995 7 71 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 1034 73 m + p-Xylene 106 10 359 10 359 (1 021) 104947 1000 00 1034 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6 78 Bromoform 173 10 988 10 988 (1 085) 216848 1000 00 993 6 79 Bromoform 173 10 988 10 988 (1 085) 262915 1000 00 993 7 87 Sylenes (total) 106 11 11 11 11 11 11 11 (1.097) 3083472 1000 00 993 8 83 1,2,2,2 Tetrachloroethane 136 11 536 (0 934) 727962 1000 00 993 8 84 1,2,3-Trachloropethane 136 11 538 (1 5092) 793112 1000 00 993 8 84 1,2,3-Trachloropethane 136 11 538 (1 5092) 793112 1000 00 993 8 85 4-Chlorotoluene 126 11 536 (1 5094) 9400 940 940 940 940 940 940 940 940 94	· ·	75	6 403	6 403 (0 915)	788065	1000 00	998.2
45 1,2-Dichloroethane 62 6 770 (6 968) 788825 1000 00 1005 47 Trichloroethene 130 7 381 7 381 (1.055) 607870 1000 00 1026 49 1,2 Dichloropropane 63 7 675 7 675 (1.057) 787044 1000 00 1006 50 Dibromomethane 93 7 797 7 797 (1 114) 328434 1000.00 995 4 53 Bromodichloromethane 83 7 974 7 974 (1 140) 711520 1000 00 1103 57 C18-1,3-Dichloropropene 75 8 439 8 439 (1.206) 1073632 1000 00 1077 58 4 Methyl-2-Pentanone 43 8.635 8 635 (0 853) 1041451 1000.00 974 6 60 Toluene 91 8 714 8 714 (0 861) 3022024 1000 00 964 8 61 trans-1,3-Dichloropropene 75 9 032 9 032 (0 892) 945299 1000 00 1065 63 1,3 Dichloropropane 76 9.380 9 380 (0 926) 1034970 1000 00 982 5 64 1,12-Trichloroethane 97 9.215 9 215 (0 910) 600359 1000 00 993 6 65 Tetrachloroethane 164 9 234 9 234 (0 912) 421277 1000 00 993 6 65 Tetrachloroethane 164 9 234 9 234 (0 912) 421277 1000 00 1017 66 2-Rexanone 43 9.466 9 466 (0 935) 728700 1000 00 993 7 66 1,2-Dibromochloromethane 129 9.570 9 570 (0 945) 465815 1000 00 993 7 70 Chlorobenzene 107 9 686 9 686 (0 957) 455815 1000 00 1034 67 Dibromochloromethane 107 9 686 9 686 (0 957) 525193 1000 00 993 7 70 Chlorobenzene 112 10 157 (10 157 (1 003) 1905796 1000 00 993 7 71 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 1012) 491937 1000 00 1034 72 Ethylbenzene 106 10 236 10.236 (1 011) 1042947 1000 00 993 7 73 m + p-Xylane 106 10 359 10 159 (1.023) 2451464 2000 00 1371 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6 75 Etyrene 104 10 787 10 787 (1 065) 2116848 1000 00 993 8 76 Styrene 105 11 111 11 111 (1 10.097) 3083472 1000 00 993 7 79 Bromoform 173 10 988 10 988 11 085) 262915 1000 00 993 8 71 1,2,3-Tirchloropropane 126 11 526 11 526 (0 925) 855544 1000 00 993 8 71 1,2,3-Tirchloropropane 126 11 526 11 526 (0 925) 855544 1000 00 993 8 71 1,2,3-Tirchloropropane 126 11 636 11.636 (0 934) 777962 1000 00 993 8 71 1,2,3-Tirchloropropane 120 11 526 11 526 (0 925) 793112 1000 00 993 8 71 1,2,3-Tirchloropropane 120 11 526 11 526 (0 925) 793112 1000 00 993 8 71 1,2,3-Tirchloropropane 120 11 526 11 526 (0 925) 793112		117	6 342	6 342 (0 906)	486506	1000 00	1092
47 Trichloroethene 130 7 381 (1.055) 607870 1000 00 1026 49 1,2 Dichloropropane 63 7 675 7 675 (1.097) 787044 1000 00 1006 50 Dibromomethane 93 7 797 7 797 (1 114) 328434 1000.00 995 4 53 Bromodichloromethane 83 7 974 7 974 (1 140) 711520 1000 00 1103 57 Cis-1,3-Dichloropropene 75 8 439 8 439 (1.206) 1073632 1000 00 1077 58 4 Methyl-2-Pentanone 43 8.635 8 635 (0 853) 1041451 1000.00 974 6 60 Toluene 91 8 714 8 714 (0 861) 3022024 1000 00 964 8 61 trans-1,3-Dichloropropene 75 9 032 9 032 (0 892) 945299 1000 00 1065 63 1,3 Dichloropropane 76 9.380 9 380 (0 996) 1034970 1000 00 982 5 64 1,1,2-Trichloroethane 97 9.215 9 215 (0 910) 600359 1000 00 1065 65 Tetrachloroethane 164 9 234 9 234 (0 912) 421277 1000 00 1017 66 2-Bexanone 43 9.466 9 466 (0 935) 728700 1000 00 993 6 65 Tetrachloroethane 129 9.570 9 570 (0 945) 465815 1000 00 1152 66 1,2-Dibromochane 107 9 686 9 686 (0 957) 525393 1000 00 1152 66 1,2-Dibromochane 112 10 157 10 157 (1 003) 1905796 1000.00 993 7 70 Chlorobenzene 106 10 236 10.236 (1 011) 1042947 1000 00 983 5 71 1,1,1,2-Tetrachloroethane 106 10 236 10.236 (1 011) 1042947 1000 00 951.4 73 m + p-Xylene 106 10 756 (1 062) 1297590 1000 00 993 6 75 Stylenes (total) 106 75 Stylenes (total) 106 76 Styrene 104 10 787 10 787 (1 065) 2116848 1000 00 993 7 78 Stylenes (total) 106 79 Bromoform 173 10 988 10 988 11 085) 262915 1000 00 993 7 79 Bromoform 173 10 988 10 988 11 085) 262915 1000 00 993 6 79 Bromoform 173 10 988 10 988 11 085) 262915 1000 00 993 6 81 n-Propylbenzene 156 11 453 11 453 (0 919) 670508 1000 00 993 8 83 1,1,2,2 Tetrachloroethane 83 11 490 11 490 (0 922) 793112 1000 00 993 8 84 1,2,3-Trichloropropane 126 11 536 11 538 (0 926) 207925 1000 00 993 8 85 4-Chlorotoluene 126 11 558 11 752 (0 943) 728366 1000 00 993 99 99 99 99 99 99 99 99 99 99 99 99		78	6 63 <b>6</b>	6.636 (0 948)	2803750	. 1000 00	994 6
49 1.2 Dichloropropane 63 7 675 7 675 (1.097) 787044 1000 00 1006 50 Dibromomethane 93 7 797 7 797 (1 114) 328434 1000 00 1006 50 Dibromomethane 93 7 797 7 797 (1 114) 328434 1000 00 1103 57 cia-1,3-Dichloropropene 75 8 439 8 439 (1.206) 1073632 1000 00 1077 58 4 Methyl-2-Pentanone 43 8.635 8 635 (0 853) 1041451 1000.00 974 6 60 Toluene 91 8 714 8 714 (0 861) 3022024 1000 00 964 8 61 trans-1,3-Dichloropropene 75 9 032 9 032 (0 892) 945299 1000 00 1065 63 1,3 Dichloropropane 76 9.380 9 380 (0 926) 1034970 1000 00 982 5 64 1,1,2-Trichloroethane 97 9.215 9 215 (0 910) 600359 1000 00 932 6 65 Tetrachloroethene 164 9 234 9 234 (0 912) 421277 1000 00 943 1 66 7 Dibromochloromethane 129 9.570 9 570 (0 945) 465815 1000 00 995 7 7 0 Chloromethane 129 9.570 9 570 (0 945) 465815 1000 00 995 7 7 0 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 933 5 6 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 931 5 7 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 933 5 7 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 933 5 7 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 0 102) 491937 1000 00 1034 7 1 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 0 102) 491937 1000 00 951 6 10 7 10 7 10 7 10 7 10 7 10 7 10 7 1		62	6 770	6 770 (0 968)	788825	1000 00	1005
50 Dibromomethane 93 7 797 7 797 (1 114) 328434 1000.00 995 4 53 Bromodichloromethane 83 7 974 7 974 (1 140) 711520 1000 00 1103 57 C18-1,3-Dichloropropene 75 8 439 8 439 (1.206) 1073632 1000 00 1077 58 4 Methyl-2-Pentanone 43 8.635 8 635 (0 853) 1041451 1000.00 974 6 60 Toluene 91 8 714 8 714 (0 861) 3022024 1000 00 964 8 61 trans-1,3-Dichloropropene 75 9 032 9 032 (0 892) 945299 1000 00 1065 63 1,3 Dichloropropane 76 9.380 9 380 (0 996) 1044970 1000 00 982 5 64 1,1,2-Trichloroethane 97 9.215 9 215 (0 910) 600359 1000 00 983 6 65 Tetrachloroethene 164 9 214 9 214 0912 421277 1000 00 1017 66 2-Hexanone 43 9.466 9 466 (0 935) 728700 1000 00 943 1 67 Dibromochloromethane 129 9.570 9 570 (0 945) 465815 1000 00 985 7 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 993 5 71 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 1014 72 Ethylbenzene 106 10 236 10.236 (1 011) 1042947 1000 00 1871 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6 75 Xylenes (total) 106 10 787 (1 0 787 (1 0 65) 2116848 1000 00 1242 78 Isopropylbenzene 105 11 111 1111 (1 1,097) 308472 1000 00 1242 78 Bromoform 173 10 988 10 988 10 985 10 26915 1000 00 993 6 79 Bromochane 126 11 526 (1 526 (0 925) 856544 1000 00 993 7 83 1,1,2,2 Tetrachloroethane 131 138 11 526 (0 925) 856544 1000 00 993 8 83 1,1,2,2 Tetrachloropenae 126 11 538 (1 538 (0 934) 77962 1000 00 993 8 83 1,1,2,2 Tetrachloropenae 126 11 538 (1 526 (0 925) 856544 1000 00 993 8 84 1,2,3-Trichloropropane 126 11 538 (1 538 (0 934) 77962 1000 00 993 8 84 1,2,3-Trichloropropane 105 11 710 11 710 (0 940) 2450711 1000 00 990 9		130	7 381	7 381 (1-055)	607870	1000 00	1026
53 Bromodichloromethane  53 7 974 7 974 (1 1140) 711520 1000 00 1103 57 c18-1,3-Dichloropropene  75 8 439 8 439 (1.206) 1073632 1000 00 1077 58 4 Methyl-2-Pentanone  43 8.635 8 635 (0 853) 1041451 1000.00 974 6 60 Toluene  91 8 714 8 714 (0 861) 3022024 1000 00 964 8 61 trans-1,3-Dichloropropene  75 9 032 9 032 (0 892) 945299 1000 00 1065 63 1,3 Dichloropropane  76 9.380 9 380 (0 996) 1034970 1000 00 993 6 64 1,1,2-Trichloroethane  97 9.215 9 215 (0 910) 600359 1000 00 993 6 65 Tetrachloroethene  164 9 234 9 234 (0 912) 421277 1000 00 993 6 65 Tetrachloroethene  164 9 234 9 9.570 (0 945) 465815 1000 00 993 7 66 7 Dibromochloromethane  107 9 686 9 686 (0 957) 525393 1000 00 1152 68 1,2-Dihromochane  107 9 686 9 686 (0 957) 525393 1000 00 993 7 70 Chlorobenzene  112 10 157 10 157 (1 003) 1905796 1000.00 993 7 71 1,1,1,2-Tetrachloroethane  131 10.249 10 249 (1 012) 491937 1000 00 993 7 72 Ethylbenzene  106 10 359 10 359 (1.023) 2451464 2000 00 1871 74 Xylene-0  106 10 756 10 756 (1 065) 2116848 1000 00 993 6 75 Styrene  104 10 787 10 787 (1 065) 2116848 1000 00 993 6 75 Styrene  105 11 111 11 111 11 111 (1.097) 308472 1000 00 993 8 76 Styrene  106 10 787 10 787 (1 065) 2116848 1000 00 993 8 78 Isopropylbenzene  106 10 787 10 787 (1 065) 2116848 1000 00 993 8 79 Bromoform  173 10 988 10 988 10 983 262915 1000 00 1242 78 Isopropylbenzene  105 11 111 11 111 (1.097) 308472 1000 00 993 8 8 1a.Propylbenzene  106 11 526 11 526 (0 925) 856544 1000 00 993 8 81 1,2,2,2 Tetrachloroethane  83 11 490 11 490 (0 922) 79311 1000 00 993 8 81 1,2,2,3-Trichloropropane  107 11 538 11 538 (0 926) 207925 1000 00 994 9 87 147 Bruthloropropane  107 11 538 11 538 (0 926) 207925 1000 00 994 9 87 147 Bruthloropropane  107 11 118 119 11 110 11 110 (0 940) 2450711 1000 00 993 9 88 13 1,3,5-Trimethylbenzene  108 11 111 11 11 11 11 11 11 11 10 10 0 940) 940 940 940 940 940 940 940 940 940 940	49 1,2 Dichloropropane	63	7 675	7 675 (1.097)	787044	1000 00	1006
57 C18-1,3-D1chloropropene 75 8 439 8 439 (1.206) 1073632 1000 00 1077 58 4 Methyl-2-Pentanone 43 8.635 8 635 (0 853) 1041451 1000.00 974 6 60 Toluene 91 8 714 (0 861) 3022024 1000 00 964 8 61 trans-1,3-D1chloropropene 75 9 032 9 032 (0 892) 945299 1000 00 1065 63 1,3 D1chloropropane 76 9.380 9 380 (0 926) 1034970 1000 00 982 5 64 1,1,2-Trichloroethane 97 9.215 9 215 (0 910) 600359 1000 00 993 6 65 Tetrachloroethene 164 9 234 9 234 (0 912) 421277 1000 00 1017 66 2-Hexanone 43 9.466 9 466 (0 935) 728700 1000 00 993 6 67 Dibromochloromethane 129 9.570 9 570 (0 945) 465815 1000 00 993 7 68 1,2-D1bromocthane 107 9 686 9 686 (0 957) 525393 1000 00 993 7 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 993 5 68 1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 1034 72 Ethylbenzene 106 10 236 10.236 (1 011) 1042947 1000 00 551.4 73 m + p-Xylene 106 10 359 10 359 (1.023) 2451464 2000 00 1871 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6 75 Xylenes (total) 106 76 Styrene 104 10 787 10 787 (1 065) 2116848 1000 00 993 6 77 Bromoform 173 10 988 10 988 (1 085) 262915 1000 00 993 7 78 Bromoform 173 10 988 10 988 (1 085) 262915 1000 00 1242 78 Taopropylbenzene 105 11 111 111 11 111 (1.097) 3083472 1000 00 993 7 8 Taopropylbenzene 105 11 111 111 (1 1077) 3083472 1000 00 993 8 8 1 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 993 8 8 1 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 993 8 8 1 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 993 8 8 1 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 993 8 8 1 n-Propylbenzene 126 11 636 11 636 11 636 (0 934) 777962 1000 00 993 8 8 1 n-Propylbenzene 126 11 636 11 636 (1 636) 11 630 (0 934) 777962 1000 00 993 8 8 1 n-Propylbenzene 126 11 636 11 636 (1 636) 12 636 (1 0 934) 777962 1000 00 993 8 8 1 n-Propylbenzene 126 11 526 11 526 (0 925) 856544 1000 00 993 8 8 1 n-Propylbenzene 126 11 752 11 752 (0 943) 728366 1000 00 996 9	· -	93	7 797	7 797 (1 114)	328434	1000.00	995 4
8 4 Methyl-2-Pentanone 43 8.635 8 635 (0 833) 1041451 1000.00 974 6 60 Toluene 91 8 714 (0 861) 3022024 1000 00 974 6 60 Toluene 91 8 714 (0 861) 3022024 1000 00 974 6 60 Toluene 91 8 714 (0 861) 3022024 1000 00 974 6 60 Toluene 75 9 032 9 032 (0 892) 945299 1000 00 1065 63 1,3 Dichloropropane 76 9.380 9 380 (0 926) 1034970 1000 00 982 5 64 1,1,2-Trichloroethane 97 9.215 9 215 (0 910) 600359 1000 00 993 6 65 Tetrachloroethene 164 9 234 9 234 (0 912) 421277 1000 00 1017 66 2-Hexanone 43 9.466 9 466 (0 935) 728700 1000 00 943 1 67 Dibromochloromethane 129 9.570 9 570 (0 945) 465815 1000 00 993 5 7 70 Chloromochane 107 9 686 9 686 (0 957) 528393 1000 00 993 7 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 993 7 71 1,1,1,2-Tetrachloroethane 111 10.249 10 249 (1 012) 491937 1000 00 1034 72 Ethylbenzene 106 10 236 10.236 (1 011) 1042947 1000 00 951.4 73 m + p-Xylene 106 10 359 10 359 (1.023) 2451464 2000 00 1871 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6 8 7 75 Xylenes (total) 106	53 Bromodichloromethane	83	7 974	7 974 (1 140)	711520	1000 00	1103
60 Toluene 91 8 714 8 714 (0 861) 3022024 1000 00 964 8 61 trans-1,3-Dichloropropene 75 9 032 9 032 (0 892) 945299 1000 00 1065 63 1,3 Dichloropropane 76 9.380 9 380 (0 926) 1034970 1000 00 982 5 64 1,1,2-Trichloroethane 97 9.215 9 215 (0 910) 600359 1000 00 993 6 7 Tetrachloroethane 164 9 234 9 234 (0 912) 421277 1000 00 1017 65 2-Hexanone 43 9.466 (0 935) 728700 1000 00 943 1 67 Dibromochloromethane 129 9.570 9 570 (0 945) 465815 1000 00 1152 68 1,2-Dibromochlane 107 9 686 9 686 (0 957) 525393 1000 00 993 5 6 8 1,2-Dibromochlane 112 10 157 10 157 (1 003) 1905796 1000.00 943 5 7 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 1034 22 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2	57 cis-1,3-Dichloropropene	75	8 439	8 439 (1.206)	1073632	1000 00	1077
61 trans-1,3-Dichloropropene 75 9 032 9 032 (0 892) 945299 1000 00 1065 63 1,3 Dichloropropane 76 9.380 9 380 (0 926) 1034970 1000 00 982 5 64 1,1,2-Trichloroethane 97 9.215 9 215 (0 910) 600359 1000 00 993 6 7 7 7 7 7 7 8 7 8 8 8 8 8 8 8 8 8 8 8		43	8.635	8 635 (0 853)	1041451	1000.00	974 6
63 1.3 Dichloropropane 76 9.380 9 380 (0 926) 1034970 1000 00 982 5 64 1,1,2-Trichloroethane 97 9.215 9 215 (0 910) 600359 1000 00 983 6 65 Tetrachloroethane 164 9 234 9 234 (0 912) 421277 1000 00 1017 66 2-Hexanone 43 9.466 9 466 (0 935) 728700 1000 00 943 1 67 Dibromochloromethane 129 9.570 9 570 (0 945) 465815 1000 00 1152 68 1,2-Dibromochloromethane 107 9 686 9 686 (0 957) 525393 1000 00 995 7 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 983 5 71 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 951.4 11,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 951.4 11,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 951.4 11,1,1,2-Tetrachloroethane 134 10.249 10 236 10.236 (1 011) 1042947 1000 00 951.4 11,1,1,2-Tetrachloroethane 106 10 359 10 359 (1.023) 2451464 2000 00 1871 11,1,1,2-Tetrachloroethane 106 10 359 10 359 (1.023) 2451464 2000 00 1871 11,1,1,2-Tetrachloroethane 106 10 756 10 756 (1 062) 1297590 1000 00 993 6 1	60 Toluene	91	8 714	8 714 (0 861)	3022024	1000 00	964 8
64 1,1,2-Trichloroethane 97 9.215 9 215 (0 910) 600359 1000 00 993 6 65 Tetrachloroethane 164 9 234 9 234 (0 912) 421277 1000 00 1017 66 2-Hexanone 43 9.466 9 466 (0 935) 728700 1000 00 943 1 67 Dibromochloromethane 129 9.570 9 570 (0 945) 465815 1000 00 943 1 67 Dibromochloromethane 107 9 686 9 686 (0 957) 525393 1000 00 995 7 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 983 5 71 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 951 4 72 Ethylbenzene 106 10 236 10.236 (1 011) 1042947 1000 00 951 4 73 m + p-Xylene 106 10 359 10 359 (1.023) 2451464 2000 00 1871 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6 7 75 Xylenes (total) 106 37 10 787 (1 065) 2116848 1000 00 963 8 75 Xylenes (total) 106 37 10 988 10 988 (1 085) 262915 1000 00 975 2 8 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	61 trans-1,3-Dichloropropene	75	9 032	9 032 (0 892)	945299	1000 00	1065
65 Tetrachloroethene 164 9 234 9 234 (0 912) 421277 1000 00 1017 66 2-Hexanone 43 9.466 9 466 (0 935) 728700 1000 00 943 1 67 Dibromochloromethane 129 9.570 9 570 (0 945) 465815 1000 00 1152 68 1,2-Dibromochane 107 9 686 9 686 (0 957) 525393 1000 00 995 7 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 983 5 71 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 1034 72 Ethylbenzene 106 10 236 10.236 (1 011) 1042947 1000 00 951.4 73 m + p-Xylene 106 10 359 10 359 (1.023) 2451464 2000 00 1871 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6 75 Xylenes (total) 106	63 1,3 Dichloropropane	76	9.380	9 380 (0 926)	1034970	1000 00	982 5
66 2-Hexanone	64 1,1,2-Trichloroethane	97	9.215	9 215 (0 910)	600359	1000 00	993 6
67 Dibromochloromethane 129 9.570 9 570 (0 945) 465815 1000 00 1152 68 1,2-Dibromocthane 107 9 686 9 686 (0 957) 525393 1000 00 995 7 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 983 5 71 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 1034 72 Ethylbenzene 106 10 236 10.236 (1 011) 1042947 1000 00 951.4 73 m + p-Xylene 106 10 359 10 359 (1.023) 2451464 2000 00 1871 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6 75 Xylenes (total) 106	65 Tetrachloroethene	164	9 234	9 234 (0 912)	421277	1000 00	1017
68 1,2-Dibromocthane 107 9 686 9 686 (0 957) 525393 1000 00 995 7 70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 983 5 71 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 1034 72 Ethylbenzene 106 10 236 10.236 (1 011) 1042947 1000 00 951.4 73 m + p-Xylene 106 10 359 10 359 (1.023) 2451464 2000 00 1871 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6 75 Xylenes (total) 106 377 10 787 (1 065) 2116848 1000 00 963 8 75 Xylenes (total) 106 377 10 787 (1 065) 2116848 1000 00 963 8 77 Bromoform 173 10 988 10 988 (1 085) 262915 1000 00 1242 78 Isopropylbenzene 105 11 111 11 111 (1.097) 3083472 1000 00 975 2 79 Bromobenzene 156 11 453 11 453 (0 919) 670508 1000 00 993 8 10 1 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 993 8 8 1,1,2,2 Tetrachloroethane 83 11 490 11 490 (0 922) 793112 1000 00 964 4 12,3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4 85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 990 9		43	9.466	9 466 (0 935)	728700	1000 00	943 1
70 Chlorobenzene 112 10 157 10 157 (1 003) 1905796 1000.00 983 5 7 1 1,1,1,2-Tetrachloroethane 131 10.249 10 249 (1 012) 491937 1000 00 1034 72 Ethylbenzene 106 10 236 10.236 (1 011) 1042947 1000 00 951.4 73 m + p-Xylene 106 10 359 10 359 (1.023) 2451464 2000 00 1871 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6 75 Xylenes (total) 106 3749054 1000 00 2871 75 Xylenes (total) 106 3749054 1000 00 2871 76 Styrene 104 10 787 10 787 (1 065) 2116848 1000 00 963 8 10 988 (1 085) 262915 1000 00 1242 78 Isopropylbenzene 105 11 111 11 111 (1.097) 3083472 1000 00 975 2 79 Bromobenzene 156 11 453 11 453 (0 919) 670508 1000 00 1060 81 n-propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 992 7 82 2-Chlorotoluene 126 11 636 11.636 (0 934) 727962 1000 00 993 8 11.2,2,2 Tetrachloroethane 83 11 490 11 490 (0 922) 793112 1000 00 964 4 12.3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4 85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 990 9 87 5 2 1 5 1 5 1 5 1 5 1 5 1 5 1 5 1 5 1 5	67 Dibromochloromethane	129	9.570	9 570 (0 945)	465815	1000 00	1152
71 1,1,1,2-Tetrachloroethane	68 1,2-Dibromoethane	107	9 686	9 686 (0 957)	525393	1000 00	995 7
72 Ethylbenzene 106 10 236 10.236 (1 011) 1042947 1000 00 951.4 73 m + p-Xylene 106 10 359 10 359 (1.023) 2451464 2000 00 1871 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6  M 75 Xylenes (total) 106 3749054 1000 00 2871 76 Styrene 104 10 787 10 787 (1 065) 2116848 1000 00 963 8 77 Bromoform 173 10 988 10 988 (1 085) 262915 1000 00 1242 78 Isopropylbenzene 105 11 111 111 (1.097) 3083472 1000 00 975 2 79 Bromobenzene 156 11 453 11 453 (0 919) 670508 1000 00 1060 81 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 992 7 82 2-Chlorotoluene 126 11 636 11.636 (0 934) 727962 1000 00 993 8 83 1,1,2,2 Tetrachloropethane 83 11 490 11 490 (0 922) 793112 1000 00 964 4 85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6 86 1,3,5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	70 Chlorobenzene	112	10 157	10 157 (1 003)	1905796	1000.00	983 5
73 m + p-xylene 106 10 359 10 359 (1.023) 2451464 2000 00 1871 74 Xylene-0 106 10 756 10 756 (1 062) 1297590 1000 00 993 6  M 75 Xylenes (total) 106 3749054 1000 00 2871 76 Styrene 104 10 787 10 787 (1 065) 2116848 1000 00 963 8  77 Bromoform 173 10 988 10 988 (1 085) 262915 1000 00 1242 78 Isopropylbenzene 105 11 111 11 111 (1.097) 3083472 1000 00 975 2  79 Bromobenzene 156 11 453 11 453 (0 919) 670508 1000 00 1060  81 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 992 7  82 2-Chlorotoluene 126 11 636 11.636 (0 934) 727962 1000 00 993 8  83 1,1,2,2 Tetrachloropethane 83 11 490 11 490 (0 922) 793112 1000 00 1002  84 1.2,3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4  85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6  86 1,3,5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	71 1,1,1,2-Tetrachloroethane	131	10.249	10 249 (1 012)	491937	1000 00	1034
74 Xylene-o 106 10 756 10 756 (1 062) 1297590 1000 00 993 6  M 75 Xylenes (total) 106 3749054 1000 00 2871  76 Styrene 104 10 787 10 787 (1 065) 2116848 1000 00 963 8  77 Bromoform 173 10 988 10 988 (1 085) 262915 1000 00 1242  78 Isopropylbenzene 105 11 111 11 111 (1.097) 3083472 1000 00 975 2  79 Bromobenzene 156 11 453 11 453 (0 919) 670508 1000 00 1060  81 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 992 7  82 2-Chlorotoluene 126 11 636 11.636 (0 934) 727962 1000 00 993 8  83 1,1,2,2 Tetrachloroethane 83 11 490 11 490 (0 922) 793112 1000 00 964 4  84 1.2,3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4  85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6  86 1,3,5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	72 Ethylbenzene	106	10 236	10.236 (1 011)	1042947	1000 00	951.4
M 75 Xylenes (total) 106 3749054 1000 00 2871 76 Styrene 104 10 787 10 787 (1 065) 2116848 1000 00 963 8 77 Bromoform 173 10 988 10 988 (1 085) 262915 1000 00 1242 78 Isopropylbenzene 105 11 111 11 111 (1.097) 3083472 1000 00 975 2 79 Bromobenzene 156 11 453 11 453 (0 919) 670508 1000 00 1060 81 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 992 7 82 2-Chlorotoluene 126 11 636 11.636 (0 934) 727962 1000 00 993 8 83 1,1,2,2 Tetrachloroethane 83 11 490 11 490 (0 922) 793112 1000 00 1002 84 1.2,3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4 85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6 86 1,3,5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	73 m + p-Xylene	106	10 359	10 359 (1.023)	2451464	2000 00	1871
M 75 Xylenes (total) 106 3749054 1000 00 2871 76 Styrene 104 10 787 10 787 (1 065) 2116848 1000 00 963 8 77 Bromoform 173 10 988 10 988 (1 085) 262915 1000 00 1242 78 Isopropylbenzene 105 11 111 11 111 (1.097) 3083472 1000 00 975 2 79 Bromobenzene 156 11 453 11 453 (0 919) 670508 1000 00 1060 81 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 992 7 82 2-Chlorotoluene 126 11 636 11.636 (0 934) 727962 1000 00 993 8 83 1,2,2,2 Tetrachloroethane 83 11 490 11 490 (0 922) 793112 1000 00 1002 84 1.2,3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4 85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6 86 1,3,5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	74 Xylene-o	106	10 756	10 756 (1 062)	1297590	1000 00	
76 Styrene       104       10 787 10 787 (1 065)       2116848       1000 00       963 8         77 Bromoform       173       10 988 10 988 (1 085)       262915       1000 00       1242         78 Isopropylbenzene       105       11 111 11 11 (1.097)       3083472       1000 00       975 2         79 Bromobenzene       156       11 453 11 453 (0 919)       670508 1000 00       1060         81 n-Propylbenzene       120       11 526 11 526 (0 925)       856544 1000 00       992 7         82 2-Chlorotoluene       126       11 636 11.636 (0 934)       727962 1000 00       993 8         83 1,1,2,2 Tetrachloroethane       83       11 490 11 490 (0 922)       793112 1000 00       1002         84 1,2,3-Trichloropropane       110       11 538 11 538 (0 926)       207925 1000 00       964 4         85 4-Chlorotoluene       126       11 752 11 752 (0 943)       728366 1000 00       961 6         86 1,3,5-Trimethylbenzene       105       11 710 11 710 (0 940)       2450711 1000 00       990 9	M 75 Xylenes (total)	106			3749054		
77 Bromoform 173 10 988 10 988 (1 085) 262915 1000 00 1242 78 Isopropylbenzene 105 11 111 11 111 (1.097) 3083472 1000 00 975 2 79 Bromobenzene 156 11 453 11 453 (0 919) 670508 1000 00 1060 81 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 992 7 82 2-Chlorotoluene 126 11 636 11.636 (0 934) 727962 1000 00 993 8 83 1,1,2,2 Tetrachloroethane 83 11 490 11 490 (0 922) 793112 1000 00 1002 84 1,2,3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4 85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6 86 1,3,5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	76 Styrene	104	10 787	10 787 (1 065)			
78 Isopropylbenzene       105       11 111 11 111 (1.097) 3083472 1000 00 975 2         79 Bromobenzene       156       11 453 11 453 (0 919) 670508 1000 00 1060         81 n-Propylbenzene       120       11 526 11 526 (0 925) 856544 1000 00 992 7         82 2-Chlorotoluene       126       11 636 11.636 (0 934) 727962 1000 00 993 8         83 1,1,2,2 Tetrachloroethane       83       11 490 11 490 (0 922) 793112 1000 00 1002         84 1,2,3-Trichloropropane       110       11 538 11 538 (0 926) 207925 1000 00 964 4         85 4-Chlorotoluene       126       11 752 11 752 (0 943) 728366 1000 00 961 6         86 1,3,5-Trimethylbenzene       105       11 710 11 710 (0 940) 2450711 1000 00 990 9	77 Bromoform	173	10 988	10 988 (1 085)			
79 Bromobenzene 156 11 453 11 453 (0 919) 670508 1000 00 1060 81 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 992 7 82 2-Chlorotoluene 126 11 636 11.636 (0 934) 727962 1000 00 993 8 83 1,1,2,2 Tetrachloroethane 83 11 490 11 490 (0 922) 793112 1000 00 1002 84 1,2,3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4 85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6 86 1,3,5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	78 Isopropylbenzene	105	11 111				
81 n-Propylbenzene 120 11 526 11 526 (0 925) 856544 1000 00 992 7 82 2-Chlorotoluene 126 11 636 11.636 (0 934) 727962 1000 00 993 8 83 1,1,2,2 Tetrachloroethane 83 11 490 11 490 (0 922) 793112 1000 00 1002 84 1,2,3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4 85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6 86 1,3,5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	79 Bromobenzene	156	11 453	11 453 (0 919)			
82 2-Chlorotoluene 126 11 636 11.636 (0 934) 727962 1000 00 993 8 83 1,1,2,2 Tetrachloroethane 83 11 490 11 490 (0 922) 793112 1000 00 1002 84 1.2,3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4 85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6 86 1,3,5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	81 n-Propylbenzene	120	11 526				
83 1,1,2,2 Tetrachloroethane 83 11 490 11 490 (0 922) 793112 1000 00 1002 84 1.2,3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4 85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6 86 1,3,5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	82 2-Chlorotoluene	126	11 636				
84 1.2,3-Trichloropropane 110 11 538 11 538 (0 926) 207925 1000 00 964 4 85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6 86 1.3,5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	83 1,1,2,2 Tetrachloroethane	83					
85 4-Chlorotoluene 126 11 752 11 752 (0 943) 728366 1000 00 961 6 86 1.3,5-Trimethylbenzene 105 11 710 (1 940) 2450711 1000 00 990 9	84 1,2,3-Trichloropropane	110					
86 1.3.5-Trimethylbenzene 105 11 710 11 710 (0 940) 2450711 1000 00 990 9	85 4-Chlorotoluene	126	11 752				
87 tart Dutylhongona	86 1,3,5-Trimethylbenzene	105	11 710				
	87 tert-Butylbenzene	119	12 015	12 015 (0 964)			

Page 3

# 657 174

						AMOUN	TS
		QUANT SIG				CAL-AMT	ON-COL
Compo	unds	MASS	RT	EXP RT REL RT	RESPONS2	( ng)	(ng)
	=======================================	====	2 2	===== ======	=======	======	c=====#
88	1,2,4 Trimetnylbenzene	105	12 083	12 083 (0 970)	2491491	1000 00	986 9
89	sec-Butylbenzene	105	12 242	12 242 (0.982)	3366610	1000 00	977 0
90	4-Isopropyltoluene	119	12.388	12 388 (0 994)	2322311	1000 00	903 9
91	1,3-Dichlorobenzene	146	12 382	12 382 (0 994)	1170519	1000 00	928 0
93	1,4-Dichlorobenzene	146	12.486	12 486 (1 002)	1259741	1000 00	979 6
94	n-Butylbenzene	91	12.798	12 798 (1 027)	2952832	1000 00	1003
95	1,2-Dichlorobenzene	146	12 859	12 859 (1 032)	1193990	1000 00	1006
96	1,2-Dibromo-3-chloropropane	157	13 660	13 660 (1 096)	105902	1000 00	1338
97	1,2,4-Tricalorobenzene	180	14 461	14 461 (1 160)	588023	1000 00	1073
98	Hexachlorobutadiene	225	14.583	14 583 (1 170)	293669	1000 00	1108
99	Naphthalene	128	14 711	14 711 (1 181)	1312547	1000 00	1070
100	1,2,3-Trichlorobenzene	180	14 962	14 962 (1 201)	410033	1000 00	944 6

# VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: 40826D

Instrument ID: HP4 Calibration Date: 05/17/00 Time: 1108

Heated Purge: (Y/N) N Init. Calib. Times: 1108 1303

GC Column: DB 624 ID: 0.18 (mm)

			MĪN	1 1	MAX
COMPOUND	RRF	RRF50	RRF	%D	%D
	========	======	=======	=====	====
Dichlorodifluoromethane	0.122	0.127	0.01	4.1	50.0
Chloromethane	0.275	0.287	0.1	4.4	50.0
Vinyl Chloride	0.260	0.262	0.01		20.0
Bromomethane	0.038	0.040	0.01		50.0
Chloroethane	0.044	0.049	0.01		50.0
Trichlorofluoromethane	0.070	0.083	0.01	18.6	50.0
1,1-Dichloroethene	0.162	0.180	0.01		
Methylene Chloride	0.200				
trans-1,2-Dichloroethene	0.194				50.0
1,1-Dichloroethane	0.397	0.424	0.1		50.0
cis-1,2-dichloroethene	0.224	0.241			50.0
Chloroform	0.311	0.332	0.01		
Bromochloromethane	0.088				
1,1,1-Trichloroethane	0.230				50.0
Carbon Tetrachloride	0.145				50.0
1,2-Dichloroethane	0.254				50.0
Benzene	0.913				50.0
Trichloroethene	0.192				50.0
1,2-Dichloropropane	0.253				20.0
Bromodichloromethane	0.209				50.0
cis-1,3-Dichloropropene	0.323				50.0
Toluene	4.483				20.0
trans-1,3-Dichloropropene	1.270				50.0
1,1,2-Trichloroethane	0.865				50.0
Tetrachloroethene	0.593				50.0
Dibromochloromethane	_      0.578				50.0
Chlorobenzene	2.773				50.0
Ethylbenzene	1.569				20.0
Styrene	3.144				50.0
Bromoform	0.303				50.0
1,1,2,2-Tetrachloroethane_	0.805				50.0
1,3-Dichlorobenzene	1.284			8.5	50.0
1,4-Dichlorobenzene	1.309				50.0
1,2-Dichlorobenzene	1.208				50.0
Dibromomethane	0.107				50.0
1,2-Dibromoethane	0.755				50.0
1,1,1,2-Tetrachloroethane	0.681	0.753	0.01	10.6	50.0

page 1 of 2

FORM VII VOA

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: 40826D

Instrument ID: HP4 Calibration Date: 05/17/00 Time: 1108

Heated Purge: (Y/N) N Init. Calib. Times: 1108 1303

GC Column: DB 624 ID: 0.18 (mm)

1,2,3-Trichloropropane		222	DDDEA	MIN	%D	MAX %D
1,2,3-Trichloropropane	COMPOUND	RRF	RRF50	RRF		1
1,2-Dibromo-3-chloropropane						5 I
2,2-Dichloropropane						
1,1-Dichloropropene				· ·		
1.508						
n-Propylbenzene						
Bromobenzene						
1,3,5-Trimethylbenzene       2.517       2.670       0.01       6.1       50.0         2-Chlorotoluene       0.745       0.791       0.01       6.2       50.0         4-Chlorotoluene       0.771       0.827       0.01       7.3       50.0         tert-Butylbenzene       2.151       2.296       0.01       6.7       50.0         1,2,4-Trimethylbenzene       2.569       2.742       0.01       6.7       50.0         sec-Butylbenzene       3.507       3.736       0.01       6.5       50.0         4-Isopropyltoluene       2.615       2.881       0.01       10.2       50.0         n-Butylbenzene       2.995       3.137       0.01       4.7       50.0         n-Butylbenzene       0.558       0.516       0.01       7.5       50.0         1,2,4-Trichlorobenzene       0.270       0.243       0.01       10.0       50.0         Naphthalene       1.248       1.061       0.01       15.0       50.0         1,2,3-Trichlorobenzene       0.442       0.357       0.01       19.2       50.0         Acetone       0.103       0.098       0.01       4.8       50.0         Carbon Disulfide						
2-Chlorotoluene						
4-Chlorotoluene       0.771       0.827       0.01       7.3       50.0         tert-Butylbenzene       2.151       2.296       0.01       6.7       50.0         1,2,4-Trimethylbenzene       2.569       2.742       0.01       6.7       50.0         sec-Butylbenzene       3.507       3.736       0.01       6.5       50.0         4-Isopropyltoluene       2.615       2.881       0.01       10.2       50.0         n-Butylbenzene       2.995       3.137       0.01       4.7       50.0         1,2,4-Trichlorobenzene       0.558       0.516       0.01       7.5       50.0         Hexachlorobutadiene       0.270       0.243       0.01       10.0       50.0         Naphthalene       1.248       1.061       0.01       15.0       50.0         1,2,3-Trichlorobenzene       0.442       0.357       0.01       19.2       50.0         Acetone       0.103       0.098       0.01       4.8       50.0         Carbon Disulfide       0.494       0.548       0.01       10.9       50.0         2-Butanone       1.562       0.01       1.5       50.0         4-Methyl-2-Pentanone       1.529       <						
tert-Butylbenzene         2.151         2.296         0.01         6.7         50.0           1,2,4-Trimethylbenzene         2.569         2.742         0.01         6.7         50.0           sec-Butylbenzene         3.507         3.736         0.01         6.5         50.0           4-Isopropyltoluene         2.615         2.881         0.01         10.2         50.0           n-Butylbenzene         2.995         3.137         0.01         4.7         50.0           1,2,4-Trichlorobenzene         0.558         0.516         0.01         7.5         50.0           Hexachlorobutadiene         0.270         0.243         0.01         10.0         50.0           Naphthalene         1.248         1.061         0.01         15.0         50.0           1,2,3-Trichlorobenzene         0.442         0.357         0.01         19.2         50.0           Acetone         0.103         0.098         0.01         4.8         50.0           Carbon Disulfide         0.494         0.548         0.01         10.9         50.0           2-Butanone         1.529         1.562         0.01         1.5         50.0           4-Methyl -2-Pentanone         1.06						
1,2,4-Trimethylbenzene       2.569       2.742       0.01       6.7       50.0         sec-Butylbenzene       3.507       3.736       0.01       6.5       50.0         4-Isopropyltoluene       2.615       2.881       0.01       10.2       50.0         n-Butylbenzene       2.995       3.137       0.01       4.7       50.0         1,2,4-Trichlorobenzene       0.558       0.516       0.01       7.5       50.0         Hexachlorobutadiene       0.270       0.243       0.01       10.0       50.0         Naphthalene       1.248       1.061       0.01       15.0       50.0         1,2,3-Trichlorobenzene       0.442       0.357       0.01       19.2       50.0         Acetone       0.103       0.098       0.01       4.8       50.0         Carbon Disulfide       0.494       0.548       0.01       10.9       50.0         2-Butanone       1.529       1.562       0.01       1.1       50.0         4-Methyl-2-Pentanone       1.06       1.089       0.01       1.5       50.0         2-Hexanone       0.475       0.506       0.01       6.5       50.0         Isopropylbenzene       4.52						
sec-Butylbenzene       3.507       3.736       0.01       6.5       50.0         4-Isopropyltoluene       2.615       2.881       0.01       10.2       50.0         n-Butylbenzene       2.995       3.137       0.01       4.7       50.0         1,2,4-Trichlorobenzene       0.558       0.516       0.01       7.5       50.0         Hexachlorobutadiene       0.270       0.243       0.01       10.0       50.0         Naphthalene       1.248       1.061       0.01       15.0       50.0         1,2,3-Trichlorobenzene       0.442       0.357       0.01       19.2       50.0         Acetone       0.103       0.098       0.01       4.8       50.0         Carbon Disulfide       0.494       0.548       0.01       10.9       50.0         2-Butanone       0.178       0.176       0.01       1.1       50.0         4-Methyl-2-Pentanone       1.529       1.562       0.01       1.5       50.0         2-Hexanone       4.525       4.947       0.01       1.5       50.0         Methyl tert-butyl ether       0.256       0.506       0.01       6.5       50.0         1,2-Dichloroethene (total)						
4-Isopropyltoluene       2.615       2.881       0.01       10.2       50.0         n-Butylbenzene       2.995       3.137       0.01       4.7       50.0         1,2,4-Trichlorobenzene       0.558       0.516       0.01       7.5       50.0         Hexachlorobutadiene       0.270       0.243       0.01       10.0       50.0         Naphthalene       1.248       1.061       0.01       15.0       50.0         1,2,3-Trichlorobenzene       0.442       0.357       0.01       19.2       50.0         Acetone       0.103       0.098       0.01       4.8       50.0         Carbon Disulfide       0.494       0.548       0.01       10.9       50.0         2-Butanone       0.178       0.176       0.01       1.1       50.0         4-Methyl-2-Pentanone       1.529       1.562       0.01       1.5       50.0         2-Hexanone       0.475       0.506       0.01       6.5       50.0         Methyl tert-butyl ether       0.475       0.506       0.01       6.5       50.0         1,2-Dichloroethene (total)       1.869       2.004       0.01       7.2       50.0         2,20       1.50						
n-Butylbenzene       2.995       3.137       0.01       4.7       50.00         1,2,4-Trichlorobenzene       0.558       0.516       0.01       7.5       50.00         Hexachlorobutadiene       0.270       0.243       0.01       10.0       50.00         Naphthalene       1.248       1.061       0.01       15.0       50.00         1,2,3-Trichlorobenzene       0.442       0.357       0.01       19.2       50.00         Acetone       0.103       0.098       0.01       4.8       50.00         Carbon Disulfide       0.494       0.548       0.01       10.9       50.00         2-Butanone       0.178       0.176       0.01       1.1       50.00         4-Methyl-2-Pentanone       1.529       1.562       0.01       2.2       50.00         2-Hexanone       1.106       1.089       0.01       1.5       50.00         Methyl tert-butyl ether       0.475       0.506       0.01       6.5       50.00         I,2-Dichloroethene (total)       0.209       0.228       0.01       9.1       50.00         Xylenes (total)       1.869       2.004       0.01       7.2       50.00         Dibromofluorome						
1,2,4-Trichlorobenzene       0.558       0.516       0.01       7.5       50.0         Hexachlorobutadiene       0.270       0.243       0.01       10.0       50.0         Naphthalene       1.248       1.061       0.01       15.0       50.0         1,2,3-Trichlorobenzene       0.442       0.357       0.01       19.2       50.0         Acetone       0.103       0.098       0.01       4.8       50.0         Carbon Disulfide       0.494       0.548       0.01       10.9       50.0         2-Butanone       0.178       0.176       0.01       1.1       50.0         4-Methyl-2-Pentanone       1.529       1.562       0.01       2.2       50.0         2-Hexanone       1.106       1.089       0.01       1.5       50.0         Methyl tert-butyl ether       0.475       0.506       0.01       6.5       50.0         Isopropylbenzene       4.525       4.947       0.01       9.3       50.0         1,2-Dichloroethene (total)       1.869       2.004       0.01       7.2       50.0         2-2-2-2-2-2-2-2-2-2-2-2-2-2-2-2-2-2-2-						
Hexachlorobutadiene       0.270       0.243       0.01       10.0       50.0         Naphthalene       1.248       1.061       0.01       15.0       50.0         1,2,3-Trichlorobenzene       0.442       0.357       0.01       19.2       50.0         Acetone       0.103       0.098       0.01       4.8       50.0         Carbon Disulfide       0.494       0.548       0.01       10.9       50.0         2-Butanone       0.178       0.176       0.01       1.1       50.0         4-Methyl-2-Pentanone       1.529       1.562       0.01       2.2       50.0         2-Hexanone       1.106       1.089       0.01       1.5       50.0         Methyl tert-butyl ether       0.475       0.506       0.01       6.5       50.0         Isopropylbenzene       4.525       4.947       0.01       9.3       50.0         Xylenes (total)       1.869       2.004       0.01       7.2       50.0         Dibromofluoromethane       0.171       0.170       0.01       0.6       50.0         1,2-Dichloroethane-d4       0.234       0.235       0.01       0.4       50.0         Toluene-d8       4.199 </td <td>1 3 4 Trichlorchongone</td> <td></td> <td></td> <td></td> <td></td> <td></td>	1 3 4 Trichlorchongone					
Naphthalene       1.248       1.061       0.01       15.0       50.0         1,2,3-Trichlorobenzene       0.442       0.357       0.01       19.2       50.0         Acetone       0.103       0.098       0.01       4.8       50.0         Carbon Disulfide       0.494       0.548       0.01       10.9       50.0         2-Butanone       0.178       0.176       0.01       1.1       50.0         4-Methyl-2-Pentanone       1.529       1.562       0.01       2.2       50.0         2-Hexanone       1.106       1.089       0.01       1.5       50.0         Methyl tert-butyl ether       0.475       0.506       0.01       6.5       50.0         Isopropylbenzene       4.525       4.947       0.01       9.3       50.0         1,2-Dichloroethene (total)       1.869       2.004       0.01       7.2       50.0         Xylenes (total)       1.869       2.004       0.01       7.2       50.0         Dibromofluoromethane       0.234       0.235       0.01       0.4       50.0         Toluene-d8       4.199       4.201       0.01       0.0       50.0	Userable rebutadions					
1,2,3-Trichlorobenzene       0.442       0.357       0.01       19.2       50.0         Acetone       0.103       0.098       0.01       4.8       50.0         Carbon Disulfide       0.494       0.548       0.01       10.9       50.0         2-Butanone       0.178       0.176       0.01       1.1       50.0         4-Methyl-2-Pentanone       1.529       1.562       0.01       2.2       50.0         2-Hexanone       1.106       1.089       0.01       1.5       50.0         Methyl tert-butyl ether       0.475       0.506       0.01       6.5       50.0         Isopropylbenzene       4.525       4.947       0.01       9.3       50.0         1,2-Dichloroethene (total)       0.209       0.228       0.01       9.1       50.0         Xylenes (total)       1.869       2.004       0.01       7.2       50.0         1,2-Dichloroethane       0.171       0.170       0.01       0.6       50.0         1,2-Dichloroethane-d4       0.234       0.235       0.01       0.4       50.0         Toluene-d8       4.199       4.201       0.01       0.0       50.0						
Acetone       0.103       0.098       0.01       4.8 50.0         Carbon Disulfide       0.494       0.548       0.01       10.9 50.0         2-Butanone       0.178       0.176       0.01       1.1 50.0         4-Methyl-2-Pentanone       1.529       1.562       0.01       2.2 50.0         2-Hexanone       1.106       1.089       0.01       1.5 50.0         Methyl tert-butyl ether       0.475       0.506       0.01       6.5 50.0         Isopropylbenzene       4.525       4.947       0.01       9.3 50.0         1,2-Dichloroethene (total)       0.209       0.228       0.01       9.1 50.0         Xylenes (total)       1.869       2.004       0.01       7.2 50.0         1,2-Dichloroethane       0.171       0.170       0.01       0.6 50.0         1,2-Dichloroethane-d4       0.234       0.235       0.01       0.4 50.0         Toluene-d8       4.199       4.201       0.01       0.0 50.0						
Carbon Disulfide       0.494       0.548       0.01       10.9       50.0         2-Butanone       0.178       0.176       0.01       1.1       50.0         4-Methyl-2-Pentanone       1.529       1.562       0.01       2.2       50.0         2-Hexanone       1.106       1.089       0.01       1.5       50.0         Methyl tert-butyl ether       0.475       0.506       0.01       6.5       50.0         Isopropylbenzene       4.525       4.947       0.01       9.3       50.0         1,2-Dichloroethene (total)       0.209       0.228       0.01       9.1       50.0         Xylenes (total)       1.869       2.004       0.01       7.2       50.0         1,2-Dichloroethane       0.171       0.170       0.01       0.6       50.0         1,2-Dichloroethane-d4       0.234       0.235       0.01       0.4       50.0         1,2-Dichloroethane-d8       4.199       4.201       0.01       0.0       50.0						
2-Butanone       0.178       0.176       0.01       1.1       50.0         4-Methyl-2-Pentanone       1.529       1.562       0.01       2.2       50.0         2-Hexanone       1.106       1.089       0.01       1.5       50.0         Methyl tert-butyl ether       0.475       0.506       0.01       6.5       50.0         Isopropylbenzene       4.525       4.947       0.01       9.3       50.0         1,2-Dichloroethene (total)       0.209       0.228       0.01       9.1       50.0         Xylenes (total)       1.869       2.004       0.01       7.2       50.0         1,2-Dichloroethane       0.171       0.170       0.01       0.6       50.0         1,2-Dichloroethane-d4       0.234       0.235       0.01       0.4       50.0         1,2-Dichloroethane-d8       4.199       4.201       0.01       0.0       50.0						
4-Methyl-2-Pentanone       1.529       1.562       0.01       2.2       50.0         2-Hexanone       1.106       1.089       0.01       1.5       50.0         Methyl tert-butyl ether       0.475       0.506       0.01       6.5       50.0         Isopropylbenzene       4.525       4.947       0.01       9.3       50.0         1,2-Dichloroethene (total)       0.209       0.228       0.01       9.1       50.0         Xylenes (total)       1.869       2.004       0.01       7.2       50.0         2						
2-Hexanone       1.106       1.089       0.01       1.5 50.0         Methyl tert-butyl ether       0.475       0.506       0.01       6.5 50.0         Isopropylbenzene       4.525       4.947       0.01       9.3 50.0         1,2-Dichloroethene (total)       0.209       0.228       0.01       9.1 50.0         Xylenes (total)       1.869       2.004       0.01       7.2 50.0         Dibromofluoromethane       0.171       0.170       0.01       0.6 50.0         1,2-Dichloroethane-d4       0.234       0.235       0.01       0.4 50.0         Toluene-d8       4.199       4.201       0.01       0.0 50.0						
Methyl tert-butyl ether       0.475       0.506       0.01       6.5 50.0         Isopropylbenzene       4.525       4.947       0.01       9.3 50.0         1,2-Dichloroethene (total)						
Isopropylbenzene       4.525       4.947       0.01       9.3       50.0         1,2-Dichloroethene (total)       0.209       0.228       0.01       9.1       50.0         Xylenes (total)       1.869       2.004       0.01       7.2       50.0         Dibromofluoromethane       0.171       0.170       0.01       0.6       50.0         1,2-Dichloroethane-d4       0.234       0.235       0.01       0.4       50.0         Toluene-d8       4.199       4.201       0.01       0.0       50.0						
1,2-Dichloroethene (total)       0.209       0.228       0.01       9.1       50.0         Xylenes (total)       1.869       2.004       0.01       7.2       50.0         Dibromofluoromethane       0.171       0.170       0.01       0.6       50.0         1,2-Dichloroethane-d4       0.234       0.235       0.01       0.4       50.0         Toluene-d8       4.199       4.201       0.01       0.0       50.0					9.3	50.0
Xylenes (total)       1.869       2.004       0.01       7.2       50.0         Dibromofluoromethane       0.171       0.170       0.01       0.6       50.0         1,2-Dichloroethane-d4       0.234       0.235       0.01       0.4       50.0         Toluene-d8       4.199       4.201       0.01       0.0       50.0					9.1	50.0
Dibromofluoromethane       0.171       0.170       0.01       0.6       50.0         1,2-Dichloroethane-d4       0.234       0.235       0.01       0.4       50.0         Toluene-d8       4.199       4.201       0.01       0.0       50.0					7.2	50.0
1,2-Dichloroethane-d4     0.234     0.235     0.01     0.4 50.0       Toluene-d8     4.199     4.201     0.01     0.0 50.0		========	=======	======	=====	====
1,2-Dichloroethane-d4     0.234     0.235     0.01     0.4 50.0       Toluene-d8     4.199     4.201     0.01     0.0 50.0	Dibromofluoromethane	0.171	0.170	0.01	0.6	50.0
Toluene-d8 4.199 4.201 0.01 0.0 50.0					0.4	50.0
					0.0	50.0
					1.3	50.0
· · · · · · · · · · · · · · · · · · ·					Ì	<u> </u>

page 2 of 2

FORM VII VOA

Data File: /var/chem/hp4.i/40517d.b/cc40517.d

Report Date: 05/17/2000

#### CONTINUING CALIBRATION COMPOUNDS PERCENT DRIFT REPORT

Instrument ID: hp4.i Injection Date: 17-MAY-2000 11:08

Lab File ID: cc40517.d

Lab Sample ID: vstd50 Method File: /var/chem/hp4.i/40517d.b/8260bh2o.m Analysis Type: WATER

	EXPECTED	Measured		MAX
COMPOUND	CONC.	CONC	₹D	₽D.
		=======================================		
0 Xylene-o	250.0000	268 0833		50 (
0 m + p-Xylene	500.0000		•	50.0
1 Dichlorodifluoromethane	250 0000			50.0
2 Chloromethane	250.0000	260 9784		50 (
3 Vinyl Chloride	250.0000	252 2334	0 9	20.
4 Bromomethane	250.0000	267 7116	7 1	50
5 Chloroethane	250.0000	281 4583	12 6	50
6 Trichlorofluoromethane	250.0000	297 0377		50
9 1,1-Dichloroethene	250 0000	277 3090	10 9	20.
10 Methylene Chloride	250 0000	279 1715	11 7	50
13 trans-1,2-Dichloroethene	250.0000	278 4390	11.4	50.
15 1,1-Dichloroethane	250 0000	266 9890	6.8	50.
17 cis-1,2-dichloroethene	250 0000	269 0701	7 6	50.
18 Chloroform	250.0000	267.0103	6 8	20.
19 Bromochloromethane	250 0000	266 0261	6.4	50
20 1,1,1-Trichloroethane	250.0000	267 4681	7 0	50.
21 Carbon Tetrachloride	250.0000	275 4105	10.2	50.
23 1,2-Dichloroethane	250.0000	269.7169	7 9	50
24 Benzene	250 0000	268 0992	7.2	50.
26 Trichloroethene	250 0000	268 8808	7.6	50.
27 1,2-Dichloropropane	250 0000	268.6392	7.5	20
28 Bromodichloromethane	[ 250.0000]	275 8652	10.3	50.
31 cis-1,3-Dichloropropene	250.0000	275 3741	10 1	50.
33 Toluene	250 0000	267 6848	7 1	20
34 trans-1,3-Dichloropropene	250.0000	271.6155	8 6	50.
36 1,1,2-Trichloroethane	250 0000	265 8106	6.3	50.
37 Tetrachloroethene	250 0000	266.1906	6 5	50
38 Dibromochloromethane	250.0000	273.7660	9 5	50
40 Chlorobenzene	250.0000	266 4739	6 6	50
41 Ethylbenzene	250.0000	268.8007	7 5	20
44 Styrene	250 0000	275.8051	10 3	50.
45 Bromoform	250.0000	276.9511	10.8	50.
46 1,1,2,2-Tetrachloroethane	250.0000j	268 4579	74	50
48 1,3-Dichlorobenzene	250.0000	271 2924	8 5	50
49 1,4-Dichlorobenzene	250.0000	262 9850	5.2	50.
50 1,2-Dichlorobenzene	250.0000	264 5568	5 8[	50.
60 Dibromomethane	250 0000	276.7437	10.7	50
63 1,2-Dibromoethane	250 0000	268 3789	7.4	50.
64 1,1,1,2-Tetrachloroethane	1 250 00001	276 7052	10.7	50

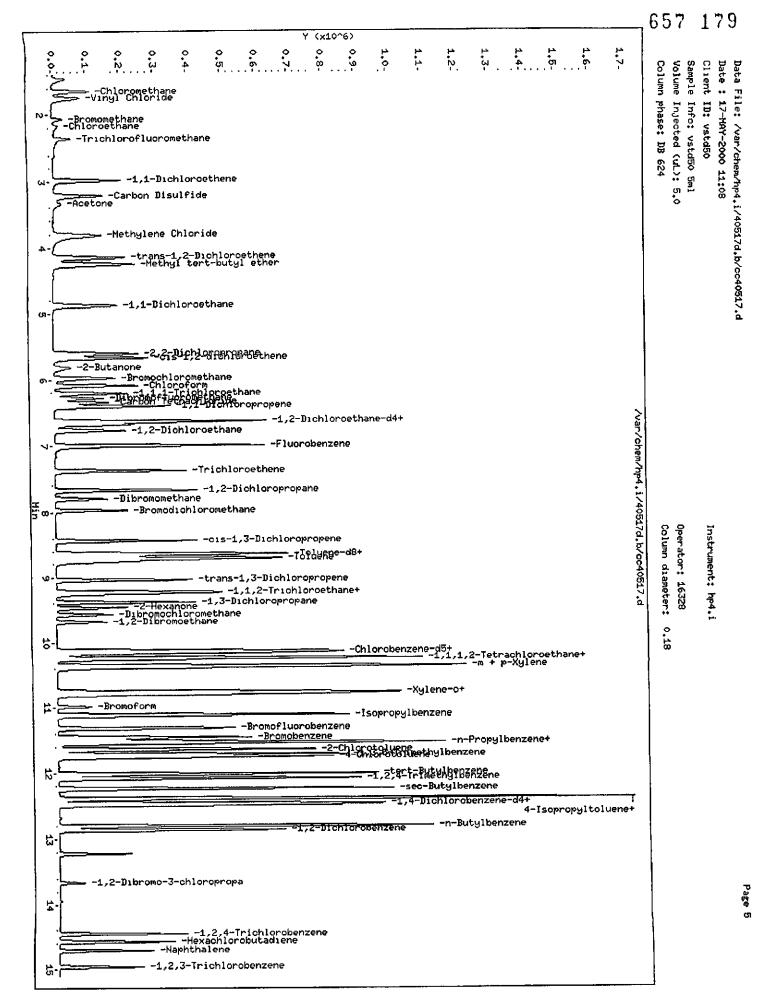
Data File: /var/chem/hp4.i/40517d.b/cc40517.d Report Date: 05/17/2000

### CONTINUING CALIBRATION COMPOUNDS PERCENT DRIFT REPORT

Instrument ID: hp4.i Lab File ID: cc40517.d Analysis Type: WATER Injection Date: 17-MAY-2000 11:08

Lab Sample ID: vstd50
Method File: /var/chem/hp4.i/40517d.b/8260bh2

1			EXPECTED	MEASURED	1	MAX
	CO	1POUND	CONC.	CONC	%D	%D
=	====	***************************************			3=3===	
1	65	1,2,3-Trichloropropane	250.0000	270.3583	8.1	50 0
1	69	1,2-Dibromo-3-chloropropane	250 0000	250 5436	0 2	50 0
1	74	2,2-Dichloropropane	250 0000	265,0125	6 0	50.0
	75	1,1-Dichloropropene	250 0000	269 0169	7 6	50.0
1	77	1,3-Dichloropropane	250 0000	266 6257	6.7	50 0
1	83	n-Propylbenzene	250 0000	264 7772	5.9	50 0
}	84	Bromobenzene	250 0000	261 7646	4.7	50 0
1	85	1,3,5-Trimethylbenzene	250.0000	265 1801	6.1	50.0
1	86	2-Chlorotoluene	250 0000	265 3799	6.2	50 0
1	87	4-Chlorotoluene	250 0000	268.2931	7.3	50.0
1	88	tert-Butylbenzene	250.0000	266 8661	6.7	50 0
1	89	1,2,4-Trimethylbenzene	250.0000	266 8274	6.7	50 0
	90	sec-Butylbenzene	250.0000	266 3363	6.5	50 0
1	91	4-Isopropyltoluene	250.0000	275 4893	10.2	50.0
1	94	n-Butylbenzene	250 0000	261 9074	4 8	50.0
1	95	1,2,4-Trichlorobenzene	250.0000	231 3182	7 5	50.0
1	96	Hexachlorobutadiene	250.0000	224 8516	10.1	50.0
Ι	97	Naphthalene	250 0000	212 5767	15 0	50 0
I	98	1,2,3-Trichlorobenzene	250 0000	201 8390	19 3	50 0
Ι	101	Chlorobenzene-d5	250 0000	250.0000	0.0	50.0
j	102	1,4-Dichlorobenzene-d4	250.0000	250 0000	0.0	50.0
Ι	106	Acetone	250.0000	237 7850	4.9	50.0
Τ	107	Carbon Disulfide	250.0000	277.4393	11 0	50.0
1	108	2-Butanone	250.0000	247 2264	1.1	50.0
	110	4-Methyl-2-Pentanone	250.0000	255 4008	2.2	50 0
1	111	2-Hexanone	250.0000	246 1886	1.5	50.0
ļ	134	Methyl tert-butyl ether	250 0000	266.3779	6.6	50 0
1	137	Fluorobenzene	250.0000	250.0000	0.0	50.0
ŀ	144	Isopropylbenzene	25 <b>0</b> 0000	273.2772	9.3	50 0
İ	153	1,2-Dichloroethene (total)	500 0000	546.8472	9.4	50 0
1	154	Xylenes (total)	750.0000	814.3590	8 5	50 0
İ	149	Dibromofluoromethane	250 0000	248 1923	0 7	50 0
Ì	150	1,2-Dichloroethane-d4	250.0000	251 3176	0.5	50 0
ţ	151	Toluene-d8	[ 250.0000]	250.0962	0.0	50.0
}	152	Bromofluorobenzene	250 0000	253 1797	1 3	50 0
1_						



Page 1

KsInla

Data File: /var/chem/hp4.i/40517d.b/cc40517.d

Report Date: 17-May-2000 14:06

657 180

#### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file : /var/chem/hp4.i/40517d.b/cc40517.d

Lab Smp Id: vstd50 Client Smp ID: vstd50

Inj Date : 17-MAY-2000 11:08

Operator: 16328 Inst ID: hp4.i

Smp Info : vstd50 5ml

Misc Info : ,40517d.b, med1000.m

Comment :

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m

Meth Date: 17-May-2000 14:06 dudeckk Quant Type: ISTD Cal Date: 17-MAY-2000 13:03 Cal File: 1e40517.d

Als bottle: 4 Calibration Sample, Level: 3

Dil Factor: 1.00000 Integrator: HP RTE

Integrator: HP RTE Compound Sublist: 2-padep.sub

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

Name	Value	Description
DF Vf Vo Vi	5.000 5.000	Dilution Factor FinalVolume Sample Volume Purge Volume

						AMOUN	TS
		QUANT SIG				CAL-AMT	ON-COL
Com	pounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
				*****	****	#==== <b>=</b>	======
*	46 Fluorobenzene	96	6 994	6.994 (1 000)	660677		
*	69 Chlorobenzene-d5	119	10 124	10.124 (1 000)	149779		
•	92 1,4-Dichlorobenzene-d4	152	12.460	12 460 (1 000)	227254		
\$	39 Dibromofluoromethane	113	6 285	6 285 (0 899)	112189	250.000	248.2
s	43 1,2-Dichloroethane-d4	65	6 683	6 683 (0 955)	155579	250.000	251 3
\$	59 Toluene-d8	98	8.645	8 645 (0.854)	629253	250 000	250.1
\$	80 Bromofluorobenzene	95	11 310	11 310 (1 117)	246086	250.000	253.2
	1 Dichlorodifluoromethane	85	1 450	1.450 (0.207)	84060	250.000	261.0
	2 Chloromethane	50	1.645	1.645 (0.235)	189504	250.000	261 0
	3 Vinyl Chloride	62	1 743	1 743 (0 249)	173382	250.000	252 2
	4 Bromomethane	94	2.061	2 061 (0 295)	26679	250.000	267.7
	5 Chloroethane	64	2.165	2.165 (0.309)	32667	250 000	281 4
	6 Trichlorofluoromethane	101	2 354	2.354 (0 337)	54753	250 000	297 0
	12 1.1-Dichloroethene	96	2 978	2 978 (0.426)	118932	250.000	277 3
	13 Acetone	43	3 351	3 351 (0 479)	64935	250 000	237 В
	<del></del>						

Data File: /var/chem/hp4.i/40517d.b/cc40517.d Report Date: 17-May-2000 14:06

					AMOUN	TS
	QUANT SIG				CAL-AMT	ON-COL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
克耳达 经自然 医医性动脉 化苯基乙烷 经基础 医中中中毒		= =	BCE=== =================================		******	=====
15 Carbon Disulfide	76	3 216	3 216 (0 460)	362443	250 000	277.4
18 Methylene Chloride	84	3.815	3 815 (0 545)	147271	250.000	279.2
19 trans-1,2-Dichloroethene	96	4.152	4 152 (0.594)	142814	250.000	278 4
20 Methyl tert-butyl ether	73	4.249	4 249 (0 608)	334257	250 000	266 4
24 1,1-Dichloroethane	63	4.867	4 867 (0 696)	280382	250 000	267.0
27 2,2-Dichloropropane	77	5.613	5 613 (0 802)	167972	250 000	265.0
28 cis-1,2-dichloroethene	96	5 674	5.674 (0 811)	158989	250 000	269.1
M 29 1,2-Dichloroethene (total)	100			301803	500.000	546.8
30 Bromochloromethane	128	5.973	5 973 (0 854)	62236	250.000	266.0
31 2-Butanone	43	5 821	5.821 (0 832)	116124	250 000	247.2
37 Chloroform	83	6 090	6.090 (0 871)	219279	250 000	267.0
38 1,1,1-Trichloroethane	97	6 212	6 212 (0.888)	162494	250.000	267.5
40 1,1-Dichloropropene	75	6.407	6.407 (0.916)	181812	250.000	269.0
41 Carbon Tetrachloride	117	6.352	6.352 (0 908)	105078	250.0 <b>00</b>	275.4
42 Benzene	78	6.640	6.640 (0 949)	646947	250.000	268 1
45 1,2-Dichloroethane	62	6.774	6 774 (0.969)	181162	250.000	269 7
47 Trichloroethene	130	7.379	7 379 (1.055)	136342	250 000	268 9
49 1,2-Dichloropropane	63	7.673	7.673 (1.097)	179867	250.000	268 6
50 Dibromomethane	93	7 795	7.795 (1.114)	78170	250 000	276 7
53 Bromodichloromethane	83	7 972	7 972 (1.140)	152326	250 000	275 9
57 cis-1,3-Dichloropropene	75	8.437	8 437 (1 206)	234977	250 000	275 4
58 4-Methyl-2-Pentanone	43	8.627	8.627 (0 852)	234014	250.000	255 4
60 Toluene	91	8 712	8.712 (0.861)	718952	250 000	267 7
61 trans-1,3-Dichloropropene	75	9.030	9.030 (0.892)	206628	250.000	271 6
63 1,3-Dichloropropane	76	9.379	9 379 (0.926)	240837	250 000	266 6
64 1,1,2-Trichloroethane	97	9 207	9 207 (0.909)	137716	250 000	265 8
65 Tetrachloroethene	164	9 232	9 232 (0.912)	94563	250.000	266.2
66 2-Hexanone	43	9.458	9.458 (0.934)	163100	250.000	246 2
67 Dibromochloromethane	129	9.568	9 568 (0.945)	94873	250 000	273 8
68 1,2-Dibromoethane	107	9 678	9.678 (0 956)	121424	250.000	268 4
70 Chlorobenzene	112	10 149	10 149 (1 002)	442746	250.000	266.5
71 1,1,1,2-Tetrachloroethane	131	10 247	10 247 (1.012)	112857	250.000	276.7
72 Ethylbenzene	106	10 234	10.234 (1 011)	25 <b>265</b> 1	250 000	268 8
73 m + p-Xylene	106	10.363	10 363 (1 024)	61171B	500.000	544.6
74 Xylene-o	106	10.754	10 754 (1 062)	300199	250.000	268.1
M 75 Xylenes (total)	106			911917	250.000	814.4
76 Styrene	104	10 785		519441	250 000	275 8
77 Bromoform	173	10 986		50250	250.000	277.0
78 Isopropylbenzene	105	11 109		740920	250.000	273 3
79 Bromobenzene	156	11.445		153219	250 000	261.8
81 n-Propylbenzene	120	11.524		211369	250.000	264 8
82 2-Chlorotoluene	126	11 634		179841	250.000	265.4
83 1,1,2,2-Tetrachloroethane	83	11.488		196523	250 000	268.4
84 1,2,3-Trichloropropane	110	11.537		53924	250 000	270.4
85 4-Chlorotoluene	126	11 750		188006	250.000	268 3
86 1,3,5-Trimethylbenzene	105	11.708		606790	250 000	265.2
87 tert-Butylbenzene	119	12 013	12 013 (0 964)	521697	250 000	266 9

# 657 182

						AMOUN	TS
		QUANT SIG				CAL-AMT	ON-COL
Compo	unds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
30305		====	==		=======	****	
88	1,2,4-Trimethylbenzene	105	12 081	12.081 (0 970)	623203	250 000	266.8
89	sec-Butylbenzene	105	12,233	12.233 (0 982)	849111	250 000	266.3
90	4-Isopropyltoluene	119	12 386	12.386 (0.994)	654793	250.000	275.5
91	1,3-Dichlorobenzene	146	12 386	12 386 (0.994)	316567	250.000	271.3
93	1,4-Dichlorobenzene	146	12 484	12.484 (1 002)	312890	250.000	263.0
94	n-Butylbenzene	91	12 796	12 796 (1 027)	713004	250.000	261 9
95	1,2-Dichlorobenzene	146	12 857	12.857 (1.032)	290497	250.000	264 6
96	1,2-Dibromo-3-chloropropane	157	13.664	13 664 (1 097)	18343	250 000	<b>250</b> 5
97	1,2,4-Trichlorobenzene	180	14 459	14 459 (1 160)	117260	250.000	231 3
98	Hexachlorobutadiene	225	14.581	14 581 (1 170)	55149	250 000	224 8
99	Naphthalene	128	14.709	14 709 (1 181)	241145	250 000	212.6
100	1.2.3-Trichlorobenzene	180	14.960	14.960 (1 201)	81058	250 000	201.8

Page 3

## GC/MS VOLATILE QC DATA

Page 2

Data File: /var/chem/hp4.1/40517d.b/bf40517.d

Date : 17-MAY-2000 10:41

Client ID: 50NGBFB

Sample Info: bfb 1ul

Volume Injected (uL): 1.0

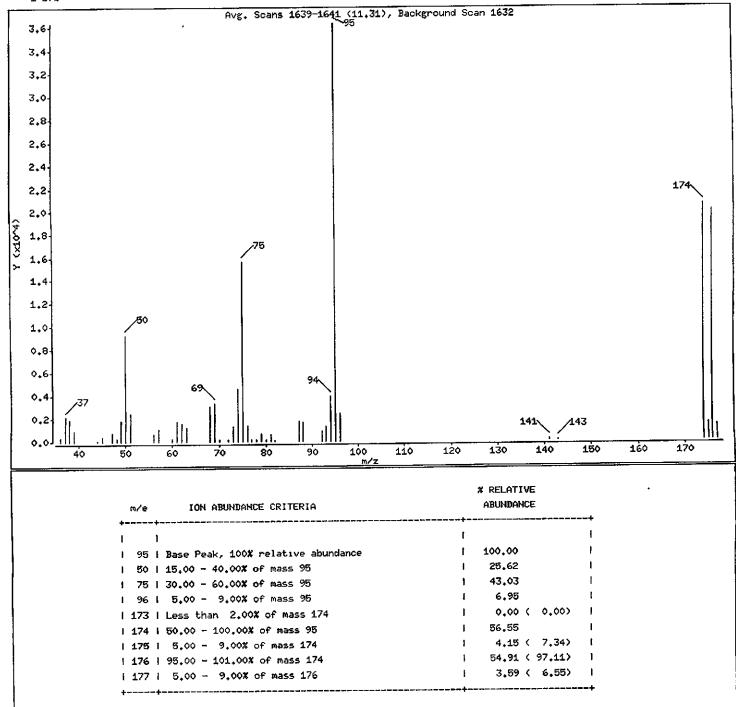
Column phase: DB624 20m

Instrument: hp4.1

Operator: 16328

Column diameter: 0.18

1 bfb



Page 3

Data File: /var/chem/hp4.1/40517d.b/bf40517.d

Date : 17-HAY-2000 10:41

Client ID: 50NGBFB

Instrument: hp4.1

Sample Info: bfb 1ul

Volume Injected (uL): 1.0

Operator: 16328

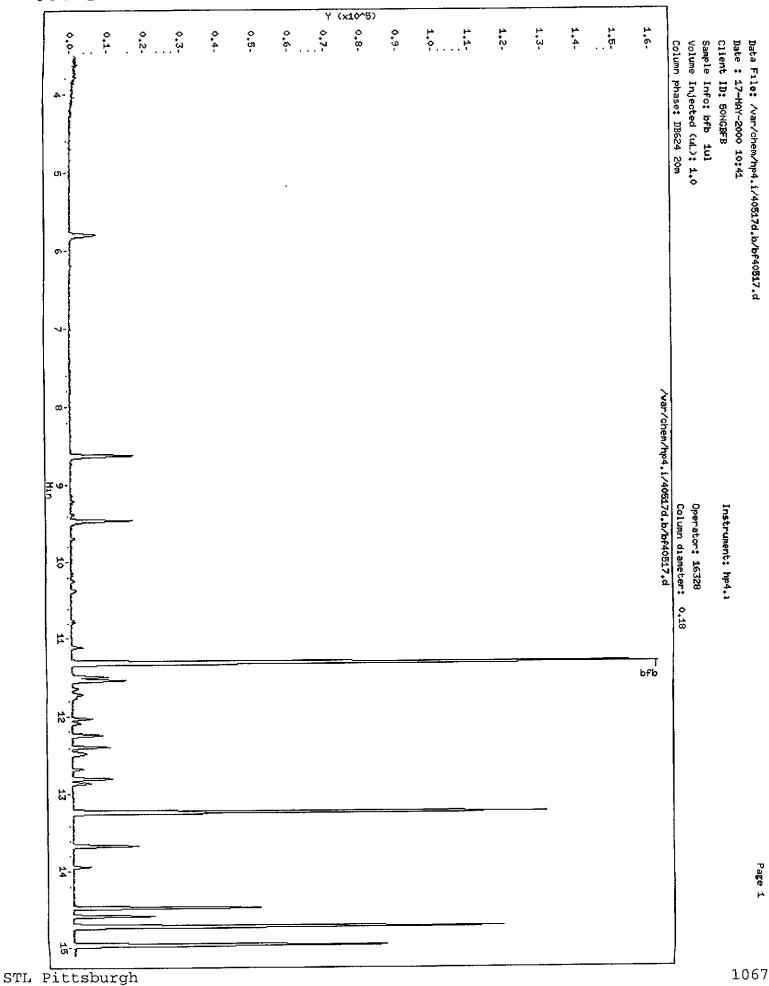
Column phase: DB624 20m Column diameter: 0.18

Data File: bf40517.d

Spectrum: Avg. Scans 1639-1641 (11.31), Background Scan 1632

Location of Maximum: 95.00 Number of points: 44

	m/z	Y		m/z	Y		m/z	Y	m/z	Y	
1	36,00	370	- <b>+</b> -	57.00	1123	•	76.00	1413		36352	•
1	37.00	2167	1	60.00	247	ı	77.00	173 !	96.00	2526	ı
1	38.00	1907	ı	61.00	1740	ı	78.00	147 I	141.00	158	ı
1	39.00	887	1	62.00	1593	1	79.00	638 I	143.00	85	1
1	44,00	93	1	63.00	1219	1	80,00	141 l	174.00	20560	ı
+-			-+-			+-		·			-+
1	45.00	386	ŀ	68.00	3087	1	81.00	618	175.00	1509	1
ı	47.00	760	1	69.00	3325	1	82.00	83	176,00	19960	1
1	48.00	251	i	70.00	208	ł	87,00	1752	177.00	1307	ı
1	49.00	1854	ı	72.00	145	I	88,00	1691			١
1	50,00	9314	ŧ	73,00	1351	ì	92,00	911			ı
+-			-+-			-+-					-+
1	51.00	2548	I	74,00	4618	I	93.00	1356			1
ŧ	56,00	627	I	75.00	15644	1	94,00	3961	l		1
+-			-+-			+-					-+



## UXB INTERNATIONAL METHOD BLANK COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E170000 101

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DD95L101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

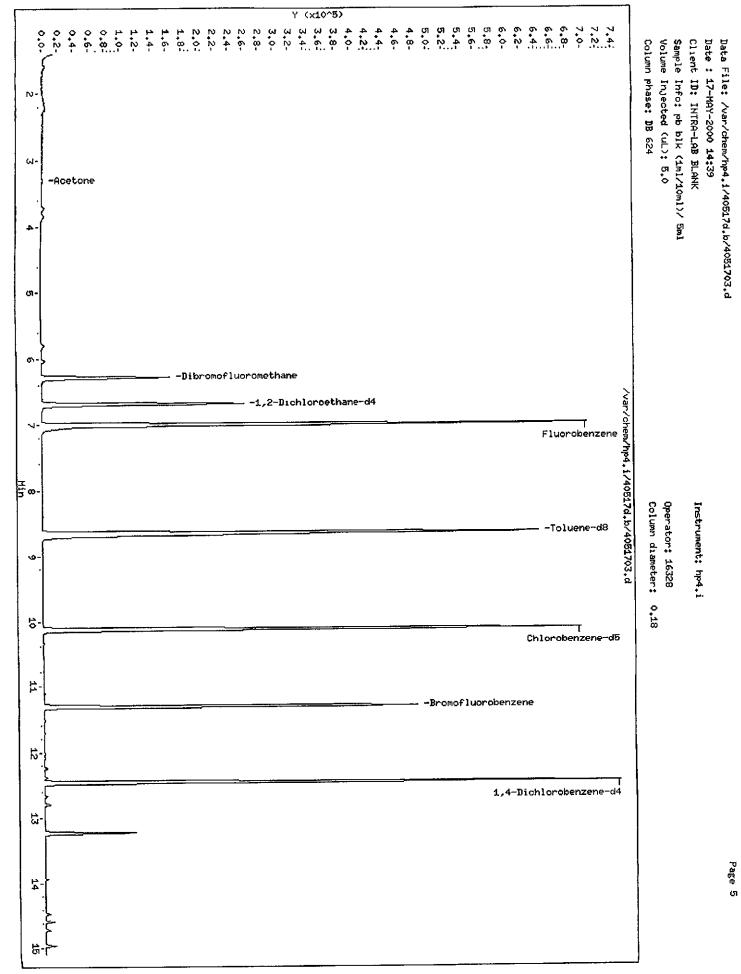
Moisture %:NA

QC Batch: 0138262

Client Sample Id: INTRA-LAB BLANK

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L o	r ug/kg) mg/L 💢	2
71-43-2	Benzene	0.050	U
78-93-3	2-Butanone	0.050	<u> </u>
56-23-5	Carbon tetrachloride	0.050	<u>u</u>
108-90-7	Chlorobenzene	0.050	<u>  u                                   </u>
67-66-3	Chloroform	0.050	<u>""</u>
107-06-2	1,2-Dichloroethane	0.050	<u> </u>
75-35-4	1,1-Dichloroethene	0.050	<u>ات</u> ا
127-18-4	Tetrachloroethene	0.050	<u> </u>
79-01-6	Trichloroethene	0.050	ן ט
75-01-4	Vinyl chloride	0.050	<u>                                     </u>



Page 1

Data File: /var/chem/hp4.i/40517d.b/4051703.d

Report Date: 17-May-2000 14:59

### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file : /var/chem/hp4.i/40517d.b/4051703.d

Lab Smp Id: DD95L101 Client Smp ID: INTRA-LAB BLANK

Inj Date : 17-MAY-2000 14:39 Operator : 16328 Inst ID: hp4.i

Smp Info : pb blk (1ml/10ml)/5ml Misc Info : dd951101,40517d.b,med1000.m

Comment

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m Meth Date : 17-May-2000 14:06 dudeckk Quant 7 Cal Date : 17-MAY-2000 13:03 Cal Fil Als bottle: 11 QC Samp Quant Type: ISTD Cal File: 1e40517.d QC Sample: METHOD BLANK

Dil Factor: 1.00000

Compound Sublist: tclp.sub Integrator: HP RTE

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

Name	Value	Description
DF	1.000	Dilution Factor
Vf	5.000	FinalVolume
Vo	5.000	Sample Volume
Vi	5.000	Purge Volume

D1)5/17/00

						CONCENTRA	TIONS
		QUANT SIG				ON-COLUMN	FINAL
Co	mpounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( UG/L)
21.0	0000000000000000000000			ROUNDO WREEKE	========	*****	
*	46 Fluorobenzene	96	7.006	6.994 (1 000)	792190		
•	69 Chlorobenzene-d5	119	10 124	10 124 (1 000)	154881		
*	92 1,4-Dichlorobenzene-d4	152	12.459	12 460 (1 000)	210752		
\$	43 1,2-Dichloroethane-d4	65	6.688	6 683 (0 955)	192994	260 001	52.00
\$	39 Dibromofluoromethane	113	6 285	6 285 (0 897)	133376	246.080	49.22
\$	59 Toluene-d8	98	8 651	8 645 (0 854)	650896	250.176	50 04
\$	80 Bromofluorobenzene	95	11 304	11 310 (1.117)	225309	224 168	44.83
	3 Vinyl Chloride	62.00	Com	pound Not Detect	ed.		
	12 1,1-Dichloroethene	96 00	Com	pound Not Detect	ed.		
	31 2-Butanone	43 00	Com	pound Not Detect	ed		
	37 Chloroform	83.00	Con	pound Not Detect	eđ.		
	41 Carbon Tetrachloride	117.00	Con	pound Not Detect	ed.		
	42 Benzene	78.00	Con	pound Not Detect	ed.		
	45 1,2-Dichloroethane	62.00	Соп	pound Not Detect	ed		
	47 Trichloroethene	130.00	Con	pound Not Detect	ed		

657 190

Data File: /var/chem/hp4.i/40517d.b/4051703.d

Report Date: 17-May-2000 14:59

CONCENTRATIONS QUANT SIG ON-COLUMN FINAL ( ng) (UG/L) EXP RT REL RT RESPONSE MASS RT Compounds ##=====**#** --------------==== 164.00 Compound Not Detected. 65 Tetrachloroethene 112 00 Compound Not Detected. 70 Chlorobenzene

Page 2

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:COE170000 262

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Date Received: 05/13/00 Work Order: DDA08101 Date Extracted:05/17/00 Dilution factor: 1 Date Analyzed: 05/17/00

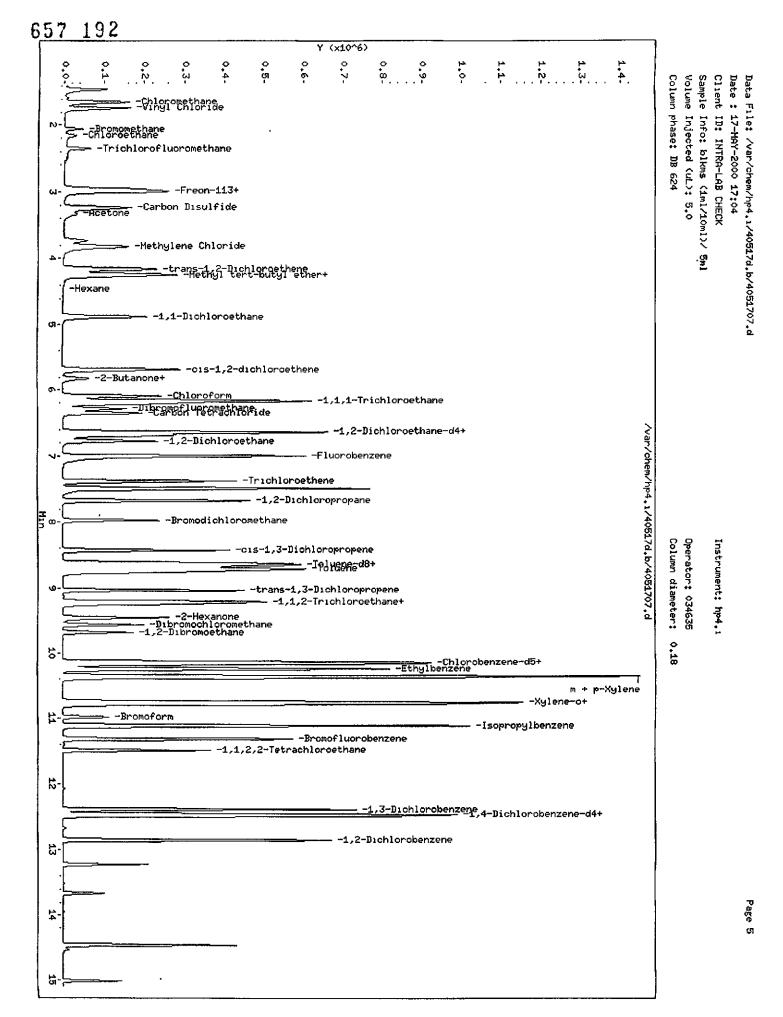
Moisture %:NA

QC Batch: 0138262

Client Sample Id: CHECK SAMPLE

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L o	r ug/kg) mg/L Q
71-43-2	Benzene	0.546
78-93-3	2-Butanone	0.503
56-23-5	Carbon tetrachloride	0.567
108-90-7	Chlorobenzene	0.543
67-66-3	Chloroform	0.540
107-06-2	1,2-Dichloroethane	0.552
75-35-4	1,1-Dichloroethene	0.510
127-18-4	Tetrachloroethene	0.566
79-01-6	Trichloroethene	0.545
75-01-4	Vinyl chloride	0.627



Data File: /var/chem/hp4.i/40517d.b/4051707.d

Report Date: 17-May-2000 17:22

Page 1

#### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file : /var/chem/hp4.i/40517d.b/4051707.d

Lab Smp Id: DDA08101 Client Smp ID: INTRA-LAB CHECK

Inj Date : 17-MAY-2000 17:04

Operator : 034635 Inst ID: hp4.i

Smp Info : blkms (1ml/10ml)/5ml

Misc Info : dda08101,40517d.b, med1000.m

Comment

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m Meth Date : 17-May-2000 14:06 dudeckk Quant 1 Quant Type: ISTD Cal File: 1e40517.d Cal Date : 17-MAY-2000 13:03 Als bottle: 12 QC Sample: METHOD SPIKE

Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: tclp.sub

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

Concentration	FOIMULA: AME	· · DF · (VI/VO)/VI	, \\(\alpha\)
Name	Value	Description	0115/17/10
DF Vf		Dilution Factor FinalVolume	pn /
Vo	5.000	Sample Volume	,
Vi	5.000	Purqe Volume	

						CONCENTRA	ATIONS
		QUANT SIG				ON-COLUMN	FINAL
Compo	unds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( UG/L)
	<b>医医院电影</b> 电电子 医二甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基	====		244220	******		
* 46	Fluorobenzene	96	7 002	6 994 (1.000)	743134		
* 69	Chlorobenzene-d5	119	10.126	10.124 (1.000)	163960		
* 92	1,4-Dichlorobenzene-d4	152	12 455	12 460 (1 000)	229414		
\$ 43	1,2-Dichloroethane-d4	65	6.684	6 683 (0 955)	179634	257.978	51.60
\$ 39	Dibromofluoromethane	113	6.293	6.285 (0.899)	129802	255.295	51.06
\$ 59	Toluene-d8	98	8.653	8.645 (0.855)	718005	260 689	52 14
\$ 80	Bromofluorobenzene	95	11 312	11 310 (1.117)	273445	256 995	51 40
3	Vinyl Chloride	62	1.738	1.743 (0.248)	242371	313.474	62.69
12	1,1-Dichloroethene	96	2 998	2 978 (0 428)	122965	254 899	50.98
31	2-Butanone	43	5 816	5 821 (0.831)	132801	251 360	50 27
37	Chloroform	83	6.091	6 090 (0.870)	249445	270.040	54 01
41	Carbon Tetrachloride	117	6.354	6 352 (0.907)	121574	283 290	56.66
42	Benzene	78	6 641	6.640 (0.948)	740709	272 B96	54.58
45	1,2-Dichloroethane	62	6 776	6.774 (0 968)	208627	276 143	55 23
47	Trichloroethene	130	7 381	7.379 (1.054)	155412	272 481	54 50

657 194 Data File: /var/chem/hp4.i/40517d.b/4051707.d Report Date: 17-May-2000 17:22

Page 2

					CONCENTRA	RIONS
	QUANT SIG				ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	(UG/L)
はなりはは存在時間間間の発展的なればなりでもませる。		===		2222227	=======	*******
65 Tetrachloroethene	164	9 239	9.232 (0 912)	109969	282.784	56 56
70 Chlorobenzene	112	10 156	10.149 (1.003)	493927	271 566	54 31

#### UXB INTERNATIONAL MATRIX SPIKE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc.

SDG Number:

Matrix: (soil/water) SOLID

Lab Sample ID: C0E130194 001

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Work Order: DD6A410P Dilution factor: 1

Moisture %:10

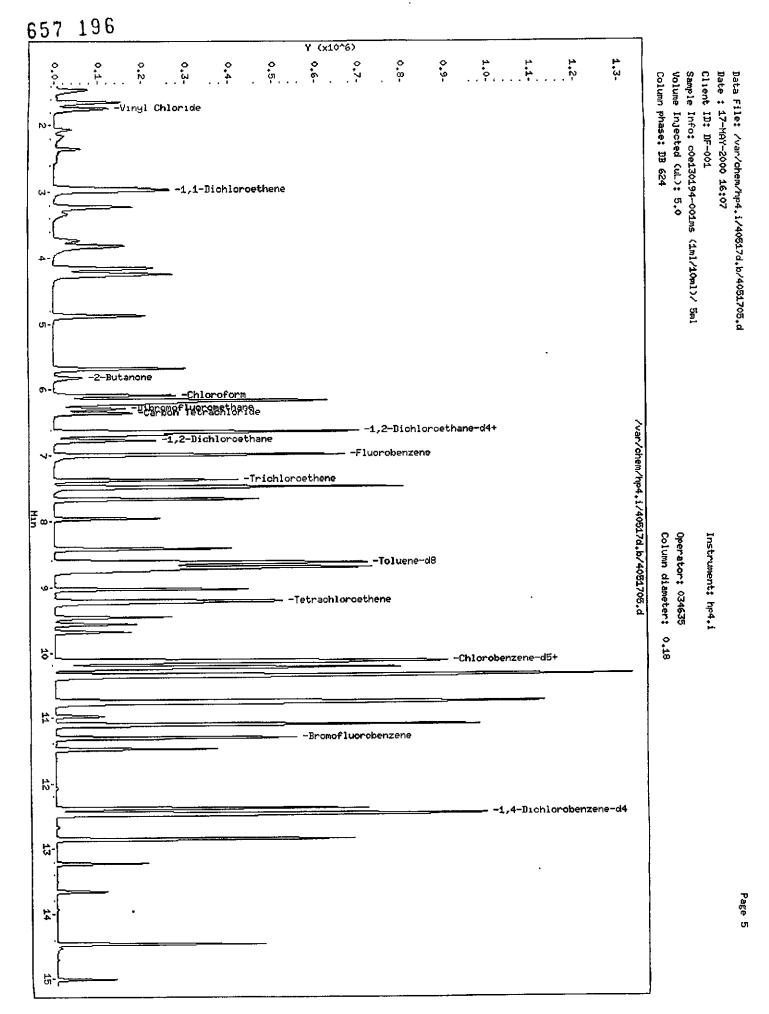
Date Received: 05/13/00 Date Extracted:05/17/00 Date Analyzed: 05/17/00

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 A

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L o	or ug/kg) mg/L Q
71-43-2	Benzene	0.544
78-93-3	2-Butanone	0.532
56-23-5	Carbon tetrachloride	0.563
108-90-7	Chlorobenzene	0.546
67-66-3	Chloroform	0.536
107-06-2	1,2-Dichloroethane	0.552
75-35-4	1,1-Dichloroethene	0.518
127-18-4	Tetrachloroethene	0.555
79-01-6	Trichloroethene	0.545
75-01-4	Vinyl chloride	0.635



Page 1

Data File: /var/chem/hp4.i/40517d.b/4051705.d

Report Date: 17-May-2000 16:52

#### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file : /var/chem/hp4.i/40517d.b/4051705.d

Lab Smp Id: DD6A410P Client Smp ID: DF-001

Inst ID: hp4.i

Inj Date : 17-MAY-2000 16:07

Operator : 034635

Smp Info : c0e130194-001ms (1ml/10ml) / 5ml Misc Info : dd6a410p,40517d.b,med1000.m

Comment

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m Meth Date : 17-May-2000 14:06 dudeckk Quant 7 Cal Date : 17-MAY-2000 13:03 Cal Fil Als bottle: 13 QC Samp Quant Type: ISTD Cal File: 1e40517.d

QC Sample: MS

Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: tclp.sub

Target Version: 3.40 Processing Host: hpuxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

Name	Value	Description
DF	1.000	Dilution Factor
Vf	5.000	FinalVolume
Vo	5.000	Sample Volume
Vi	5.000	Purge Volume

						CONCENTRA	TIONS
		QUANT SIG				ON - COLUMN	FINAL
Co	ompounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( UG/L)
41		PL 18 18 18	**	******	~~~~		
*	46 Fluorobenzene	96	6.995	6.994 (1.000)	736767		
*	69 Chlorobenzene-d5	119	10.125	10.124 (1.000)	159002		
*	92 1,4-Dichlorobenzene-d4	152	12 460	12 460 (1 000)	229690		
\$	43 1,2-Dichloroethane-d4	65	6 683	6 683 (0 955)	176693	255 947	51 19
\$	39 Dibromofluoromethane	113	6.286	6.285 (0.899)	129740	257.378	51.48
\$	59 Toluene-d8	98	8.645	8.645 (0.854)	696789	260.874	52.17
\$	80 Bromofluorobenzene	95	11.311	11.310 (1 117)	264207	256.056	51.21
	3 Vinyl Chloride	62	1.743	1.743 (0.249)	243216	317.285	63.46
	12 1,1-Dichloroethene	96	2 991	2.978 (0.428)	123871	258.997	51.80
	31 2-Butanone	43	5 815	5 821 (0.831)	139356	266.046	53.21
	37 Chloroform	83	6.090	6.090 (0 871)	245577	268.150	53 63
	41 Carbon Tetrachloride	117	6 353	6.352 (0.908)	119678	281.282	56.26
	42 Benzene	78	6.640	6.640 (0.949)	731722	271 914	54.30
	45 1,2-Dichloroethane	62	6 775	6.774 (0.969)	206618	275 847	55 17
	47 Trichloroethene	130	7 380	7 379 (1 055)	153999	272 337	54 47

Data File: /var/chem/hp4.i/40517d.b/4051705.d

Report Date: 17-May-2000 16:52

CONCENTRATIONS QUANT SIG ON-COLUMN FINAL (ng) MASS EXP RT REL RT RESPONSE ( UG/L) RT Compounds -----====== ---------\*\*\*\* 65 Tetrachloroethene 164 9.232 9.232 (0 912) 104638 277.466 55.49 70 Chlorobenzene 112 10.149 10 149 (1 002) 481292 272 871 54.57

Page 2

#### UXB INTERNATIONAL MATRIX SPIKE DUPLICATE COMPOUNDS

657 199

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID

Lab Sample ID: COE130194 001

Method: SW846 8260B

Volatile Organics, GC/MS (8260B)

Sample WT/Vol: 5 / mL Work Order: DD6A410Q Dilution factor: 1

Date Received: 05/13/00 Date Extracted:05/17/00 Date Analyzed: 05/17/00

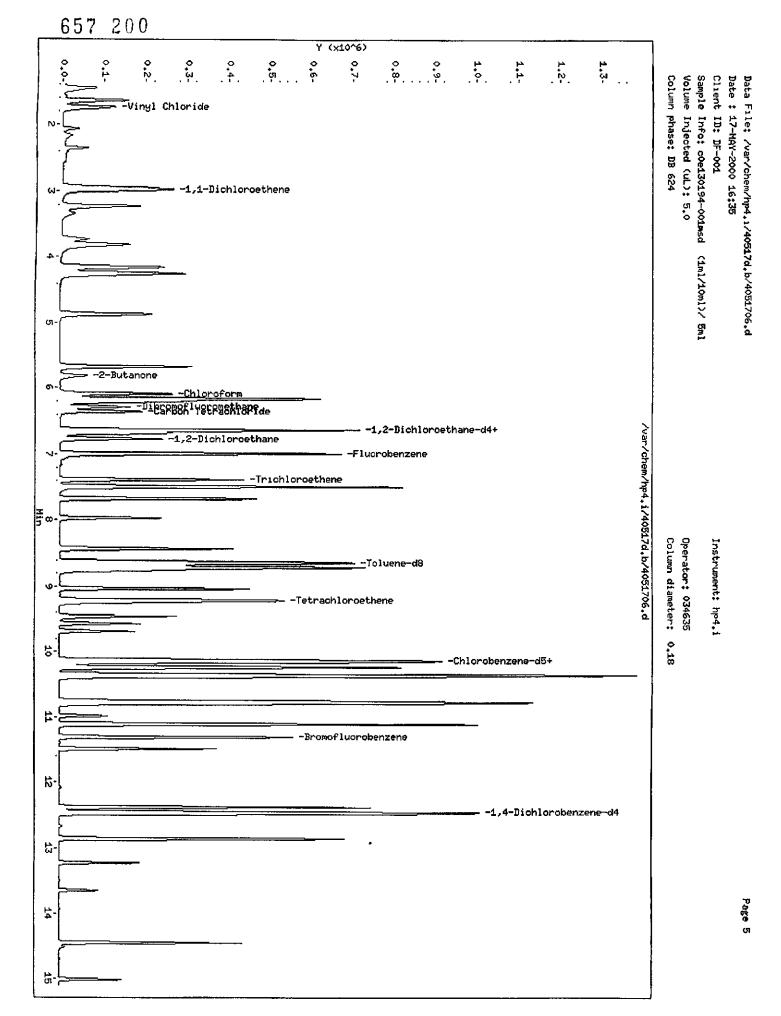
Moisture %:10

QC Batch: 0138262

Client Sample Id: DF/S1/0133/SDC/001 A

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L o	or ug/kg) mg/L Q
71-43-2	Benzene	0.552
78-93-3	2-Butanone	0.526
56-23-5	Carbon tetrachloride	0.571
108-90-7	Chlorobenzene	0.550
67-66-3	Chloroform	0.541
107-06-2	1,2-Dichloroethane	0.561
75-35-4	1,1-Dichloroethene	0.522
127-18-4	Tetrachloroethene	0.567
79-01-6	Trichloroethene	0.551
75-01-4	Vinyl chloride	0.638



657 201 Page 1

Data File: /var/chem/hp4.i/40517d.b/4051706.d

Report Date: 17-May-2000 16:52

#### STL-PITTSBURGH

VOLATILE REPORT SW-846 Method

Data file: /var/chem/hp4.i/40517d.b/4051706.d

Lab Smp Id: DD6A4100 Client Smp ID: DF-001

Inj Date : 17-MAY-2000 16:35

Operator: 034635 Inst ID: hp4.i

Smp Info : c0e130194-001msd (1ml/10ml)/5ml

Misc Info : dd6a410q,40517d.b,med1000.m

Comment :

Method : /var/chem/hp4.i/40517d.b/8260bh2o.m

Meth Date: 17-May-2000 14:06 dudeckk Quant Type: ISTD Cal File: 1e40517.d Cal Date : 17-MAY-2000 13:03

Als bottle: 14 QC Sample: MSD

Dil Factor: 1.00000 Integrator: HP RTE Compound Sublist: tclp.sub

Target Version: 3.40 Processing Host: houxcs21

Concentration Formula: Amt \* DF \* (Vf/Vo)/Vi

Name	Value	Description	. ~
DF Vf Vo Vi	5.000 5.000	Dilution Factor FinalVolume Sample Volume Purge Volume	2/4/2m

						CONCENTRATIONS			
		QUANT SIG	۵			ON-COLUMN	FINAL		
Co	Ompounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	(UG/L)		
24 7	R 双 经 名 名 名 名 立 共 数 名 名 世 B 数 色 世 共 G 名 兰 世 日	25 N G		<b>**********</b>		*******			
*	46 Fluorobenzene	96	6.996	6.994 (1 000)	739353				
*	69 Chlorobenzene-d5	119	10 126	10 124 (1.000)	161550				
*	92 1,4-Dichlorobenzene-d4	152	12.461	12.460 (1.000)	232577				
\$	43 1,2-Dichloroethane-d4	65	6.684	6.683 (0.955)	178328	257 412	51 48		
\$	39 Dibromofluoromethane	113	6 286	6.285 (0.899)	129037	255.088	51.02		
\$	59 Toluene-d8	98	8.646	8.645 (0.854)	696313	256.584	51 32		
\$	80 Bromofluorobenzene	95	11.312	11.310 (1 117)	266398	254.107	50.82		
	3 Vinyl Chloride	62	1.744	1.743 (0.249)	245302	318.887	63.78		
	12 1,1-Dichloroethene	96	2 985	2.978 (0.427)	125297	261.062	52.21		
	31 2-Butanone	43	5 816	5 821 (0 831)	138346	263.194	52 64		
	37 Chloroform	83	6.091	6.090 (0.871)	248439	270.326	54.06		
	41 Carbon Tetrachloride	117	6.354	6.352 (0.908)	121909	285.524	57.10		
	42 Benzene	78	6.641	6.640 (0.949)	745179	275 946	55.19		
	45 1,2-Dichloroethane	62	6.775	6.774 (0.969)	210766	280 401	56 08		
	47 Trichloroethene	130	7.381	7.379 (1.055)	156296	275 433	55 09		

657 202

Data File: /var/chem/hp4.i/40517d.b/4051706.d Report Date: 17-May-2000 16:52

					CONCENTRA	ATIONS
	QUANT SIG				ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	(UG/L)
<b>西班森市</b> 在中央共享经验中国中国中国共和国国际企业		==	03535 BED356	****	355566	
65 Tetrachloroethene	164	9.233	9.232 (0 912)	108643	283 542	56.71
70 Chlorobenzene	112	10 150	10.149 (1.002)	492380	274.754	54 95

Page 2

# GC/MS VOLATILE MISCELLANEOUS

Environmental Services

# Quanterra Environmental Services, Pittsburgh **GC/MS Logsheet**

Sid 192-176-6 Instrument ID: Reviewed by: Sid: 182-180-10 Som Std.: Methanol Lot No .: 192-177-F VOC Date: J-621-081 KAN

22.	20.	19.	18. 1051712	17. 451711	16. <i>本の门o</i>	15. 4051709	14. 4051708	13. 4051707	12. 4051706	11. 405/705	10. 4051764	9. 405/703	8. 40s(702	7. 4051701	6. LE40517	5. 1 D-40517	4. 184017	3. 1A40517		shal 8£40517	
			(DE 130199-065	1200-12014-009	COE 130197-003	008130194-002	(6E130 MA-00)	BILMS	COE (30194-00) Me	COS 130 194-001 mg	SWAR	PB BLK	PB BUK	からた	ort Gas A	\sh\0.12\cdots	くおりか	VSTDS	৴৽য়৽৽	\$F\$	Sample No.
			([mL (0mc)   5mm )	Chillian I ral	( Im (1 Day 1/5 m) 5	( Jing 10mc) 5th 5	[mx/10ml; 5ml/ 5]		(INV//me//sme)			(1-4/10-1) SI S	(12/10-1) 5-0 5	7.5	27	2	5-4	5.2	S-L		
			M.L.	200	Q IN															1401	

### CIEIL Leachate    Continue	10. (16t) S/1/2m 20:20
### Doi: 10.00  ### Doi: 10.00	(6)
### Date   Continue	1.00
### Citent 15   Prinsburgh, PA 15238 Ph Soffer 4 4,60   Pol 280      Lab Sample ID #   Citent ID   Weight   Vessel   Exercation Fluid   Date   Time      Co	6
### Date   Date	Location Date
### Description Fit Way Ph Boffe 4 4,60	
### Date   Department   Date   Stopped   Stopp	
HE Leachate  450 William Pin Way Ph. Boffer, 4 4,60  fethod 1311)  Lab Sample 1D # Client ID   Weight   1	
### PA SO William Pin Way Ph Coffee 7 7.00  Method 1311)  Lab Sample ID # Client ID   Weight   Vesset   Extraction Finid   Date    C. DE 130   94   BIK 4.92   N.   N.   N.   Solid   Date    C. DE 130   94   A.   A.   A.   A.   A.   A.    DOG   6.64   B.   B.    DOG   6.64   B.    DOG   6.64   B.    DOG   6.64   B.    DOG   6.64   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   6.65   B.    DOG   C.   D.    DOG   C.    DOG   C.   D.    DOG   C.    DOG   C.   D.    DOG   C.	
### Client ID   Weight   Vesset   Extraction Finid   Date	
### A50 William Pit Way Ph Boffer 4 4,60  Method 1311)  Lab Sample ID # Client ID   Weight	
### Worksheet   Quanterra Incorporated   450 William Pin Way Ph. Boffer 7 7:00   400   Method 1311)   Client ID   Weight   Vessel   Extraction Fluid   10:01	
### Date   Chamtera Incorporated   450 William Pin Way Ph. Boffer 7 7.00   Polishurgh, PA 15238   Ph. Boffer 7 7.00   Polishurgh Ph. A 15238   Ph. Boffer 7 7.00   Polishurgh Ph. Isoffer 7 7.00   Polishurgh Ph. Isoffer 7 7.00   Polishurgh Ph. Isoffer 7 7.00   Polishurgh Ph. Boffer	
### Vorksheet   Quanterra Incorporated   450 William Pin Way Pl   Boffer   1 7.00   10.01	
### Vorksheet   Quanterra Incorporated   450 William Pitt Way Ph Boffer   4 4,50   Polishurgh, PA 15238 Ph Boffer   7 7.00   Polishurgh   Ph 15238 Ph Boffer   7 7.00   Polishurgh   Ph 15238 Ph Boffer   10 10,0   Polishurgh   Ph 15238 Ph Boffer   10 10,0   Polishurgh   Ph 15238 Ph Boffer   10 10,0   Polishurgh   Poli	
### Worksheet   Quantera Incorporated   450 William Pin Way Ph Boffer 4 4,60   Pittsburgh, PA 15238 Ph Boffer 7 7,00   Pol Boffer 7 7,00   Pol Boffer 7 7,00   Pol Boffer 10 10,01   Ph Boffer 10 10,0	
### THE Leachate    Aso William Pin Way Ph Briffer I 7:00   Political Pittsburgh, PA 15238 Ph Briffer I 7:00   Political Pittsburgh, PA 15238 Ph Briffer I 7:00   Political Ph Briffer I 7	
### THE Leachate  ### Vorksheet  ### Vorksheet  ### Vorksheet  ### Vorksheet  ### Vorksheet  ### Pittsburgh, PA 15238 ph buffer 7 7,000  ### Pittsburgh, PA 15238 ph buffer 7 7,000  ### Pittsburgh, PA 15238 ph buffer 7 7,000  ### Pittsburgh, PA 15238 ph buffer 10 10,01  ### Client ID  ### Client ID  ### Verset  ### Extraction Fluid  ### No. / Volume  ### Date  ### COE130194 & 16.60  ### OO3 6,60  ### OO4 6,660  ### OO4 6,660  ### OO4 6,660  ### OO4 6,660  ### OO4 6,660  ### OO4 6,660  ### OO4 6,660  ### OO4 6,660  #### OO5 OO5 OO5 OO5 OO5 OO5 OO5 OO5 OO5	
### THE Leachate  #### Aso William Pin Way Ph Buffer 4 4,60  #### Worksheet  ##################################	
### Client ID   Weight   Vo. 100   Vo. 1311)    Lab Sample ID #   Client ID   Weight   Vessel   Extraction Fluid   No. / Volume   Date	
### Leachate  Onanterra Incorporated  Fitsburgh, PA 15238 Ph Boffer 7 7,000  Method 1311)  Lab Sample ID # Client ID Weight Vessel Extraction Fluid  COE 130 194 81K 4,92 — JJ TLLP # 1 5/16/2000  OO4 6,62 J B	
HE Leachate  Quanterra Incorporated  450 William Pin Way Ph Buffer 4 4,60  Method 1311)  Lab Sample ID # Client ID Weight 1'essel Extraction Fluid  (gm)  COE 130 194 BIK 4,92 — JJ TELP # 1 5/16/2000 C  002 6,64   N  004 6,68	
HE Leachate  Quanterra Incorporated  Vorksheet  Vorksheet  Pittsburgh, PA 15238  Ph Buffer 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffer 7 7 7,000  Ph Buffe	V
THE Leachate       Quanterra Incorporated         Vorksheet       450 William Pin Way Ph Soffer 7 7.00         Method 1311)       Pittsburgh, PA 15238 Ph Soffer 7 7.00         Method 1311)       Client ID       Weight (gm)       Passet Extraction Fluid No. / Volume Date         COE 130 194       BIK 4.93       35.0       7.2       TELP # 1       5/16/2000       Common Com	
THE Leachate         Quanterra Incorporated           Vorksheet         450 William Pin Way Ph Buffer 7 7,00         4,60           Method 1311)         Piusburgh, PA 15238 Ph Buffer 7 7,00         Police Police           Lab Sample ID #         Client ID         Weight (gm)         Vesset No. / Volume Pate           COE 130 194         BIK 4,92         7         35,0         72         Telf # 1         5/16/2000	
hate  Quanterra Incorporated  450 William Pitt Way Ph Boffer 4 4,60  Pittsburgh, PA 15238 Ph Boffer 7 7,00  Ph Boffer 7 7,00  Ph Boffer 7 7,00  Ph Boffer 7 7,00  Ph Boffer 7 7,00  Ph Boffer 10 10,01	1
Pittsburgh, PA 15238 Ph Buffer 4 4,60 Pittsburgh, PA 15238 Ph Buffer 7 7,00 Ph Buffer 10 10,0)  Th Buffer 10 10,0)	
hate  Quanterra Incorporated 450 William Pin Way Ph Buffer 4,60  Piusburgh, PA 15238 Ph Buffer 7 7,00  Ph Buffer 10 10.01	
Quanterra Incorporated 450 William Pin Way Pl, Buffer 4 4,60 Pittsburgh, PA 15238 Pl, C. F. 7	P01280)
	,

C000067

PSRO24

5/16/00

4:52:05 NT

SAMPLE CUSTODIAN REMOVAL REQUEST

PAGE 001

REQUESTED BY: TROUTE

METHOD: QK

Volatile Organics, GC/MS (82608)

STORAGE LOCATION	WORK ORDER #	PICKED CNTR#	CONTROL #	CLIENT #	ANALYSIS	LOTID	SMP#	<u>SFX</u>	MATRIX DESCRIPTION		Q1 <u>R(</u>	TY QT	
2F CLP1	DD6A4-1-01		233721	399411	A-58-QK	COE130194	001	SOLID		•	0	3	1
2F CLP1	DD6A5-1-01	<del></del>	233722	399411	A-58-QK	COE130194	002	SOLID		e'	0	3	1
2F CLP1	DD6A6-101		233723	399411	A-58-QK	COE130194	003	SOLID			0	3	1
2F CLP1	DD6A7-1-01		233724	399411	A-58-QK	COE130194	004	SOLID			0	3	1

B Front	B Front Deel	DATE/TIME 5/16/2000 07:20 5/16/2000 10:05
•		

GC/MS SEMIVOLATILE DATA

# GC/MS SEMIVOLATILE QC SUMMARY

Lab Name: Severn Trent Laboratories, Inc.

Client: UXB INTERNATIONAL

Lab Code: QESPIT

QESSDG:

Lot #: C0E130194

Ī	CLIENT ID.	SRG01	SRG02	SRG03	SRG04	SRG05	SRG06	TOT OUT
Ī		=======	======	======	======	======	=====	======
01	DF/S1/0133/SDC/001 A	51	56	78	45	47	_56	00
02	DF/S1/0133/SDC/001 B	55	<u>  57                                   </u>	75	49	_51	_59	<u></u>
03	DF/S1/0133/SDC/001 C	47	46	66	42	_45	50	<u>00                                   </u>
04	DF/S1/0133/SDC/001 D	53	54	67	48	51	59	00
	METHOD BLK. DDF2D101	64	68	67	63	_66	66	<u>00</u>
06		65	72	69	61	63	67	00
07	DF/S1/0133/SDC/001 A D	54	58	73	47	51	63	00
	DF/S1/0133/SDC/001 A S	52	57	71	46	49	60	00

SURROGAT	res	QC LIMITS
SRG01	= Nitrobenzene-d5	( 32-112)
SRG02	= 2-Fluorobiphenyl	( 30-110)
SRG03	= Terphenyl-d14	( 10-144)
SRG04	= 2-Fluorophenol	( 13-110)
SRG05	= Phenol-d5	( 10-113)
SRG06	= 2,4,6-Tribromophenol	( 21-122)

<sup>#</sup> Column to be used to flag recovery values

<sup>\*</sup> Values outside of required QC Limits

D System monitoring Compound diluted out

657 210 SW846 8270C MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A

Level: (low/med) LOW

Lot #: C0E130194

WO #: DD6A411G BATCH: 0140269

COMPOUND	SPIKE ADDED (mg/L )	SAMPLE CONCENT. (mg/L )	MS CONCENT. (mg/L )	MS % REC	LIMITS REC	QUAL
	= =======	=======	======	=====		========
1,4-Dichlorobenzene	0.250	ND	0.127	51	18- 110	
2,4-Dinitrotoluene	0.250	ND	0.118	47	31- 131	
Hexachlorobenzene	0.250	ND	0.174	<u> 70_</u>	36- 132	
Hexachlorobutadiene	0.250	ND	0.130	52	18- 116	
Hexachloroethane	0.250	ND	0.118	47_	18- 110	. l
Nitrobenzene	0.250	ND	0.131	52_	10- 211	.1
Pentachlorophenol	0.250	ND	0.138	55	10- 140	·
	0.250	ND	0.157	63	10- 148	
Pyridine	0.250	ND	0.142	57	24- 143	1
2,4,5-Trichlorophenol	0.250	ND	0.140	56	36- 135	
2,4,6-Trichlorophenol Cresols (total)	0.750	ND	0.375	50	25- 144	

# Column to be used to flag recovery and RPD values with an asterisk * Values outside of QC limits	
RPD: 0 out of 0 outside limits  Spike Recovery: 0 out of 11 outside limits	
COMMENTS:	

NOTES (S):

Lab Name. Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A

Level: (low/med) LOW

Lot #: C0E130194

WO #: DD6A411H BATCH: 0140269

	SPIKE ADDED	MSD CONCENT.	MSD %	ş	QC I	LIMITS	
COMPOUND	(mg/L )	(mg/L )	REC	RPD	RPD	REC	QUAL
		========	=====	======	====		=======
1,4-Dichlorobenzene	0.250	0.131	53	3.2	36	18- 110	
2,4-Dinitrotoluene	0.250	0.120	48	1.4	32	31- 131	
Hexachlorobenzene	0.250	0.181	72	4.1	22	36- 132	
Hexachlorobutadiene	0.250	0.135	54	3.6	32	18- 116	
Hexachloroethane	0.250	0.122	49	3.4	33	18- 110	
Nitrobenzene	0.250	0.136	54	3.8	50	10- 211	
Pentachlorophenol	0.250	0.143	57	3.6	56	10- 140	
Pyridine	0.250	0.164	66	4.5	65	10- 148	
2,4,5-Trichlorophenol	0.250	0.146	58	2.4	22	24- 143	
2,4,6-Trichlorophenol	0.250	0.143	57	1.7	27	36- 135	
Cresols (total)	0.750	0.391	52	4.2	33	25- 144	

# Column to be used to flag recovery and RPD values with an asterisk * Values outside of QC limits
RPD: 0 out of 11 outside limits Spike Recovery: 0 out of 11 outside limits
COMMENTS:

NOTES (S):

# 657 212 SW846 8270C CHECK SAMPLE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Lot #: C0E190000

WO #: DDF2D102 BATCH: 0140269

COMPOUND	SPIKE ADDED (mg/L )	SAMPLE CONCENT. (mg/L)	% REC	QC LIMITS REC	    QUAL  =========
1,4-Dichlorobenzene	0.250	0.173	69	28- 110	 
2,4-Dinitrotoluene	0.250	0.136	54	47- 131	
Hexachlorobenzene	0.250	0.191	77	57- 128	
Hexachlorobutadiene	0.250	0.179	72	36- 116	
Hexachloroethane	0.250	0.170	68	30- 110	
Nitrobenzene	0.250	0.168	67	45- 130	.
Pentachlorophenol	0.250	0.162	65	10- 140	.
Pyridine	0.250	0.272	109	10- 148	.
2,4,5-Trichlorophenol	0.250	0.171	69	41- 125	.
2,4,6-Trichlorophenol	0.250	0.173	69	46- 135	.
Cresols (total)	.0.750	0.489	65	29- 144	.

Values outside	of QC	limits			
Spike Recovery:	0	out of	11	outside limit	s
Spike kecovery:	<u></u>	Out OI		0-30	
COMMENTS:				_	

NOTES(S):

#### SW846 8270C METHOD BLANK SUMMARY

BLANK WORKORDER NO. DDF2D101

Lab Name: Severn Trent Laboratories, Inc.

Lab Code: QESPIT

SDG Number:

Lab File ID: S0522006.

Lot Number: C0E130194

Date Analyzed: 05/22/00

Time Analyzed: 12:05

Matrix: SOLID

Date Extracted:05/18/00

GC Column: DB5MS ID: .25

Extraction Method: 1311/3520C

Instrument ID: 71

Level: (low/med) LOW

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, LCS, LCSD, MS, MSD.

_		SAMPLE	LAB	DATE	TIME
	CLIENT ID.	WORK ORDER #	FILE ID	ANALYZED	ANALYZED
		~~~~~~~	=======================================		**======
01	CHECK SAMPLE	DDF2D102 C	S0522007.	05/22/00	12:41
02	DF/S1/0133/SDC/001 A	DD6A4102	S0522008.	05/22/00	13:17
03		DD6A411G S	S0522009.	05/22/00	13:53
04		DD6A411H D	S0522010.	05/22/00	14:29
05	DF/S1/0133/SDC/001 B	DD6A5102	S0522011.	05/22/00	15:05
	DF/S1/0133/SDC/001 C	DD6A6102	S0522012.	05/22/00	15:41
	DF/S1/0133/SDC/001 D	DD6A7102	S0522013.	05/22/00	16:17
08	· ————————————————————————————————————			<b>.</b>	
09					l
10				· .	
11					
12				l	
13				l	<u></u>
14				l	
15				l	l
16				l	
17				·	
18			.	<u></u>	1\
19			,		ll
20			<u> </u>	İ	1
21				l	l1
22				.l <u></u>	\ <u></u>
23			, I	.	l
24				.	
25				\ <u></u>	[
26					l[
27	· · · · · · · · · · · · · · · · · · ·			1	l
28	· ————————————————————————————————————			.}	.
25	·			]	.
30	! ————————————————————————————————————			.	

COMMENTS:	
	FORM TV

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK

657 214

DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name:

Lab Code: Case No.: SAS No.: SDG No.

Contract:

Lab File ID: S0519DF2

DFTPP Injection Date: 05/19/00

Instrument ID: 71

DFTPP Injection Time: 1144

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
=====		======================================
51	30.0 - 60.0% of mass 198	39.6
68	Less than 2.0% of mass 69	0.0 ( 0.0)1
69	Mass 69 relative abundance	39.6
70	Less than 2.0% of mass 69	0.0 ( 0.0)1
127	40.0 - 60.0% of mass 198	49.8
197	Less than 1.0% of mass 198	0.0
198	Base Peak, 100% relative abundance	100.0
199	5.0 to 9.0% of mass 198	6.7
275	10.0 - 30.0% of mass 198	23.3
365	Greater than 1.0% of mass 198	3.83
441	Present, but less than mass 443	14.4
442	Greater than 40.0% of mass 198	88.3
443	17.0 - 23.0% of mass 442	16.3 (18.4)2
	1-Value is % mass 69 2-Value is % mass	

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA	LAB	LAB	DATE	TIME
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED	ANALYZED
	==========		<b>========</b>	=========	========
01	SSTD050	SSTD50	S0519CC3	05/19/00	1320
02	SSTD020	SSTD20	S0519CC2	05/19/00	1356
03	SSTD080	SSTD80	S0519CC4	05/19/00	1432
04	SSTD120	SSTD120	S0519CC5	05/19/00	1508
05	SSTD160	SSTD160	S0519CC6	05/19/00	1544
06					
07					
08					
09					
10					
11					
12			<u> </u>		
13					
14					
15					
16					
17					
18					
19					
20	, <del></del>				
21		—	·		
22		]			

page 1 of 1

FORM V SV

# FORM 5 SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name:

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.: C0E130194

Lab File ID: S0522DF1

DFTPP Injection Date: 05/22/00

Instrument ID: 71

DFTPP Injection Time: 1109

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51 68 69 70 127 197 198 199 275 365 441 442	30.0 - 60.0% of mass 198 Less than 2.0% of mass 69 Mass 69 relative abundance Less than 2.0% of mass 69 40.0 - 60.0% of mass 198 Less than 1.0% of mass 198 Base Peak, 100% relative abundance 5.0 to 9.0% of mass 198 10.0 - 30.0% of mass 198 Greater than 1.0% of mass 198 Present, but less than mass 443 Greater than 40.0% of mass 198	36.7 0.0 ( 0.0)1 38.0 0.0 ( 0.0)1 47.9 0.0 100.0 6.9 23.1 3.59 13.7 84.8
443	17.0 - 23.0% of mass 442	16.7 (19.7)2

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA	LAB	TAD	T) A (III)	CT NOT
	,		LAB	DATE	TIME
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED	ANALYZED
			#G=====# <b>E</b>	========	========
01	SSTD050	SSTD50	S0522CCC	05/22/00	1127
02	INTRA-LAB BL	DDF2D101	S0522006	05/22/00	1205
03	INTRA-LAB CH	DDF2D102	S0522007	05/22/00	1241
04	DF/S1/0133/S	DD6A4102	S0522008	05/22/00	1317
MS 05			S0522009	05/22/00	1353
MSD 06			S0522010	05/22/00	1429
07	DF/S1/0133/S		S0522011	05/22/00	1505
08			S0522012	05/22/00	1541
09	DF/S1/0133/S		S0522012	05/22/00	1617
10	01/01/0133/0	BBORTIOZ	150322013	05/22/00	101,
11	<u> </u>		<u> </u>		l
12					
13					
14	<u> </u>		<u> </u>		
15	\				
16					
17	\				
18					
19					<del></del>
20				<del></del>	
21					
22				l ————	
22	1		l	l	1

page 1 of 1

FORM V SV

#### FORM 8 SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

657 216

Lab Name: Contract:

Lab Code: Case No.: SAS No.: SDG No.: C0E130194

Lab File ID (Standard): S0522CCC Date Analyzed: 05/22/00

Instrument ID: 71 Time Analyzed: 1127

							1
\ <del></del>		IS1 (DCB)		IS2 (NPT)		IS3 (ANT)	1
<u> </u>	ļ		#  RT #	AREA #	RT #	AREA #	RT #
l			=   =======	=======	======	=======	======
12 170	UR STD	145518	5.03	535078	6.56	265064	9.60
				1070156	7.06	530128	10.10
	LIMIT	291036		267539	6.06	132532	9.10
LOWER	LIMIT	72759	4.53	207539	0.00	152558	
=====	=====	========	=   ======	========	======		
) CLI	ENT				1		
SAMPL	e no. Ì						
=====	=====	========	:= =====	=======	======	========	======
01 INTRA-	LAB BL	161868	5.03	630828	6.57	338411	9.60
	LAB CH	161298		635000	6.56	326610	9.60
		171047		654778	6.56	330853	9.60
03 DF/S1/	0133/S		L	633807	6.57	317069	9.60
04 DF/S1/	0133/S	165477	2	635621	6.57	322693	9.60
05 DF/S1/	0133/S	MSD 165720			6.56	312452	9.60
06 DF/S1/	0133/S	162741		618658			9.59
07 DF/S1/	0133/S	171377		653524	6.56	328496	
08 DF/S1/	0133/S	155709	5.03	593533	6.56	302566	9.59
09			ĺ	<u></u>			
10					l		l
11							_
<u> </u>					\ <del></del>		
12			<del></del>	·	· [		
13			\- <del></del>		·   <del></del>	<del></del>	
14				.	.	·   <del></del>	
15			\			·	
16					.		ļ <del></del>
17				.   <u></u>		.	
18							.
19					.	.	.\
20							·
21				-			
				-	-		
22		.		_	_ I <del></del>	- I	1

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8
IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area AREA LOWER LIMIT = - 50% of internal standard area

RT UPPER LIMIT = + 0.50 minutes of internal standard RT RT LOWER LIMIT = - 0.50 minutes of internal standard RT

# Column used to flag internal standard area values with an asterisk.

\* Values outside of QC limits.

page 1 of 1

FORM VIII SV

657 217

Lab Name: Contract:

Lab Code: Case No.: SAS No.: SDG No.: C0E130194

Lab File ID (Standard): S0522CCC Date Analyzed: 05/22/00

Instrument ID: 71 Time Analyzed: 1127

ļ		IS4 (PHN)		IS5 (CRY)		IS6 (PRY)	<u> </u>
		AREA #	RT #	AREA #	RT #	AREA #	RT #
		427212	12.90	446724	19.53	599590	22 00
	12 HOUR STD UPPER LIMIT	854424	13.40	893448	20.03	1199180	22.89 23.39
	LOWER LIMIT	213606	12.40	223362	19.03	299795	22.39
	<b>a=====</b> =====	========	======	=========		======================================	======
	CLIENT						
	SAMPLE NO.			!			
01	INTRA-LAB BL	564305	12.90	527745	19.53	648566	22.89
02	INTRA-LAB CH	529319	12.90	491727	19.53	616566	22.89
03	DF/S1/0133/S	540847	12.90	508153	19.53	598482	22.89
04	DF/S1/0133/S	M5 528651	12.90	549885	19.53	656369	22.90
05	DF/S1/0133/S	MSD 522771	12.90	519636	19.53	634789	22.90
06	DF/S1/0133/S	506789	12.89	498685	19.52	606512	22.88
07	DF/S1/0133/S	524299	12.89	498960	19.52 19.52	611014 603824	22.88
08 09	DF/S1/0133/S	486337	12.89	495522	19.52	003024	22.88
10			ļ ———				\
11							
12							
13		<u> </u>					
14 15						<u></u>	<u> </u>
16			<u> </u>			·-	
17	\	\- <del></del>					
18			'				
19							
20							<u> </u>
21 22	<u> </u>	\					
44			·		l		I

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12 IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area AREA LOWER LIMIT = - 50% of internal standard area

RT UPPER LIMIT = + 0.50 minutes of internal standard RT RT LOWER LIMIT = - 0.50 minutes of internal standard RT

# Column used to flag internal standard area values with an asterisk.

\* Values outside of QC limits.

page 1 of 1

FORM VIII SV

# GC/MS SEMIVOLATILE SAMPLE DATA

#### UXB INTERNATIONAL

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 001

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DD6A4102 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

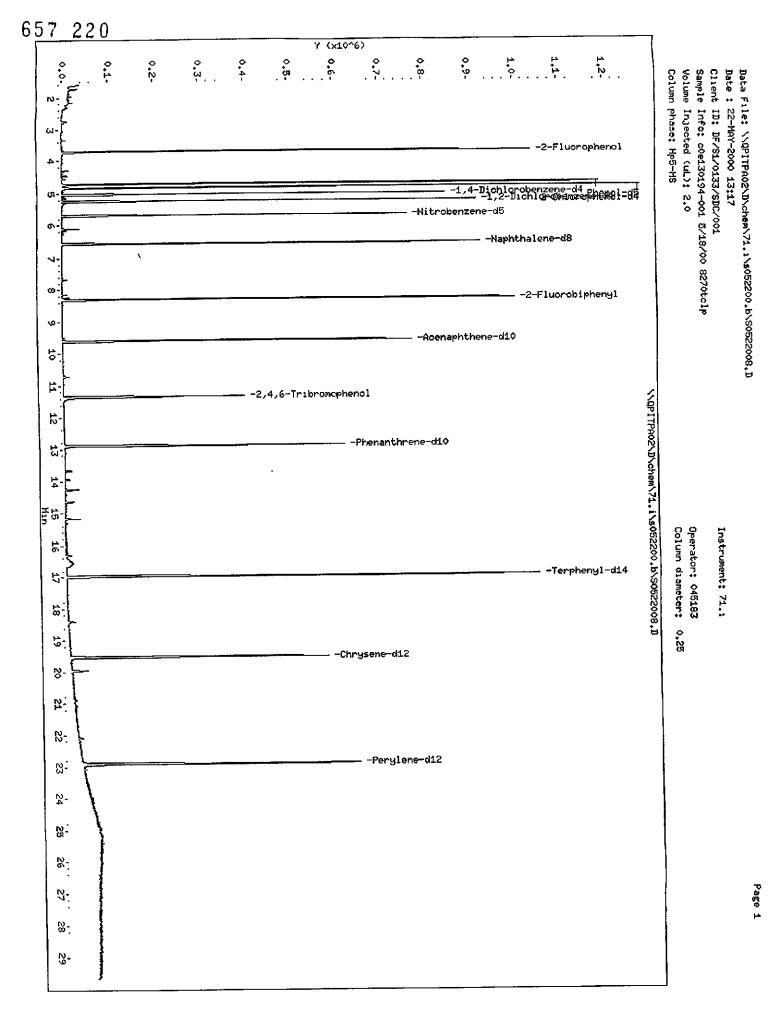
Moisture %:10

QC Batch: 0140269

Client Sample Id: DF/S1/0133/SDC/001 A

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L or	r ug/kg) mg/L	Q
106-46-7	1,4-Dichlorobenzene	0.050	<u>  U                                   </u>
121-14-2	2,4-Dinitrotoluene	0.050	<u>  u</u>
118-74-1	Hexachlorobenzene	0.050	<u>  U                                   </u>
87-68-3	Hexachlorobutadiene	0.050	<u>  u</u>
67-72-1	Hexachloroethane	0.050	<u>  U                                   </u>
98-95-3	Nitrobenzene	0.050	ַןו
87-86-5	Pentachlorophenol	0.25	<u>  u   </u>
110-86-1	Pyridine	0.10	<u>  u</u>
95-95-4	2,4,5-Trichlorophenol	0.050	<u>  </u>
88-06-2	2,4,6-Trichlorophenol	0.050	<u></u>
1319-77-3	Cresols (total)	0.050	<u>  u   </u>



Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522008.D

Report Date: 22-May-2000 13:52

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file: \\QPITPA02\D\chem\71.i\s052200.b\S0522008.D

Lab Smp Id: DD6A4102 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 22-MAY-2000 13:17

Operator : 045183 Inst ID: 71.i

Smp Info : c0e130194-001 5/18/00 8270tclp Misc Info : dd6a4102,s052200.b,8270clp.m,1-tclp.sub

Comment

Method : \\QPITPA02\D\chem\71.i\s052200.b\8270clp.m Meth Date : 22-May-2000 12:37 bachas Quant Type: I Ouant Type: ISTD Cal File: S0519CC6.D Cal Date : 19-MAY-2000 15:44

Als bottle: 10

Dil Factor: 1.00000 Integrator: HP RTE

Compound Sublist: 1-tclp.sub

Target Version: 4.03 Processing Host: PITPC050

Concentration Formula: Amt \* DF \* Uf \* Vt/(Vo \* Vi)

1/12/10

Name	Value	Description /
DF Uf Vt Vo Vi	0.001 1000.000 200.000	Dilution Factor ng unit correction factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected (uL)

						CONCENTRA	ations
	QUANT SIG					ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	( ng)	( mg/L)
				****	*****		**====
* 1 1,4-Dichlorobenzene-d4	152	5.035	5.029	(1,000)	171047	40.0000	(a)
* 2 Naphthalene-d8	136	6.563	6 562	(1.000)	654778	40.0000	(a)
* 3 Acenaphthene-dl0	164	9.597	9 596	(1 000)	330853	40.0000	(a)
* 4 Phenanthrene-d10	188	12.898	12.898	(1.000)	540847	40.0000	(a)
* 5 Chrysene-dl2	240	19 528	19.533	(1 000)	508153	40 0000	(a)
* 6 Perylene-d12	264	22.888	22.893	(1.000)	598482	40 0000	(a)
10 Pyridine	79	Con	pound No	t Detect	ed.		
28 1,4-Dichlorobenzene	146	Con	pound No	ot Detect	ed.		
M 34 Cresols, total	100	Con	pound No	ot Detect	ed.		
31 2-Methylphenol	108	Соп	pound No	ot Detect	ed		
35 4-Methylphenol	108	Con	pound No	ot Detect	ed		
38 Hexachloroethane	117	Con	npound No	ot Detect	ed		
39 Nitrobenzene	7 <b>7</b>	Cor	mpound No	ot Detect	ed.		
59 Hexachlorobutadiene	224	Cor	npound No	ot Detect	ed		

			CONCENTRA	ATIONS
	QUANT SIG		ON-COLUMN	FINAL
Compounds	MASS	RT EXP RT REL RT RESPONSE	( ng)	( mg/L)
医拉德森氏征 医医克尔氏氏 计二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十	2 A 4		8888888	
69 2,4,6-Trichlorophenol	196	Compound Not Detected.		
70 2,4,5-Trichlorophenol	196	Compound Not Detected		
91 2,4-Dinitrotoluene	165	Compound Not Detected		
113 Hexachlorobenzene	283	Compound Not Detected		
117 Pentachlorophenol	265	Compound Not Detected		
\$ 172 Nitrobenzene-d5	82	5 686 5.686 (0.866) 340374	51 4051	0 12851(a)
\$ 173 2-Fluorobiphenyl	172	8 309 8.309 (0 866) 559981	55 7 <b>967</b>	0 13949(a)
\$ 174 Terphenyl-d14	244	16 974 16.969 (0.869) 886750	78 0351	0.19509(a)
\$ 175 Phenol-dS	99	4 714 4.719 (0 936) 534844	70 0950	0.17524 (a)
\$ 176 2-Fluorophenol	112	3 704 3.709 (0 736) 381427	66.8026	0 16701(a)
\$ 177 2,4,6-Tribromophenol	330	11 344 11.338 (0 879) 112909	84.0900	0 21022(a)
\$ 178 2-Chlorophenol-d4	132	4.832 4 831 (0 960) 444824	83.2135	0.20803(a)
\$ 179 1.2-Dichlorobenzene-d4	152	5.243 5.237 (1 041) 183143	48.7647	0.12191(a)

#### QC Flag Legend

#### UXB INTERNATIONAL

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 002

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DD6A5102 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

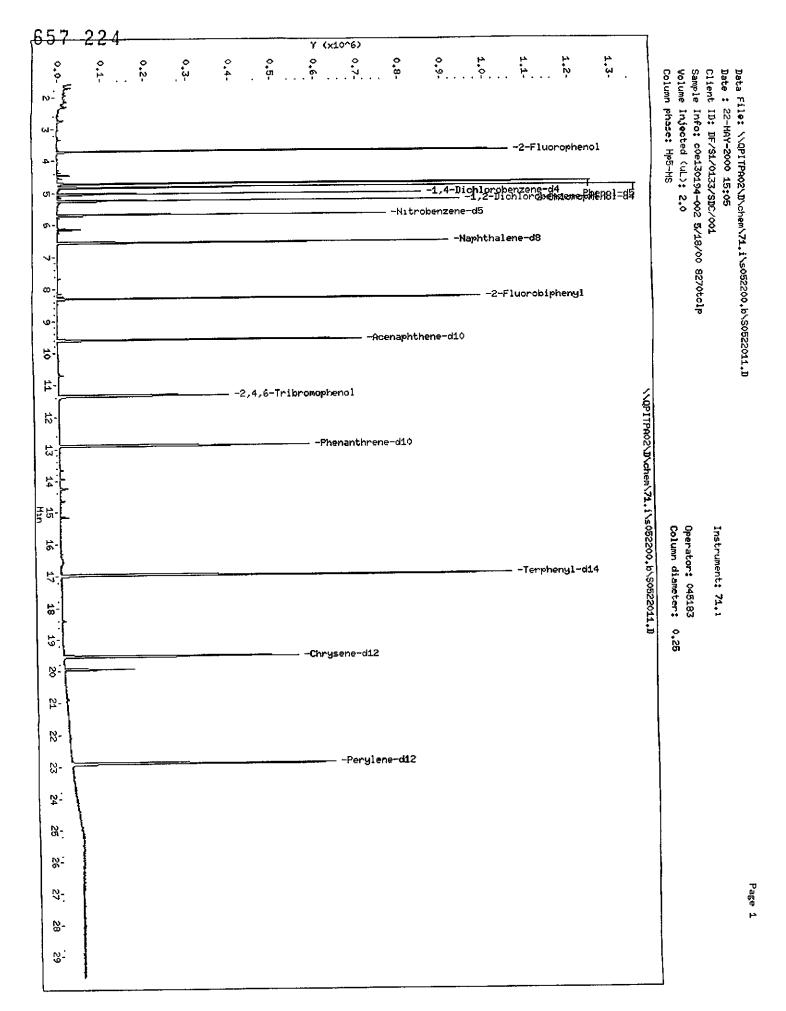
Moisture %:8.6

QC Batch: 0140269

Client Sample Id: DF/S1/0133/SDC/001 B

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L or	r ug/kg) mg/L	Q
106-46-7	1,4-Dichlorobenzene	0.050	<u>ات</u> ا
121-14-2	2,4-Dinitrotoluene	0.050	اتا_
118-74-1	Hexachlorobenzene	0.050	<u>  U                                   </u>
87-68-3	Hexachlorobutadiene	0.050	<u>  u   </u>
67-72-1	Hexachloroethane	0.050	<u>  U</u>
98-95-3	Nitrobenzene	0.050	ן
87-86-5	Pentachlorophenol	0.25	ן ט
110-86-1	Pyridine	0.10	ן
95-95-4	2,4,5-Trichlorophenol	0.050	<u>                                     </u>
88-06-2	2,4,6-Trichlorophenol	0.050	_  <u> </u>
1319-77-3	Cresols (total)	0.050	ا <u>ت</u> ا



Page 1

Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522011.D

Report Date: 22-May-2000 15:36

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file : \\QPITPA02\D\chem\71.i\s052200.b\S0522011.D

Lab Smp Id: DD6A5102 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 22-MAY-2000 15:05

Operator : 045183 Inst ID: 71.i

Smp Info : c0e130194-002 5/18/00 8270tclp

Misc Info : dd6a5102,s052200.b,8270clp.m,1-tclp.sub

Comment

Method : \\QPITPA02\D\chem\71.i\s052200.b\8270clp.m Meth Date : 22-May-2000 12:37 bachas Quant Type: I Quant Type: ISTD Cal Date : 19-MAŸ-2000 15:44 Cal File: S0519CC6.D

Als bottle: 13 Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: 1-tclp.sub

Target Version: Processing Host: PITPC050

Concentration	Formula: Amt	* DF * Uf * Vt/(Vo * Vi)	S\$1201W
Name	Value	Description	
DF	1.000	Dilution Factor	
Uf	0.001	ng unit correction factor	
Vt		Volume of final extract (uL)	
Vo	200.000	Volume of sample extracted (mi	۲)
Vi	2.000	Volume injected (uL)	

						CONCENTRA	TIONS
	QUANT SIG					ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	( ng)	( mg/L)
	B파트	<b>==</b>		000000	=====	======	
* 1 1,4-Dichlorobenzene-d4	152	5.031	5 029	(1 000)	162741	40.0000	(a)
* 2 Naphthalene-d8	136	6.564	6 562	(1.000)	618658	40.0000	(a)
* 3 Acenaphthene-d10	164	9 598	9.596	(1 000)	312452	40 0000	(a)
* 4 Phenanthrene-d10	188	12.894	12.898	(1 000)	506789	40.0000	(a)
* 5 Chrysene-d12	240	19 524	19.533	(1.000)	498685	40 0000	(a)
* 6 Perylene-d12	264	22.884	22.893	(1.000)	606512	40 0000	(a)
10 Pyridine	79	Con	npound No	ot Detect	ed		
28 1,4-Dichlorobenzene	146	Con	mpound No	ot Detect	ed.		
M 34 Cresols, total	100	Con	mpound No	ot Detect	ed.		
31 2-Methylphenol	108	Cor	npound No	ot Detect	ed		
35 4-Methylphenol	108	Cor	npound No	ot Detect	ed		
38 Hexachloroethane	117	Cor	npound No	ot Detect	ed.		
39 Nitrobenzene	77	Cot	mpound No	ot Detect	eđ.		
59 Hexachlorobutadiene	224	Cor	mpound No	ot Detect	ed.		

Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522011.D Report Date: 22-May-2000 15:36

			CONCENTRA	TIONS
	QUANT SIG		ON - COLUMN	FINAL
Compounds	RASS	RT EXP RT REL RT RESPONS	E ( ng)	( mg/L)
*******			5 5 6 5 5 5 T T T	======
69 2,4,6-Trichlorophenol	196	Compound Not Detected		
70 2,4,5-Trichlorophenol	196	Compound Not Detected		
91 2,4-Dinitrotoluene	165	Compound Not Detected.		
113 Hexachlorobenzene	283	Compound Not Detected		
117 Pentachlorophenol	265	Compound Not Detected.		
\$ 172 Nitrobenzene-d5	82	5.688 5 686 (0.867) 34269	8 54.7779	0.13694 (a)
\$ 173 2-Fluorobiphenyl	172	8.306 8.309 (0.865) 53673	2 56.6298	0.14157(a)
\$ 174 Terphenyl-d14	244	16 971 16 969 (0 869) 83971	5 75.2988	0 18825(a)
\$ 175 Phenol-d5	99	4.716 4 719 (0 937) 55520	9 76.4778	0.19119(a)
\$ 176 2-Fluorophenol	112	3 706 3.709 (0.737) 39610	7 72.9145	0 18229(a)
\$ 177 2,4,6-Tribromophenol	330	11.340 11 338 (0.879) 11163	2 88.7260	0 22182(a)
\$ 178 2-Chlorophenol-d4	132	4 828 4.831 (0.960) 44935	1 88.3507	0.22088(a)
\$ 179 1,2-Dichlorobenzene-d4	152	5 239 5.237 (1.041) 17882	2 50.0442	0 12511(a)

#### QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).

#### UXB INTERNATIONAL

Lab Name:Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E130194 003

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DD6A6102 Date Extracted:05/18/00 Date Analyzed: 05/22/00

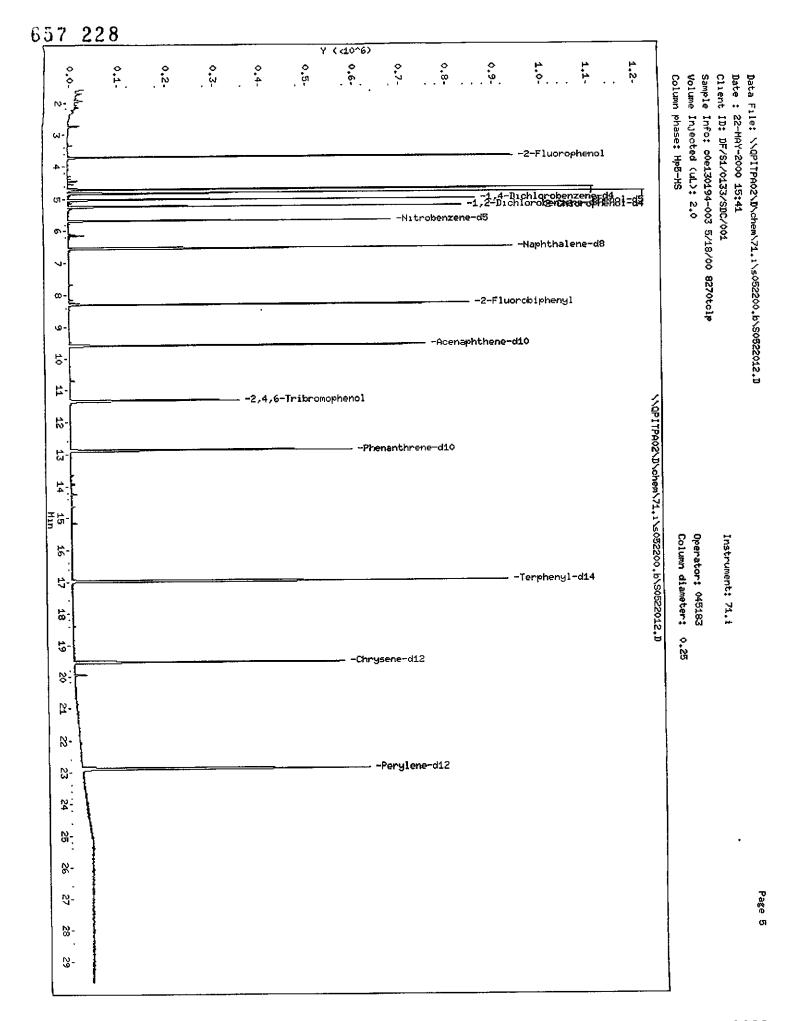
Moisture %:10

QC Batch: 0140269

Client Sample Id: DF/S1/0133/SDC/001 C

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L Q	
106-46-7	1,4-Dichlorobenzene	0.050	ַן
121-14-2	2,4-Dinitrotoluene	0.050	<u> </u>
118-74-1	Hexachlorobenzene	0.050	ַ ַ ַ ַ ַ
87-68-3	Hexachlorobutadiene	0.050	<u>U</u>
67-72-1	Hexachloroethane	0.050	<u> </u>
98-95-3	Nitrobenzene	\0.050	ַ ט
87-86-5	Pentachlorophenol	0.25	ןַּט
110-86-1	Pyridine	0.10	<u> </u>
95-95-4	2,4,5-Trichlorophenol	0.050	<u> </u>
88-06-2	2,4,6-Trichlorophenol	0.050	<u> </u>
1319-77-3	Cresols (total)	0.050	ן ע



Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522012.D Report Date: 22-May-2000 16:43

Page 1

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file: \\QPITPA02\D\chem\71.i\s052200.b\S0522012.D Lab Smp Id: DD6A6102 Client Smp ID: Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 22-MAY-2000 15:41

Operator : 045183 Inst ID: 71.i

Smp Info : c0e130194-003 5/18/00 8270tclp

Misc Info : dd6a6102,s052200.b,8270clp.m,1-tclp.sub

Comment

: \\QPITPA02\D\chem\71.i\s052200.b\8270clp.m Method Meth Date : 22-May-2000 12:37 bachas Quant Type: ISTD Cal Date : 19-MAY-2000 15:44 Cal File: S0519CC6.D

Als bottle: 14 Dil Factor: 1.00000 Integrator: HP RTE

Target Version: 4.03

Processing Host: PITPC050

Compound Sublist: 1-tclp.sub

Concentration Formula: Amt \* DF \* Uf \* Vt/(Vo \* Vi)

Name	Value	Description
DF Uf Vt Vo Vi	0.001 1000.000 200.000	Dilution Factor ng unit correction factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected (uL)

						CONCENTR	ATIONS
	QUANT SIG					ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	( ng)	( mg/L)
	***	-	=====	56550=	*******		222255
* 1 1,4-Dichlorobenzer	e-d4 152	5.029	5 029	(1.000)	171377	40 0000	(a)
* 2 Naphthalene-d9	136	6.557	6.562	(1.000)	653524	40.0000	(a)
<ul> <li>3 Acenaphthene-d10</li> </ul>	164	9.592	9 596	(1 000)	328496	40.0000	(a)
* 4 Phenanthrene-d10	188	12.893	12 898	(1.000)	524299	40 0000	(a)
* 5 Chrysene-dl2	240	19.517	19 533	(1 000)	498960	40.0000	(a)
* 6 Perylene-dl2	264	22 877	22 893	(1.000)	611014	40 0000	(a)
10 Pyridine	79	Cor	mpound N	ot Detect	ed.		
28 1,4-Dichlorobenzer	ne 146	Cor	npound No	ot Detect	eđ		
M 34 Cresols, total	100	Cor	mpound N	ot Detect	ed		
31 2-Methylphenol	108	Cor	mpound N	ot Detect	.ed.		
35 4-Methylphenol	108	Cor	mpound N	ot Detect	.ed.		
38 Hexachloroethane	117	Cot	mpound N	ot Detect	ed.		
39 Nitrobenzene	77	Cor	M baroqm	ot Detect	ed.		
59 Hexachlorobutadies	ne 224	Cor	mpound N	ot Detect	ed:		

		CONCENT	RATIONS
	QUANT SIG	ON-COLUMN	FINAL
Compounds	MASS	RT EXP RT REL RT RESPONSE ( ng)	(mg/L)
	****	三字 医医医内侧 医双双双环 异环异乙基环己烷 克尔巴亚亚亚	*****
69 2,4,6-Trichlorophenol	196	Compound Not Detected.	
70 2,4,5-Trichlorophenol	196	Compound Not Detected	
91 2,4-Dinitrotoluene	165	Compound Not Detected.	
113 Hexachlorobenzene	283	Compound Not Detected	
117 Pentachlorophenol	265	Compound Not Detected.	
\$ 172 Nitrobenzene-d5	82	5.681 5.686 (0.866) 312190 47 2390	0 11810(a)
\$ 173 2-Fluorobiphenyl	172	8.304 8 309 (0.866) 460044 46.1679	0.11542(a)
\$ 174 Terphenyl-d14	244	16 964 16.969 (0.869) 736170 65 9774	0.16494(a)
\$ 175 Phenol-d5	99	4 714 4.719 (0.937) 515789 67 4677	0 16867(a)
\$ 176 2-Fluorophenol	112	3 705 3 709 (0 737) 360711 63.0529	0.15763(a)
\$ 177 2.4.6-Tribromophenol	330	11.333 11.338 (0 879) 98392 75.5909	0.18898(a)
\$ 178 2-Chlorophenol-d4	132	4 826 4.831 (0.960) 410905 76.7204	0.19180(a)
\$ 179 1.2-Dichlorobenzene-d4	152	5.238 5.237 (1.041) 157506 41.8576	0.10464(a)

#### QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).

#### UXB INTERNATIONAL

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:COE130194 004

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DD6A7102 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

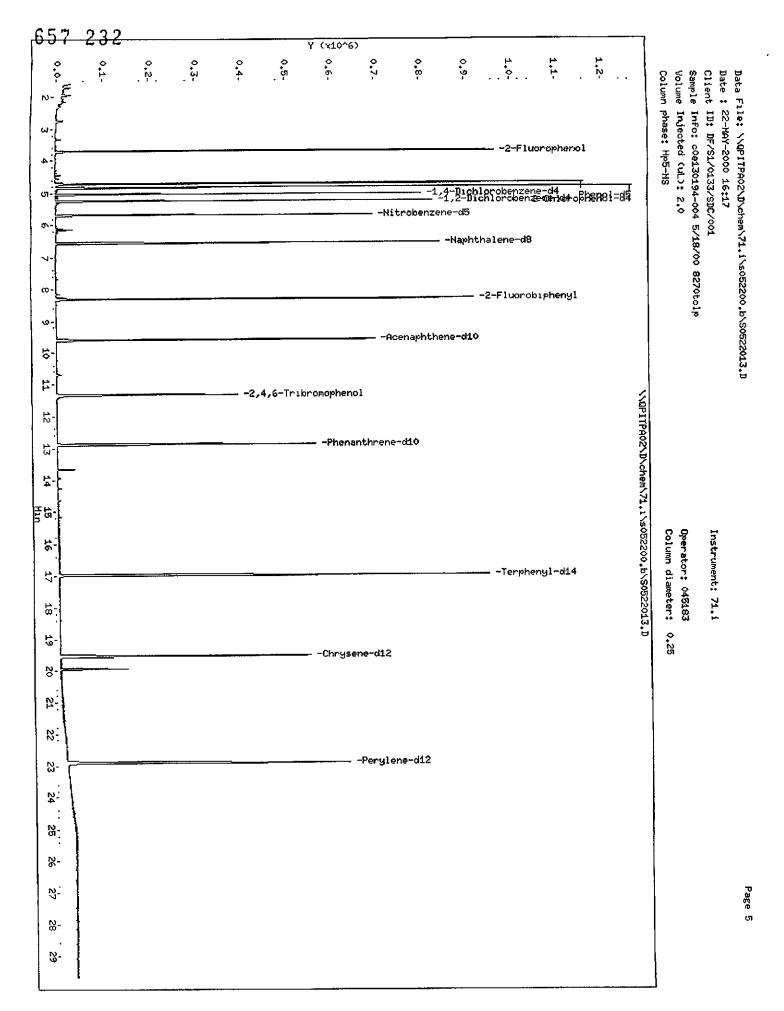
Moisture %:11

QC Batch: 0140269

Client Sample Id: DF/S1/0133/SDC/001 D

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L	Q
106-46-7	1,4-Dichlorobenzene	0.050	ַן
121-14-2	2,4-Dinitrotoluene	0.050	ט
118-74-1	Hexachlorobenzene	0.050	ן
87-68-3	Hexachlorobutadiene	0.050	ַן
67-72-1	Hexachloroethane	0.050	ַן
98-95-3	Nitrobenzene	0.050	ן ד
87-86-5	Pentachlorophenol	0.25	ַ
110-86-1	Pyridine	0.10	ן ט
95-95-4	2,4,5-Trichlorophenol	0.050	יט
88-06-2	2,4,6-Trichlorophenol	0.050	ט
1319-77-3	Cresols (total)	0.050	ן



Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522013.D

Report Date: 22-May-2000 16:49

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file : \\QPITPA02\D\chem\71.i\s052200.b\S0522013.D

Lab Smp Id: DD6A7102 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 22-MAY-2000 16:17

Operator : 045183 Inst ID: 71.i

Smp Info : c0e130194-004 5/18/00 8270tclp Misc Info : dd6a7102,s052200.b,8270clp.m,1-tclp.sub

Comment

Method : \\QPITPA02\D\chem\71.i\s052200.b\8270clp.m Meth Date : 22-May-2000 12:37 bachas Quant Type: IS Quant Type: ISTD Cal Date : 19-MAY-2000 15:44 Cal File: S0519CC6.D

Als bottle: 15

Dil Factor: 1.00000 Integrator: HP RTE Target Version: 4.03

Processing Host: PITPC050

Compound Sublist: 1-tclp.sub

Concentration Formula: Amt \* DF \* Uf \* Vt/(Vo \* Vi)

Name	Value	Description
DF Uf Vt Vo Vi	0.001 1000.000 200.000	Dilution Factor ng unit correction factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected (uL)

							CONCENTR	ATIONS
		QUANT SIG					ON-COLUMN	FINAL
Co	pmpounds	MASS	RT	EXP RT	REL RT	RESPONSE	( ng)	(mg/L)
	. 医环苯基苯基基氏 20 12 12 12 12 12 12 12 12 12 12 12 12 12			*****		****	982222	000000
*	1 1,4-Dichlorobenzene-d4	152	5.029	5.029	(1.000)	155705	40 0000	(a)
*	2 Naphthalene-d8	136	6 563	6.562	(1 000)	593533	40 0000	(a)
*	3 Acenaphthene-d10	164	9 592	9 596	(1 000)	302566	40 0000	(a)
*	4 Phenanthrene-d10	188	12.893	12 898	(1.000)	486337	40.0000	(a)
*	5 Chrysene-d12	240	19.517	19.533	(1 000)	495522	40.0000	(a)
*	6 Perylene-d12	264	22.883	22 893	(1 000)	603824	40.0000	(a)
	10 Pyridine	79	Con	pound No	ot Detect	ed		
	28 1,4-Dichlorobenzene	146	Соп	pound No	ot Detect	eđ		
М	34 Cresols, total	100	Con	pound No	ot Detect	eđ		
	31 2-Methylphenol	108	Соп	pound No	ot Detect	ed.		
	35 4-Methylphenol	108	Con	pound No	ot Detect	ed		
	38 Hexachloroethane	117	Con	pound No	ot Detect	ed.		
	39 Nitrobenzene	77	Con	pound No	ot Detect	ed.		
	59 Hexachlorobutadiene	224	Соп	mpound No	ot Detect	eđ.		

			CONCENTRATIONS
	QUANT SIG		ON-COLUMN FINAL
Compounds	MASS	RT EXP RT REL RT RESPONSE	( ng) ( mg/L)
	#===	立中 埃尔拉斯斯德 医苯苯基苯酚 电电流放应性电路	규제들위료들도 모든경크놀라요
69 2,4,6-Trichlorophenol	196	Compound Not Detected	
70 2,4,5-Trichlorophenol	196	Compound Not Detected	
91 2,4-Dinitrotoluene	165	Compound Not Detected.	
113 Hexachlorobenzene	283	Compound Not Detected.	
117 Pentachlorophenol	265	Compound Not Detected.	
\$ 172 Nitrobenzene-d5	82	5 686 5 686 (0.867) 318009	52.9832 0 13246(a)
\$ 173 2-Fluorobiphenyl	172	8.304 8 309 (0 866) 498621	54 3277 0 13582(a)
\$ 174 Terphenyl-d14	244	16 964 16 969 (0 869) 747900	67.4938 0 16873(a)
\$ 175 Phenol- <b>d5</b>	99	4 714 4.719 (0 937) 529317	76 2060 0.19052(a)
\$ 176 2-Fluorophenol	112	3 705 3 709 (0 737) 370832	71 3466 0.17837(a)
\$ 177 2,4,6-Tribromophenol	330	11 333 11 338 (0.879) 106910	88.5462 0 22136(a)
\$ 178 2-Chlorophenol-d4	132	4 826 4 831 (0.960) 419058	86.1179 0 21529(a)
\$ 179 1,2-Dichlorobenzene-d4	152	5 238 5.237 (1 041) 158520	46 3672 0 11592(a)

#### QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).

# GC/MS SEMIVOLATILE CALIBRATION DATA

# 657 236

Lab Name:

Contract:

Lab Code: Case No.: SAS No.: SDG No.:

Instrument ID: 71 Calibration Date(s): 05/19/00

Min  $\overline{RRF}$  for SPCC(#) = 0.050

Max %RSD for CCC(\*) = 30.0%

LAB FILE ID: RRF1 RRF3 =S0519CC4.D RRF4	=S05190 =S05190		RRF:		19CC3.D 19CC6.D		
	T		1	<del></del>			l
COMPOUND	RRF1	RRF2	RRF3	RRF4	RRF5	RRF	ફ RSD
Phenol	* 2.092	2.099	2.050	1.918	1.891	2.010	4.9
bis(2-Chloroethyl)ether	1.638					1.625	2.0
2-Chlorophenol	1.405		1.420				2.6
1,3-Dichlorobenzene	1.571		1.567		1.473		3.8
1,4-Dichlorobenzene	± 1.612		1.589		1.486	1.564	4.0
1,2-Dichlorobenzene	1.511	1.511	1.466	1.376	1.343	1.441	5.4
2-Methylphenol	1.365	1.395	1.386	1.326		1.358	
2,2'-oxybis(1-Chloropropane	2.103	2.119	2.067	1.967		2.038	4.1
4-Methylphenol	1.442	1.466	1.452	1.380	1.356		3.4
Hexachloroethane	0.610	0.624	0.617	0.596	0.595		2.1
Nitrobenzene	0.432	0.448	0.444	0.430	0.431	0.437	1.9
Isophorone	0.715	0.741	0.734	0.718	0.720	0.726	1.5
2-Nitrophenol	* 0.176	0.190	0.196	0.193	0.194	0.190	4.2
2,4-Dimethylphenol	0.337	0.355	0.358	0.350	0.349	0.350	2.3
bis (2-Chloroethoxy) methane	0.466	0.473	0.473	0.456	0.455	0.465	1.9
N-Nitroso-di-n-propylamine	# 1.122	1.157	1.155	1.143	1.140	1.143	1.3
2,4-Dichlorophenol	* 0.267	0.283	0.280	0.271	0.270	0.274	2.5
1,2,4-Trichlorobenzene	0.296	0.299	0.294	0.281	0.278	0.290	3.3
Naphthalene	1.103	1.110	1.085	1.017	1.003	1.064	4.7
4-Chloroaniline	0.421	0.441	0.439	0.429	0.423	0.431	2.1
	* 0.151	0.154	0.152	0.142	0.140	0.148	4.1
	* 0.302	0.324	0.323		0.323	0.319	2.9
2-Methylnaphthalene	0.720	0.730	0.716	0.680	0.670	0.703	3.7
Hexachlorocyclopentadiene	# 0.273	0.308	0.316	0.311	0.313	0.304	5.8
	* 0.316	0.330	0.331	0.327	0.326	0.326	1.8
2,4,5-Trichlorophenol	0.330	0.352	0.356	0.353	0.348	0.348	3.0
2-Chloronaphthalene	1.116	1.118	1.091	1.027	1.007	1.072	4.8
2-Nitroaniline	0.353	0.390	0.398	0.401	0.405	0.389	5.4
Dimethylphthalate	1.284	1.324	1.320	1.287	1.282	1.299	1.6
Acenaphthylene	1.828	1.878	1.861	1.784	1.758	1.822	2.8
2,6-Dinitrotoluene	0.265	0.294	0.298	0.296	0.298	0.290	4.9
3-Nitroaniline	0.331	0.367	0.374	0.372	0.374	0.364	5.0
	1.162	1.180	1.158	1.114	1.103	1.143	2.9
	# 0.094	0.140	0.158	0.175	0.187	0.151	24.0
4-Nitrophenol	0.156	0.189	0.195	0.202	0.208	0.190	10.6
Dibenzofuran	1.643	1.656	1.630	1.554	1.525	1.602	3.6
2,4-Dinitrotoluene	0.373	0.410	0.414	0.421	0.420	0.408	4.9
					J. 120	0.100	ا د ٠٠

page 1 of 3

FORM VI SV-1

1/87 Rev.

# SEMIVOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:

Contract:

657 237

Lab Code: Case No.: SAS No.: SDG No.:

Instrument ID: 71 Calibration Date(s): 05/19/00

Min  $\overline{RRF}$  for SPCC(#) = 0.050

Max %RSD for CCC( $\star$ ) = 30.0%

page 2 of 3

FORM VI SV-2

1/87 Rev.

Lab Name·

Contract:

Lab Code: Case No.: SAS No.: SDG No.:

Instrument ID: 71 Calibration Date(s): 05/19/00

Min  $\overline{RRF}$  for SPCC(#) = 0.050

Max %RSD for CCC(\*) = 30.0%

LAB FILE ID: RRF1 RRF3 =S0519CC4.D RRF4	=S05190 =S05190		RRF:		19CC3.D 19CC6.D		
COMPOUND	RRF1	RRF2	RRF3	RRF4	RR <b>F</b> 5	RRF	RSD
1,2-Diphenylhydrazine Benzidine Methyl methanesulfonate 2-Naphthylamine 7,12-dimethylbenz[a]anthrac	0.888 0.409 0.862 1.262 0.490	0.539 0.848 1.153	0.537 0.853 1.167	0.865 0.554 0.851 1.102	0.869 0.559 0.848 1.068	0.885 0.520 0.852 1.150	1.9 12.0 0.7 6.4 5.7
Nitrobenzene-d5 2-Fluorobiphenyl Terphenyl-d14 Phenol-d5 2-Fluorophenol 2,4,6-Tribromophenol 2-Chlorophenol-d4 1,2-Dichlorobenzene-d4	0.389 1.224 0.870 1.855 1.346 0.095 1.290 0.902	0.416 1.257 0.900 1.810 1.346 0.098 1.267 0.920	0.412 1.239 0.901 1.806 1.342 0.102 1.258 0.899	1.178 0.898 1.744 1.326 0.101	0.403 1.169 0.903 1.708 1.315 0.101 1.210 0.829	1.213 0.894	2.5 3.2 1.5 3.3 1.1 3.1 2.6 4.6

page 3 of 3

FORM VI SV-3

1/87 Rev.

Data File: \\QPITPA02\D\chem\71.i\s051900.b/S0519CC3.D 657 239

Report Date: 05/22/2000

# INITIAL CALIBRATION REPORT

Instrument ID: 71.i Lab File ID: S0519CC3.D Analysis Type: NONE Injection Date: 19-MAY-2000 13:20

Lab Sample ID: sstd50
Method File: \\QPITPA02\D\chem\71.i\s051900.b\

!	I
COMPOUND	%RSD
医亚二苯亚因 中国各种基础: 医医中毒性 医中毒性 化苯基磺磺酸	** 40   200525
Benzo(b)illuorantheno	j 3 :
/,12-dimetnylbonz(a)anthracen	5
Benzo(k)fluoranthene	12
Benzo(a)pyrene	j 3.
Indeno(1,2,3 cd)pyrone	1 2.
Dibenz (a, h) anthracene	İı
Benzo (q, h, 1) perylene	.   2 -

The average of all  $\mbox{RRSD}^*\mbox{s}$  in the initial calibration is 4.5

657 240

#### INITIAL CALIBRATION REPORT

Instrument ID: 71.i Lab File ID: S0519CC3.D Analysis Type: NONE

Injection Date: 19-MAY-2000 13:20

Lab Sample ID: sstd50
Method File: \\QPITPA02\D\chem\71.i\s051900.b\8270clp

COMPOUND	%RS
===== ================================	====
N-N: trosodimethylamine	1 10
Pyridine	9
Mctnyl methanesulfonate	0
2 Fluorophenol	1
Phenol ds	] 3
Phenol	4
Anuline	4
b_s(2-Chloroethyl)ether	2
2-Chlorophenol-d4	2
2-Chlorophenol	2
1,3-Dachlorobenzene	{ 3
l,4-Dichloropenzeno	4
Benzyl Aldehol	1
1,2-Dichlorobenzene-q4	4
1,2 Dicilorobenzene	5
2-Mothylpsenol	2
2,2'-oxyols(1-Chloropropane)	4
4-Methylphenol	3
N-Nitroso-di n-promylamine	1
Hexaculornethane	2
Nitrobenzene-d5	2
Nitrobenzene	1
Isophorone	1
2-Nitrophenol	4
∠,4-Dimethy phonol	2
bis(2-Chloroethoxy) methane	1
Bonzoid Acid	23
2,4-Dichlorophenol	2
1,2,4-Tricalorobenzene	3
Naphthalene	4
4-Chloroaniline	. 2
Hexachlorobutadione	4
4-Caloro-3-Methylphenol	2
2 Methylnaphthalene	3
1-Methylnaphthalene	3
Fexachlorocyclopertadiene .	5
2,4,6 frichlorophenol	1
2,4,5-Trichloropnenol	3
2-Fluorobiphenyl	i 3

2034

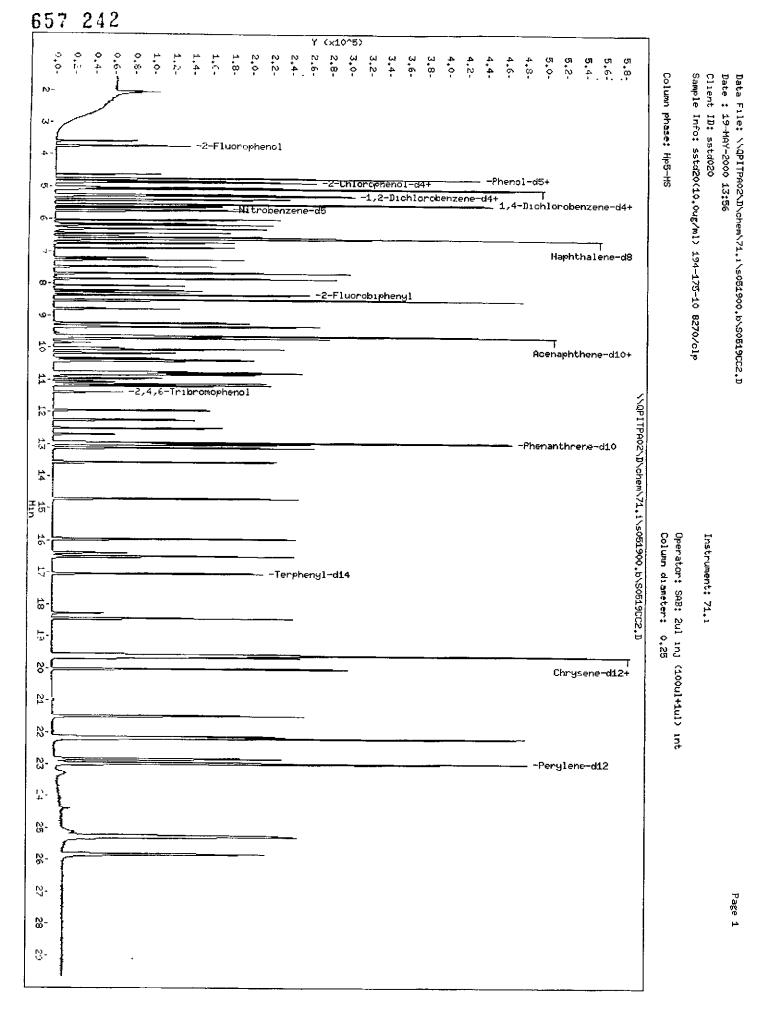
# INITIAL CALIBRATION REPORT

Instrument ID: 71.i Lab File ID: S0519CC3.D Analysis Type: NONE

Injection Date: 19-MAY-2000 13:20

Lab Sample ID: sstd50
Method File: \\QPITPA02\D\chem\71.i\s051900.b\

COMPOUND	i	*RSD
**** *********************************	* v = = =   :	
2-Chloromaphthalene	1	4 6
/-Nitroamiline	1	5 4
Dimethylphthalate	1	1 6
Acenaphthylene	1	2 8
2,6-Jinitrotoluene	1	4 9
3-Nitroamiline	}	5.0
Acenaphthene	1	2 9
2,4-Dimitropherol	1	24 (
4-Nitrophenol	1	10 6
Dibenzofuran	1	3 6
2,4 Dinitrotoluene	ĺ	4 5
2,3,5,6 Tetrachlorophenol	- 1	3.9
'2-Naphthylamine	i	6.4
7,3,4,6-Tetraculorophenol		1 7
Diethylph.nalace	ĺ	2 0
Fluorenc	- 1	3.7
4-Ciloropheryi-phenylether	i	3.3
4-Nitroamiline	- 1	5 3
4,6-D-nitro-2-methy2phenol	i	14.2
N Nitrosodiphenylamine (1)	1	3.2
1,2-Diphenylhydrazine	- 1	19
2,4,6-Tribromophenol	Ì	3 1
4-Bromophenyl-phenylether	- 1	2 9
Hexacl.lorobenzero		3 0
Pentachiorophenol	- 1	16 1
Phenanthreme	- 1	4 5
Anthracene	ĺ	4 1
Carbazole	1	2 6
Di-n-Butylphthalate	ļ	3 3
Fluorant hene	i	3 4
Benzidine	i	12 0
Pyrene	i	0 7
Terphonyd14	1	1 5
Butylbenzyiphthalare	ĺ	5 6
Benzo (a) Anthracene	}	2 0
3,3'-Dich_orobenzidine	į	6 7
Chrysenc	1	1.1
bis(2-ethylhexyl)Phthalate	i	5 3
Di-n-octylphthalate	ì	4 1



Data File: \\QPITPA02\D\chem\71.i\s051900.b\S0519CC2.D Report Date: 19-May-2000 14:32

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file: \\QPITPA02\D\chem\71.i\s051900.b\S0519CC2.D Lab Smp Id: sstd20 Client Smp ID: Client Smp ID: sstd020 Inj Date : 19-MAY-2000 13:56

Operator : SAB: 2ul inj (100ul+1ul) int Inst ID: 71.i Smp Info : sstd20(10.0ug/ml) 194-175-10 8270/clp Misc Info : sstd20,s051900.b,8270clp.m,1-82701.sub,1,1

Comment

: \\QPITPA02\D\chem\71.i\s051900.b\8270clp.m Method Meth Date : 19-May-2000 14:28 bachas Quant Type: ISTD Cal Date : 19-MAY-2000 13:56 Cal File: S0519CC2.D

Als bottle: 3 Calibration Sample, Level: 1 Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: 1-82701.sub

Target Version: 4.03 Processing Host: PITPC050

		•				AMOUN	vis
C-1	mpourids	QUANT SIG				CAL-AMT	ON-COL
	- <b>-</b>	MASS	RT	EXP RT REL RI	RESPONSE	( ng)	( ng)
	**************************************	그 한 다 프	==		======		4==== <u></u>
w	1 1,4-Dichlorobonzene-d4	152	5 072	5 072 (1 000)	95582	40 0000	
*	2 Naphthalene-d8	136	6 616	6.616 (1 000)	387208	40 0000	
*	3 Acenaphthene-dl0	164	9 661	9 661 (1.000)	222841	40.0000	
*	4 Phenanthrene dl0	188	12 973	12 973 (1 000)	408296	40 0000	
*	5 Chryseno dl2	240	19 608	19 608 (1 000)	403670	40 0000	
•	6 Porylene-d12	264	22.979	22 979 (1 000)	459326	40 0000	
	13 N-Nitrosodimetnylamine	74	2 048	2 048 (0 404)	24090	20 0000	19 435
	10 Pyrinine	79	2 070	2 070 (0 408)	45943	20 0000	20 004
	<pre>19 Vethyl methanesulfonate</pre>	80	3 582	3 582 (0 706)	41196	20 0000	20 164
	22 Amiline	93	4 768	4 768 (0 940)	124962	20.0000	20 357
	23 Phenol	94	4.762	4.762 (0 939)	99994	20 0000	19 969
	24 bis(2 Chloroethy1)ether	93	4.837	4 837 (0 954)	78305	20 0000	19.830
	25 ? Chiorophenol	128	4.880	4 880 (0 962)	67155	20 0000	19 825
	2/ 1,3 Dichlorobenzene	146	5.029	5 029 (0 992)	75082	20.0000	19 761
	28 1,4-Dichlorobenzene	146	5.088	5.088 (1 003)	77046	20 0000	19 913
	29 1,2-Dichlorobenzene	146	5 296	5 296 (1 044)	72217	20 0000	19 999
	30 Berzyl Alcohol	108	5 254	5 254 (1 036)	47626	20 0000	19 767
	31 2 Methylphonol	108	5 403	5 403 (1 065)	65219	20.0000	19 781
	32 2,2'-oxybis(1-Chloropropane)	45	5 430	5 430 (1 071)	100506	20.0000	19.922
	33 N-Natroso-di-n-propylamine	70	5 585	5.585 (1 101)	53598	20 0000	19.688
	35 4-McLhylphenol	108	5 558	5 558 (1 096)	68891	20.0000	19.888
	38 Hexachloroethane	117	5 644	5 644 (1 113)	29139	20.0000	19 833
	39 Nitropenzere	77	5 750	5.750 (0.869)	83564	20.0000	
	44 Isophorone	82	6 028	6.028 (0 911)	138454	20 0000	19 638
	45 2-Nitrophenol	139	6 146	6 146 (0 929)	34128	20 0000	19.643
	46 2,4 Dimethylpheno	107	6 194	6 194 (0 936)	65287	20.0000	19 242 19 486

			-			
					NUOMA	ris
Carry	QUANT SIG				CAL-AMT	ON-COL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	(ng)
47 bis'2-Chloroethoxyinethane	93	=± 6 317	5 313 (A ACC)		======	======
51 2,4-Dicalorophenol	162	6 445	6 317 (0 955)	90172	20 0000	19 849
52 Benzoic Acid	122	6 295	6 445 (0 974)	51708	20 2000	19 432
53 1,2,4-Trichlorobenzenc	180	6 557	6.295 (0 952)	21194	20 0000	15 930(H)
54 Naphthalene	128	6 643	6.557 (0 991) 6 643 (1 004}	57333	20.0000	19 904
55 4 Caloroaniline	127	6.750	6 750 (1 004)	213576	20 0000	19 936
o9 Nexachlorobutaniene	225	6 899	6 899 (1 043)	81495 29278	20 0000	19 534
62 4-Chloro 3-Methylphenol	107	7 476	7 476 (1 130)	58580	20 0000	19 840
65 2-Methylnaphthalenc	142	7 684	7 684 (1.161)	139480	20 0000	19 320
66 1-Methylnaphthalene	142	7.866	7 866 (1 189)	133562	20 0000 20 0000	19 873
67 Hexachlorocyclopentadiene	237	8 074	8 074 (0 836)	30461	20.0000	20 229
69 2,4,6-Trichlorophenol	196	8 224	8 224 (0 851)	35177	20.0000	18.821
/0 2,4,5-irichlorophenol	196	8 288	8 288 (0.858)	36750	20 0000	19.572 19 360
73 2-Caloronaphthalone	162	8 534	8 534 (0 883)	124387	20.0000	19 980
77 2 N_:roaniline	65	8 796	8 796 (0 910)	39306	20.0000	19 980
80 Dinethylphthalate	163	9 239	9 239 (0 956)	143039	20 0000	19 690
82 2,6-Dinitrotoluene	165	9 362	9.362 (0.969)	29521	20,0000	19 890
83 Accnaphthylene	152	9 340	9 340 (0 967)	203685	20.0000	18 909
85 3 Nutroaniline	138	9 629	9 629 (0 997)	36916	20 0000	18 987
86 Acenaphthene	153	9 725	9 725 (1.007)	129485	20 0000	19 650
87 2,4-Dinitiopmenol	184	9.848	9.848 (1 019)	10524	20 0000	16 078
89 4-Nitrophenoi	109	10 014	10 014 (1 036)	17446	20 0000	18 123
90 Dibenzofuran	168	10 062	10 062 (1 041)	183068	20 0000	19 922
93 2,4-Dinitrotoluene	165	10.179	10 179 (1 054)	41582	20 0000	19 059
95 2,3,5,6-Tetraenlorophenol	232	10.339	10 339 (1 070)	27759	20 0000	19.379
92 2,3,4,6 letrachlorophenol	232	10 436	10 436 (1 080)	29522	20 0000	19.907
96 2-Naphthylamine	143	10 404	10 404 (1 077)	140631	20 0000	20 904
9/ Diethylphthalate	149	10 772	10 772 (1 115)	144938	20.0000	19.538
98 Flucrenc	166	10 810	10 810 (1 119)	149566	20 0000	19.860
99 4-Chlorophenyl-phenylether	204	10 852	10 852 (1 123)	66595	20 0000	19.888
100 4-Natroaniline	138	10 959	10 959 (1 134)	38927	20 0000	18 883
102 4,6-Dimitro 2-methylphenol	198	11 071	11.071 (0 853)	18357	20 0000	17 503
103 N-Nitrosodiphenylamine (1)	169	11 135	11.135 (0 858)	112441	20 0000	20 429
104 1,2-Diphonyihydrazine	77	11 199	11 199 (0 863)	181300	20 0000	19 897
112 4-Bromophenyl-phenylother	248	11 958	11 958 (0.922)	36896	20 0000	19 945
113 Hexachlorobenzene	284	12.263	12 263 (0.945)	39662	20 0000	19 996
117 Pencas hlorophenol	266	12 706	12 706 (0 979)	16299	20 0000	17 037
122 Phemanthren=	178	13 021	13 021 (1 004)	219499	20 0000	20 102
123 Anthracene	178	13 128	13 128 (1.012)	222028	20 0000	19 919
126 Carbazole	167	13.566	13 566 (1.046)	210285	20.0000	19.737
130 Di-n Butylphthalate	149	14.709	14 709 (1 134)	245902	20 0000	19.245
135 Fluoranthene	202	15 965	15 965 (1 231)	224015	20 0000	19 604
136 Benzidine	184	16 392	16 392 (0 836)	82643	20.0000	17 263
137 Pyrene	202	16 510	16 510 (0 842)	238065	20.0000	19 924
144 Butylbenzylphthalate	149	18 454	18 454 (0 941)	116471	20 0000	18 988
149 3,3'-Dichlorobenzidine	252	19 624	19.624 (1 001)	75267	20 0000	18 770
150 Benzo(a) Anthracene	228	19 571	19 571 (0.998)	219875	20 0000	19 517

Data File: \\QPITPA02\D\chem\71.i\s051900.b\S0519CC2.D Report Date: 19-May-2000 14:32

					AMOUN	TS
Compounds	QUANT SIG				CAL-AMT	ON-COL
•	MASS	RT	EXP RT REL RT	RESPONSE	{ ng}	( ng)
·	****	==		4= <b>3</b> ====	======	
151 Chrysere	228	19 672	19 672 (1 003)	217527	20 0000	19 768
153 pis(2 etnylhexyl)Phthalate	149	20 030	20 030 (1 022)	164169	20 0000	
155 Di-m-octylphthalate	149	21 499	21 499 (0 936)	281772	30 0000	18 958
157 Berzo(b) f uoranLhene	252	22 151	22.151 (0 964)	238605	20 0000	19 046
158 Benzo(k)fluoranthène	252	22 204	22 204 (0 966)	287901		19 560
<pre>15) 7.12-dimethylbenz(alanthracen</pre>	256	22 204	72 204 (0 966)	112535	20 0000	20 812
16/ Benzo(a) pyrene	252	22 845	22 845 (0 994)		20 0000	20 460
169 indeno(1,2,3-cd)pyrene	276	25 255		236893	20 0000	19 649
170 Dibenz(a,h)anthracene	278	25 303	25 255 (1 099)	325005	30 0000	19 810
171 Benzo(g,h,i)perylene	276	=	25 303 (1 101)	277991	20 0000	19 848
\$ 172 Nitrobenzene-d5		25 842	25 842 (1 125)	283066	20 0000	19 604
\$ 173 2-Pluorobipnenyl	82	5.729	5 729 (0.866)	75344	20 0000	19 345
_ ·	172	8 363	8.363 (0 866)	136348	20 0000	19 732
\$ 174 Terphenyl-d14	244	17 038	17 038 (0 869)	175629	20 0000	19 656
\$ 170 Phenol-dS	99	4 746	4 746 (0 936)	88640	20 0000	20 242
5 170 2-Fluorophenol	112	3 742	3.742 (0 738)	64347	20 0000	20 001
5 177 2,4,6-Tribromopherol	330	11 402	11 402 (0 879)	19339	20 0000	19.695
\$ 178 2 Chiorophonol-d4	132	4 869	4 869 (0 960)	61637	20 0000	
\$ 179 1,2-Dichlerobenzene-d4	152	5 <b>2</b> 80	5 280 (1 041)	43125	20 0000	20 178 19 809

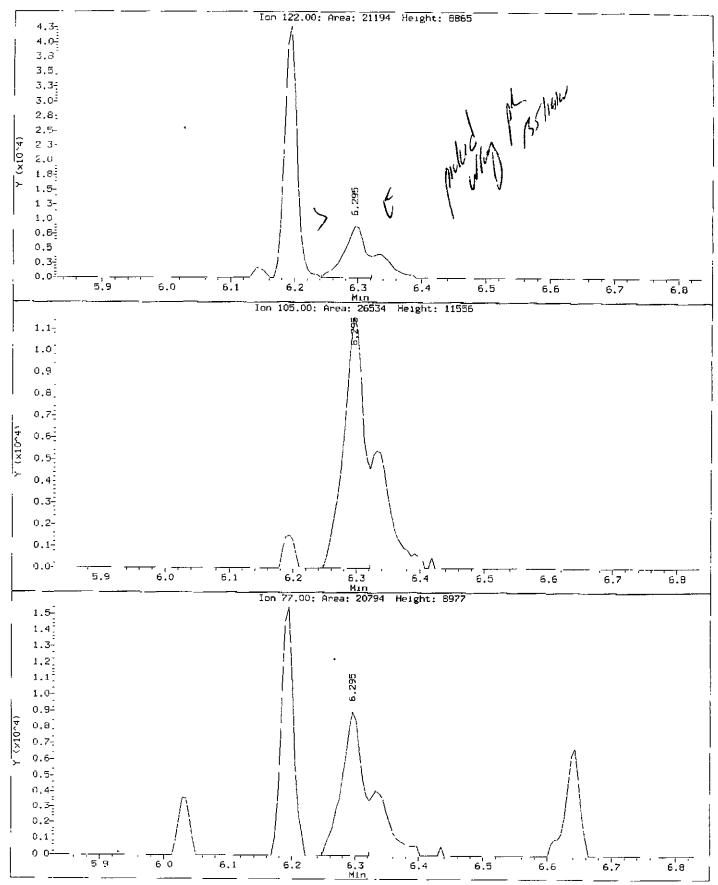
### QC Flag Legend

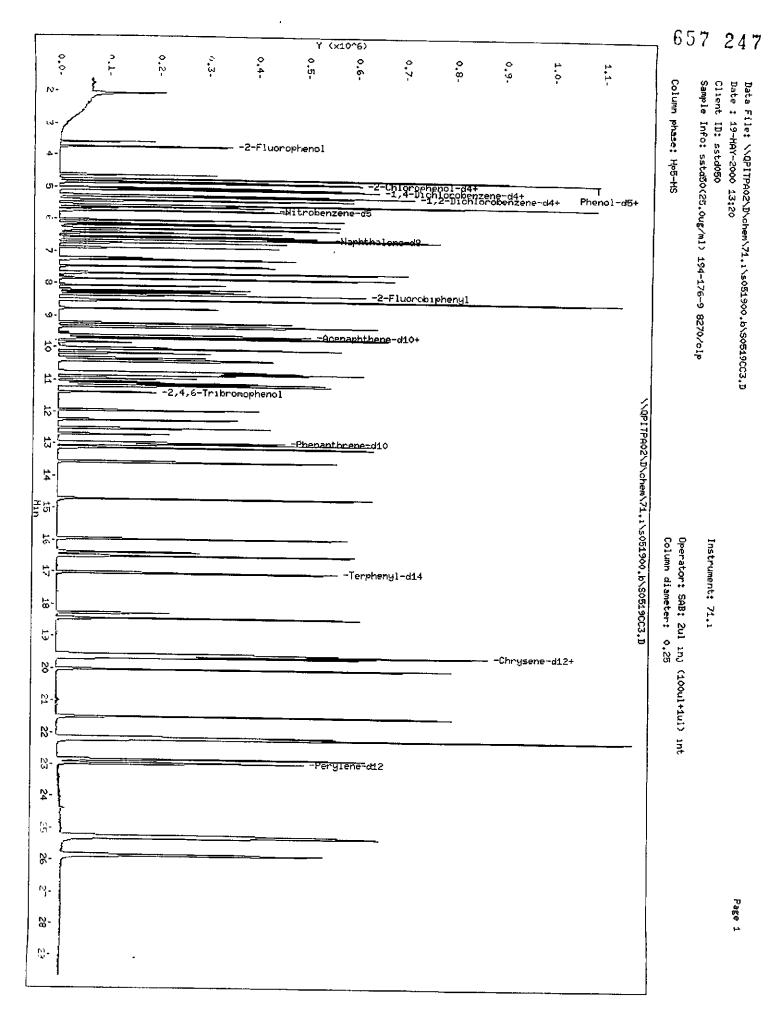
H - Operator selected an alternate compound hit.

246

Data File: \\QPITPAO2\D\chem\71.1\s051900.b\S0519CC2.D Injection Date: 19-MAY-2000 13:56 Instrument: 71.1 Client Sample ID: sstd020

Compound: Benzoic Acid CAS Number: 65-85-0





Data File: \\QPITPA02\D\chem\71.i\s051900.b\S0519CC3.D

Report Date: 19-May-2000 14:10

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file : \\QPITPA02\D\chem\71.i\s051900.b\S0519CC3.D

Lab Smp Id: sstd50 Ing Date : 19-MAY-2000 13:20 Client Smp ID: sstd050

Operator : SAB: 2ul inj (100ul+1ul) int Inst ID: 71.1

Smp Info : sstd50(25.0ug/ml) 194-176-9 8270/clp

Misc Info : sstd50,s051900.b,8270clp.m,1-82701.sub,1,2

Comment

Method : \\QPITPA02\D\chem\71.i\s051900.b\8270clp.m Meth Date : 19-May-2000 13:56 bachas Quant Type: ISTD Cal Date : 19-MAY-2000 13:20

Cal File: S0519CC3.D Als bottle: 4 Calibration Sample, Level: 2

Dil Factor: 1.00000 Integrator: HP RTE

Compound Sublist: 1-82701.sub

Target Version: 4.03 Processing Host: PITPC050

13/11/14

Page 1

					AMOUN	TS
	QUANT SIG				CAL-AMT	ON-COL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
	====	==	***=== <b>====</b>		==== t==	==== -==
1 1,4-Dichlorobenzene d4	152	5 064	5 064 (1 000)	90711	40.0000	
√ ∠ Naphthalene-d9	136	6 508	6 608 (1 000)	365257	40 0000	
3 Acenaphthene-dic	164	9 648	9 648 (1 000)	211027	40 0000	
4 Phonanthrene-dl0	188	12.960	12 960 (1 000)	395064	40 0000	
5 Chrysone-d12	240	19 600	19 600 (1 000)	396681	40 0000	
5 Perylene d12	264	22 965	22 965 (1 000)	476712	40 0000	
13 N-N trosodimethylamine	74	2 046	2 046 (0.404)	60479	50 0000	38 914
10 Pyrid-ne	79	2 051	2.051 (0 405)	108964	50 0000	40 778 (M)
19 Methyl methanesulfonate	80	3 579	3.579 (0 707)	9€149	50 0000	90.420
22 Aniline	93	4 765	4 765 (0 941)	286086	50.0000	97 333
∠3 Phenol	94	4 760	4 760 (0.940)	237977	50.0000	62 195
24 bis(2-Chloroethy1)ether	93	4 829	4 829 (0 954)	188978	50 0000	61 781
25 2-Chlorophenol	128	4 877	4 877 (0 963)	162137	50 0000	58.599
27 1,3-Dichlorobenzene	146	5 027	5 027 (0 993)	182454	50 0000	59 318
28 l,4-Dichiorobenzene	146	5.085	5 085 (1 004)	184390	50 0000	59 389
29 1,2-D-chlorobenzene	146	5 288	5.288 (1 044)	171362	50 0000	60.083
30 Pcnzyl Alcohoi	108	5 246	5 246 (1 036)	115661	50 0000	58 782 (M)
31 2-Methylphenol	108	5 395	5 395 (1 065)	158169	50 0000	58 616
32 2,8'-oxybis(1-Chloropropane)	45	5.422	5 422 (1 071)	240324	50 0000	63.171
33 N-Nitrosc-di-n propylamine	70	5.577	5.577 (1 101)	131200	50 0000	60.868
3b 4-Methylphenol	108	5 556	5.556 (1 097)	166203	50 0000	60 879
38 Hexachloroethane	117	5 636	5 636 (1.113)	70729	50 0000	59 768
39 Nitropenzene	77	5 748	5 748 (0 870)	204339	50 0000	63 049
44 isophorone	82	6 026	6 026 (0 912)	338383	50 0000	63 958
45 2-Nitrophenol	139	6 138	6 138 (0 929)	86819	50 0000	57 466
46 2,4-Dimethylphonol	107	6 186	6 186 (0 936)	162083	50,0000	58 358

					AMOUN	ITS
Compounds	QUANT SIG				CAL-AMT	ON-COL
TIBEBBEN TERRET MERSE	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
4/ bis(2 Chloroethoxy)methane	32.63	= 10	=======================================	3 # 2 2 2 E F F	====>==	===6===
51 2,4-Dichlorophenol	93	6 309	6 309 (0 955)	215890	50 0000	62 215
>2 Benzouc Acid	162	6 437	6 437 (0 974)	129066	50 0000	58 934
53 1,2,4 Trichlorobenzene	122	6 330	6 330 (0 958)	75522	50 0000	51 964(11)
54 Naphtbalene	180 128	6 549	6 549 (0 991)	136509	50 0000	59 056
55 4-Chloroaniline	128	6 635 6 736	6 635 (1.004)	506904	50 0000	61 464
53 Hexachlorobutadiene	225	6 891	6 736 (1 019)	201350	50 0000	66 461
62 4-Chloro-3-Methylphenol	107	7.473	6 891 (1 C43)	70160	50 0000	59 670
65 2-Methylnaphthalenc	142	7 676	7 473 (1 131)	147871	50 0000	59.252
66 1-Methyinaphthalene	142	7 858	7 676 (1 162)	333144	50 0000	61 897
67 Hexachlorocyclopentadiene	237	8 066	7 858 (1 189)	307834	50 0000	59 252
69 2,4,6-Trichlorophenol	196	8 210	8 066 (0 836) 8 210 (0 851)	81147	50 0000	64.899
/C 2,4,5-frichlorophenal	196	8 280	8 280 (0 858)	86921	50 0000	59 217
73 2-Chloronaphthalene	162	8 520	8 520 (O 893)	92756	50 0000	56 466
77 2 Nitroaniline	65	8 787	8 787 (0 911)	295057	50 0000	62 580
80 Dimethylphthalate	163	9 231	9 231 (0 957)	102811	50 0000	62.476
82 2,6-Dinitrotoluene	165	9 359	9.359 (0 970)	349292	50 0000	61 676
83 Acenaphthylene	152	9 332	9 332 (0 967)	77490 495539	50 0000	59 393
85 3-Nilroaniline	138	9 621	9 621 (0 997)		50 0000	61 478
86 Accnaphthere	153	9 717	9 717 (1 007)	96720 311190	50 0000	58 882
87 2,4 Dimitrophenol	184	9 840	9 840 (1 020)	37072	50 0000	61 143
89 4-Nitrophenol	109	10 011	10 011 (1 038)	49859	50 0000	51 380
90 Diberzofulan	168	10.048	10 048 (1 042)	436788	50 0000	60 101
91 2,4 Dinitrotoluene	165	10 171	10 171 (1 054)	108161	50 0000 50 0000	61 637
95 2,3,5,6-Intrachlorophenol	232	10 331	10 331 (1 071)	69954		58 791
32 2.3,4,6-Tetracolorophenol	232	10 427	10 427 (1 081)	70544	50 0000	56 640
96 2-Naphthylamine	143	10.395	10.395 (1 078)	304128	50 0000 50 0000	57 647
97 Diethylphthalate	149	10 764	10 764 (1 116)	359343	50 0000	61 776
98 Fluorene	166	10 801	10 801 (1 120)	359063	50 0000	61 954
39 4-Chlorophenyl-phonylether	204	10 839	10 839 (1 123)	159438	50 0000	61 234 61 175
100 4 Nitroaniline	138	10 962	10 962 (1,136)	103061	50 0000	59 607
102 4,6-Dinitro-2-methylphenol	198	11 063	11 063 (0 854)	57074	50 0000	54 750
103 N-Nitrosodiphenylamine (1)	169	11 127	11 127 (0 859)	260564	50 0000	65 236
104 1,2-Diphenylhydrazine	77	11 186	11 186 (0 863)	443090	50 0000	68 648
112 4-Bromophenyl-phenylether	248	11 945	11.945 (0 922)	89743	50 0000	60 742
113 Wexachlorobenzene	284	12 249	12 249 (0 945)	95977	50 0000	60 065
117 Pentachlorophenoi	266	12 693	12 693 (0 979)	53140	50 0000	49 791
122 Pachanthrenc	178	13 013	13 013 (1 004)	525579	50 0000	62 316
123 Anthracene	178	13 120	13.120 (1 012)	541431	50 0000	64 977
126 Carbazolo	167	13 558	13 558 (1 046)	522242	50 0000	62 257
130 Di-n-Butylphthalacc	149	14 701	14 701 (1 134)	611526	50 0000	62 782
135 Fluoranthene	202	15 957	15 957 (1 231)	563796	50.0000	61 445
136 Benzidine	184	16.379	16 379 (0.836)	267421	50.0000	66 685
137 Pyrene	202	16 496	16 496 (0 842)	589294	50 0000	59 349
144 Butylbonzylphthalate	149	18 441	18 441 (0 941)	316631	50 0000	59 024
149 3,3' Dichlorobenzidine	252	19 611	19 611 (1 001)	209147	50.0000	58 329
150 Berizo(a) Ant hracone	228	19 563	19 563 (0.998)	566912	50.0000	57 901

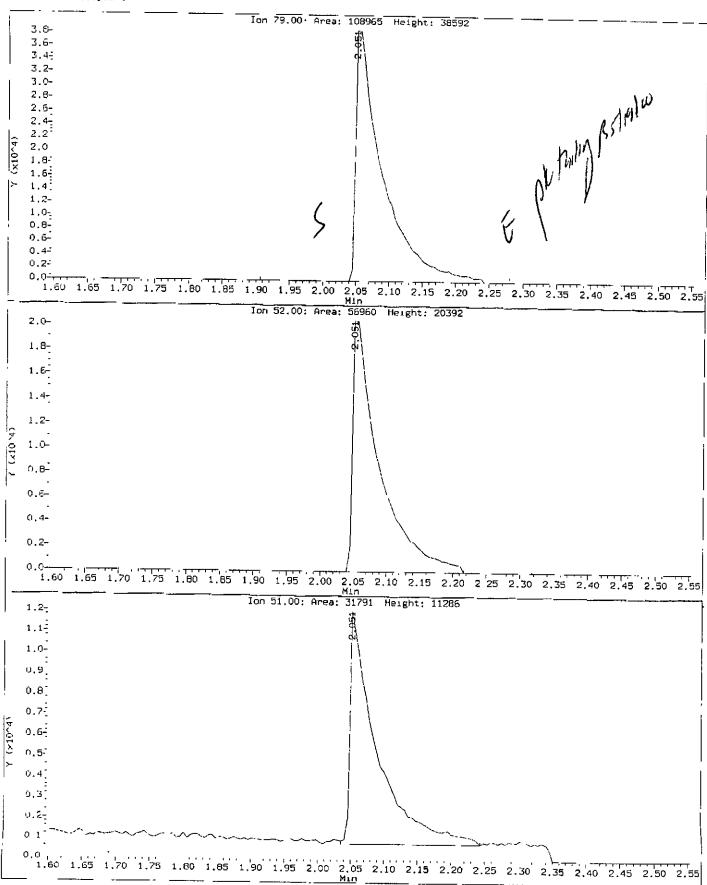
					AMOUN	TS
	QUANT SIG				CAL-AMT	ON-COL
Compounds	· MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
	====	70 00		===4====	=== -===	==+====
151 Chrysene	228	19 664	19 664 (1 003)	546969	50 0000	57 235
153 bis(2-ethylhexyl)Phthalat	e 149	20 017	20 017 (1 021)	447632	50 0000	57 473
155 Pr-n-octylpnihalate	149	21,491	21.491 (0 936)	804314	50 0000	60 277
157 Benzo(b)fluoranthene	252	22 148	22.148 (0 964)	646950	50 0000	59 036
158 Benzo(k)fluorantmenc	252	22 207	22 207 (0 967)	688682	50 0000	67 644
159 7,12-dimethylbenz(a)anthr	acen 256	22 202	22 202 (0 967)	278848	50 0000	65 952
167 Benzo(a)pyrene	25?	22 843	22 843 (0 995)	636581	50 0000	63 178
169 indeno(1,2,3-cd)pyrene	276	25 247	25 247 (1 099)	859402	50 0000	60 321
170 Dibens(a,h)anthracenc	278	25 295	25 295 (1 101)	732361	50 0000	61 072
171 Benzo(g,n,i)perylene	276	25.845	25 845 (1 125)	764160	50 0000	59 681
\$ 172 Nitiopenzene-d5	82	5.726	5 726 (0 867)	189711	50 0000	62 779
\$ 173 2-Fluorob_phenyl	172	8 355	8 355 (0 866)	331548	50 0000	60 980
\$ 174 Torphenyl-dl4	244	17 030	17 030 (0.869)	446559	50 0000	61 330
\$ 1/p Phenol-d5	99	4 744	4.744 (0 937)	205269	50 0000	62 855
\$ 176 z-Fluorophenol	112	3 734	3.734 (0 737)	152654	50.0000	57 140
\$ 177 2,4,6-Tribromophenol	330	11 389	11 389 (0 879)	48231	50 0000	<b>57</b> 752
\$ 178 2 Chlorophenol-d4	132	4 861	4 861 (0 950)	143652	50 0000	54 259
\$ 1/9 1,2 Dichlorobenzene-d4	152	5 272	5.272 (1 041)	104290	50 0000	59 318

#### QC Flag Legend

M - Compound response manually integrated.
H - Operator selected an alternate compound hit.

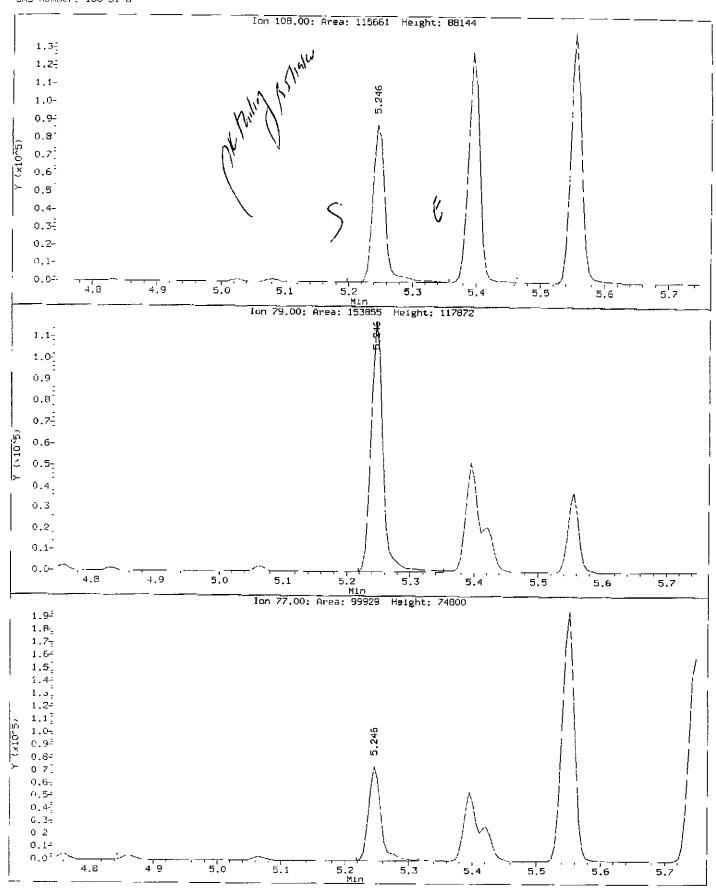
Data File: .\QPITPAO2\D\chem\71.1\s051900.b\S0519CC3.D Injection Date: 19-MAY-2000 13:20 Instrument: 71.1 Client Sample ID: sstd050

Compound Pyridine CAS Number: 110–86–1



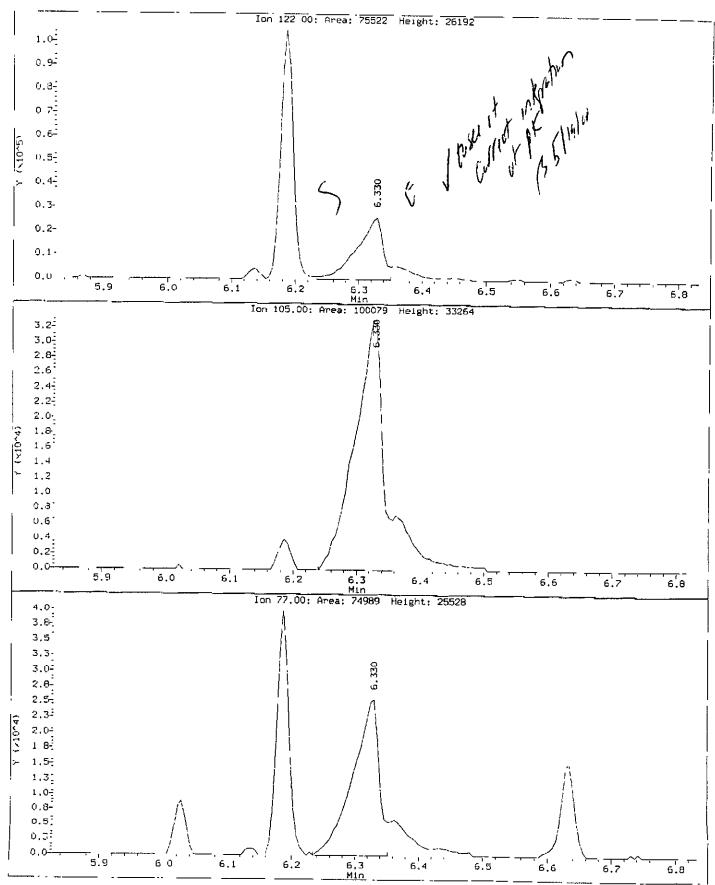
Data File: \\QPTIPA02\D\chem\71.1\s051900 b\S0519CC3.D Injectior Date: 19-MAr-2000 13:20 Instrument: 71.1 Client Sample ID: sstdU50

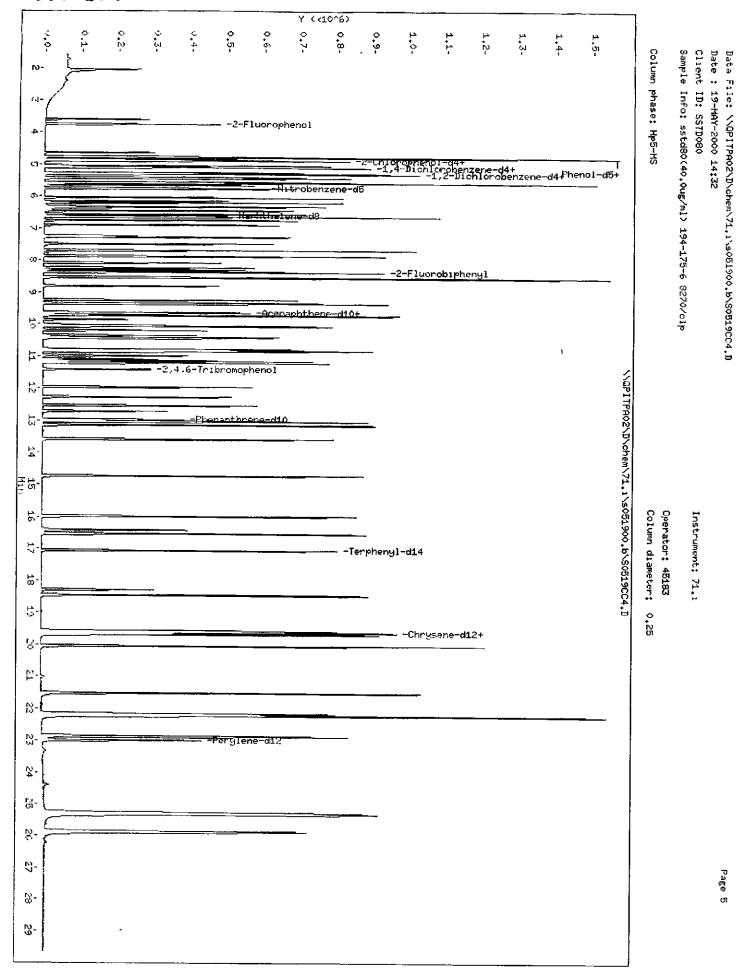
Compound: Benzyl Alcohol CAS Number: 100-51-6



Data File: \\QPfTPAO2\D\chem\71.i\s051900.b\S0519CC3.D lnjection Date. 19-MAY-2000 13:20 Instrument. 71.i Client Sample ID; sstd050

Compound: Benzolc Acid CAS Number: 65-85-0





Data File: \\QPITPA02\D\chem\71.i\s051900.b\S0519CC4.D Report Date: 22-May-2000 08:00

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file: \\QPITPA02\D\chem\71.i\s051900.b\S0519CC4.D Lab Smp Id: sstd80 Client Smp ID:

Client Smp ID: SSTD080

Inj Date : 19-MAY-2000 14:32

Operator : 45183 Inst ID: 71.i

Smp Info : sstd80(40.0ug/ml) 194-175-6 8270/clp

Misc Info : sstd80,s051900.b,8270clp.m,1-82701.sub,1,3

Comment

: \\QPITPA02\D\chem\71.i\s051900.b\8270clp.m Method Meth Date : 22-May-2000 07:54 bachas Quant Type: ISTD Cal Date : 19-MAY-2000 15:44

Cal File: S0519CC6.D Als bottle: 5

Calibration Sample, Level: 3 Dil Factor: 1.00000 Integrator: HP RTE

Compound Sublist: 1-82701.sub

Target Version: 4.03 Processing Host: PITPC050

						TS KINU
	Office and				AMOUN	TS /
Сотроина	QUANT SIG				CAL-AMT	ON-COL
	MASS	RT	EXP RT REL RT	RESPONSE	( ਹਰ੍ਹ)	( ng)
* 1 1,4-Dichioropenzene-d4	150	##	**************************************			<b>海港市</b> 连接管
* 2 Naphthalone-d8	152	5.079	5 079 (1 000)	83213	40 0000	
* 3 Acenauli hene di0	136	6.623	6 623 (1 000)	335457	40 0000	
* 4 Phenanthrene-d10	164	9 674	9 674 (1 000)	192142	40 0000	
	188	12 986	12 986 (1 000)	355838	40 0000	
* 5 Chrysene-dl2	240	19 631	19 631 (1 000)	348071	40 0000	
* 6 Perylone-dl2	264	22 992	22.992 (1 000)	426270	40 0000	
13 N Nitrosodimethylamine	74	2 051	2.051 (0 404)	85691	80 1000	80 200
10 Pyridine	79	2 056	2 056 (0 405)	160040	80 0000	80 000 (M)
19 Methyl mothane-ulforate	80	3 589	3 589 (0 707)	141969	80 0000	80 000
22 An_line	93	4 775	4 775 (0 940)	418139	80 0000	80 000
23 Phenol	94	4 775	4 775 (0 940)	341209	80 0000	80 000
24 big(2 Chloroethyl)ether	93	4 844	4 844 (0 954)	272240	80 0000	80 000
25 2-Chloropheno1	128	4 893	4 893 (0 963)	236318	80 0000	80 000
27 1,3-Dichloropenzene	146	5 037	5 037 (0 992)	260823	80 0000	80 000
28 1,4 Dichlorobanzene	146	5 096	5 096 (L 003)	264477	80,0000	80 000
29 1,2-Dichlorchenzene	146	5 304	5 304 (1 044)	243921	80 0000	80 000
30 Benzyl Alcohol	108	5 261	5 261 (1.036)	172022	80.0000	80.000
31 2-Mothylphenol	208	5 411	5 411 (1.065)	230744	80 0000	80.000
32 2,2 -oxybis(1-Chloropropane)	45	5 437	5.437 (1 070)	344031	80 0000	80 000
33 N-Nitrosc-di-n propylamine	70	5 598	5 598 (1 102)	192276	80 0000	80 000
35 4-Methylphenoi	108	5 571	5 571 (1.097)	241650	80 0000	80 000
38 Hexachloroethane	117	5.651	5 651 (1.113)	102708	80 0000	80.000
39 Nitrobenzere	77	5 763	5 763 (0 870)	297844	80 0000	80 000
44 Imophorone	82	6 041	6 041 (0 912)	492451	80 0000	80 000
45 2-Nitrophenol	139	6 153	6 153 (0 929)	131199	80 0000	80 000
46 2,4-Dimethy phonol	107	6 201	6.201 (0 936)	240195	80 0000	80 000

	QUANT SIG				AMOUN	
Compounds	MASS	RT	EXP RT REL RT	DEGEOVER	CAL-AMT	ON-COL
7	====	==	DAT KI KEL KI	RESPONSE	( ng)	( ng)
47 bis(2-Chloroethoxy)methane	93	6 330	6 330 (0 956)	317418	80 0000	80 000
51 2,4 Dichlorophenol	162	6 452	6 452 (0 974)	187968	80 0000	80 000
52 Benzold Acid	122	6 36/	6 367 (0 961)	126174	0000 08	80 000 (M
53 1.2,4-Trichloropenzone	180	6 570	6 570 (0 992)	197612	80 0000	80 000 (M
54 Naphthalene	128	6 650	6 650 (1 004)	728025	80 0000	80 000
55 4-Chloroariline	127	6 757	6 757 (1 020)	294298	80 0000	80 000
59 HexaculoroLutadieno	222	6 907	6.907 (1 043)	301668	80.0000	80 000
62 4-Chloro-3-Methylphenol	107	7 494	7 494 (1 131)	216692	80 0000	80 000
65 2-Methylmaphthalone	142	7.697	7 697 (1 162)	480114	80 0000	80 000
66 1 Methylnaphthalene	142	7 879	7 879 (1 190)	444206	80 0000	80 000
67 Nexachlorocyclopentadione	237	8 087	8.087 (0.836)	121425	80.0000	80 000
69 2,4,6-Trichlorophenol	196	8 237	8 237 (0 851)	127047	80.0000	80 000
70 2,4,5-Trichlorophenol	196	8 306	8 306 (0 859)	136886	80 0000	80 000
73 2-Chloronaphthalene	1.62	8 547	8 547 (0 883)	419213	80 0000	80 000
77 2-Nitrosmiline	65	8 814	8.814 (0 911)	153125	80 0000	80 000
80 Dimothylpathalate	163	9 257	9 257 (0 957)	507295	80 0000	
82 2,6-Dimitrofoluene	165	9 380	9 380 (0 970)	114569	80 0000	80 000
83 Acenaphthyiene	152	9 358	9 358 (0 967)	725193	80 0000	80 000
85 3-Nitroani_ne	138	9 652	9.652 (0 998)	14355B	80 0000	80 000
86 Acenaphthene	153	9 738	9 738 (1 007)	445205	80 0000	80 000
87 2,4-Dinitrophenol	184	9 866	9 866 (1 020)	60669		80 000
69 4 Nitrophenol	109	10 042	10.042 (1.038)	71944	0000 08	80 000
90 Dibenzefaran	168	10 074	10.074 (1.038)		80.0000	80 000
91 2,4-Dimitrotoluene	165	10 203	10 203 (1 055)	626236	80 0000	80 000
95 2,3,5,6-Tetrachlorophonol	232	10 253	10 357 (1 071)	159148	80 0000	80 000
92 2,3,4,6-Totrachlorophenol	232	10.459	10.459 (1 081)	104207	80 0000	80.000
96 2 Naphthylamine	143	10.433	10.427 (1 078)	104594	80.0000	80.000
97 Dietnylphthalate	149	10 790	10.427 (1 0767	448558	80 0000	80 000
98 Fluorene	166	10 /30	•	515007	80 0000	80 000
99 4-Chlorophenyl phenyletaer	204	10 865	10 828 (1 119)	511605	80.0000	80 000
1 ( 4-Nitroam ine	138	10 993	10.865 (1 123) 10 993 (1 136)	229949	80 0000	80 000
102 4,6-Dinitro-2-methylphenol	198	11 095		150312	80 0000	80 000
103 N-Nitrosodiphenviamine (1)	169	11 159	11 095 (0 854)	87275	87 0000	80 000
104 1,2 Dippenylhydrazine	77	11.218	11 259 (0 859)	376808	80 0000	80 000
112 4-Bromophenyl pleayletter	248	11 976	11 218 (0 864)	643013	80 0000	80 000
113 Hexachlorobenzene	284	12 281	11 9/6 (0 922) 12 281 (0 946)	127898	80,0000	80 000
1_7 Pertachlorophenol	266			13738/	80.0000	80 000
122 Phonanthrene	178	12 724 13 045	12 724 (0 980)	81986	80.0000	80 000
123 Antoracche	178	13 151	13 045 (1 005)	743676	80 0000	80 000
12º Carbazole	167			770576	80 0000	80 000
130 Di-n Buty_pathalate	149	13 589 14,727		743910	80 0000	80 000
_3r F uoranchene			14 727 (1 134)	920830	80.0000	80 000
136 Berzidine	202			795806	80 0000	80 000
137 Pyrene	184		16 410 (0 836)	373875	80.0000	80 000
144 Buty_nemzylphthalate	202	16.528	•	836365	80 0000	80 000
149 3,3'-D_chlorobenzidine	149	18 472	18 472 (0 941)	451147	80 0000	80 000
	252	19 642	19 642 (1 001)	294463	80.0000	80 000
150 Benzo(a)Anthracene	228	19 589	19 589 (0 998)	790186	80 0000	80 000

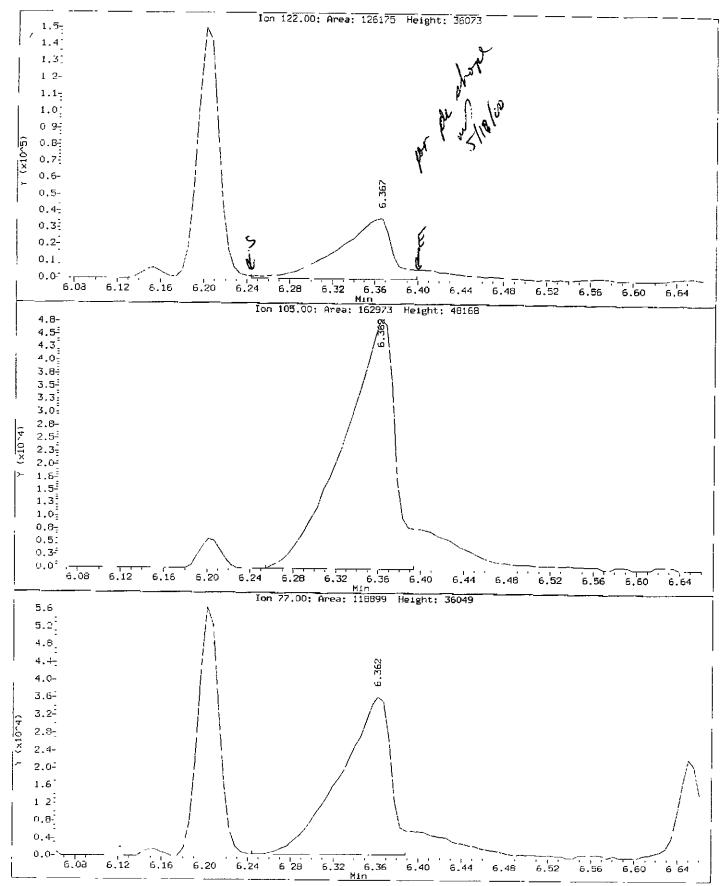
							AMOUN	TS
Compounds	QUANT SIG					CA!	L-AMT	ON-COL
•	MASS	RT	EXP RT	REL RT	RESPONSE	{	ng)	( ng)
要T (海南亚亚亚) 葡萄亚亚亚 (苏南非香油- 线性		==	=======	=\ ====	===:::::::::::::::::::::::::::::::::::	==		*=====
131 Chrysene	228	19 696	19 696	(1 003)	761847	80	0000	80 000
153 ols(2-ethylhexy*)Phthalaco	149	20.048	20 048	(1 021)	632878	80	0000	80 000
155 Di n-ostylphinalate	149	21 517	21 517	(0 936)	1119934		0000	80 000
157 Benzo(b)l_woranthene	252	22 180	22 180	(0 965)	884166		0000	80 000
158 Benzo(k) fluoranthenc	252	22 244	22 244	(0 367)	975293		0000	
159 7,12 dimethylbenz[a]an-hracen	256	22 238	22 238	(0 967)	396831		0000	80 000
167 Benzo(a)pyrene	252	22 874	22.874	(0 995)	289772		0000	80 000
169 Trácno(1,2,3-cd)pyrene	276	25 294	25 294	(1 100)	1229271			80 000
170 Dibenz (a, h) anthracene	278	25 337		(1.102)	1047029		0000	80 000
17! Benzo(g,n,1)gerylene	276	25 892		(1.126)			0000	80 000
\$ 1/2 Nitrobenzeno-d5	82	5 742			1098690		0000	80 000
\$ 173 2-Fluorobiphenyl	172	8 376		(0 867)	276632		0000	80 000
\$ 174 Terpmenyl-a14	244			(0.866)	476261	90	0000	80 000
\$ 175 Phenol-d5	-	17 057	17 057		627118	80	0000	80 000
\$ 1/6 2-Fluoropheno:	90	4 759		(0 937)	300485	80	CCOO	80 000
-	112	3 749	3 749	(0.738)	223421	60	0000	80 000
\$ 177 2,4,6-Tribrumophenol	330	11 421	11 421	(0.879)	/2637	80	0000	80 000
\$ 1/8 2-Chiorophenol-d4	132	4 877	4 877	(0 960)	209317	80	0000	80 000
\$ 179 1,2-Dich_oronanzene-a4	152	5 288	5.288	(1 041)	149611	80	0000	80 000

# QC Flag Legend

M - Compound response manually integrated.

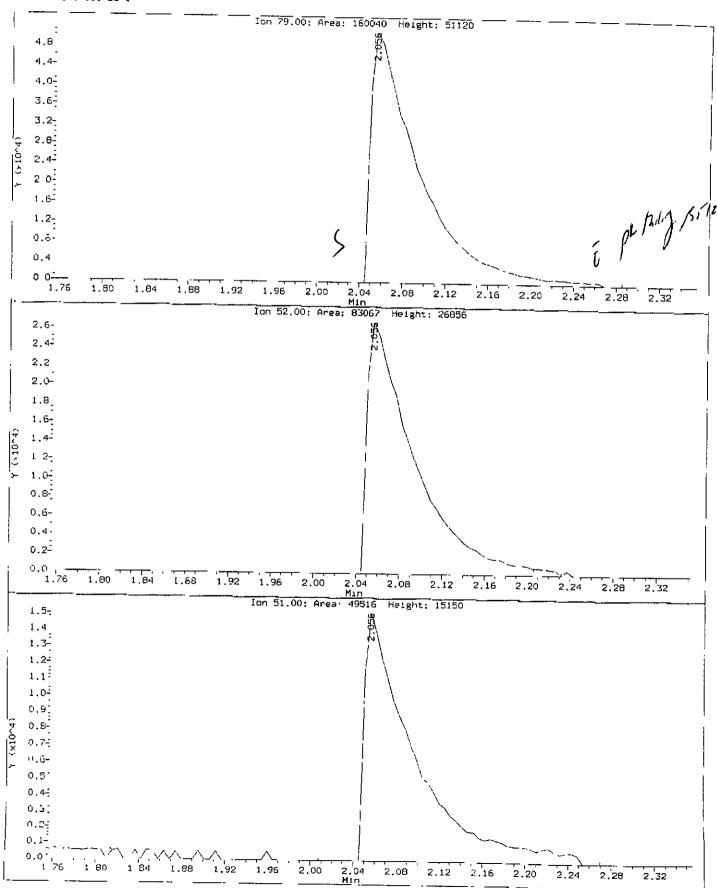
Data File: \\QPITPAO2\D\chem\71.1\s051900.b\S0519CC4.D Injection Date: 19-MAY-2000 14:32 Instrument: 71.1 Client Sample ID:

Compound, Benzoic Acid CAS Number: 65-85-0



Data File: \\OPITPAO2\D\chem\71.1\s051900.b\S0519CC4.D
Injection Date: 19-MAY-2000 14;32
Instrument: 71.1
Client Gample ID: SSTD080 ...

Compound: Pyridine CAS Number: 110-86-1



ß.

Semivolatile REPORT SW-846 Method 8270

Data file : \\QPITPA02\D\chem\71.i\s051900.b\S0519CC5.D

Lab Smp Id: sstd120 Client Smp ID: SSTD120

Inj Date : 19-MAY-2000 15:08

Operator : 45183 Inst ID: 71.i Smp Info : sstd120(60.0ug/ml) 194-175-13 8270/clp Misc Info : sstd120, s051900.b, 8270clp.m, 1-82701.sub, 1, 4

Comment

: \\QPITPA02\D\chem\71.i\s051900.b\8270clp.m Method Meth Date: 19-May-2000 16:29 bachas Quant Type: ISTD Cal Date : 19-MAY-2000 15:44

Cal File: S0519CC6.D Als bottle: 6 Calibration Sample, Level: 4

Dil Factor: 1.00000 Integrator: HP RTE

Target Version: 4.03

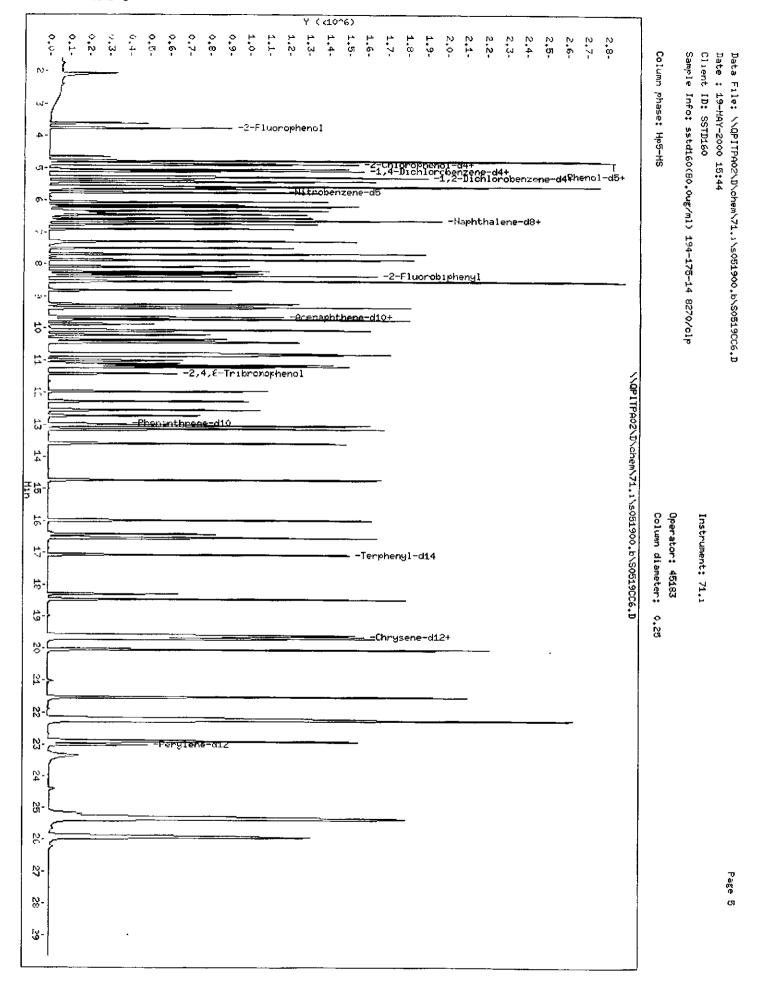
Processing Host: PITPC050

Compound Sublist: 1-82701.sub

							AMOUN	TS
,~,	Compounds	QUANT SIG					CAL AMT	ON-COL
-	•	MASS	RT	EXP RT	REL RT	RESPONSE	( ng)	( ng)
-	(秦孝孝孝,77世四四四四四7 1年3四四四年) 在秦莽年	22 × 2		=====	=====	3E3E##80	======	=======
	1,1-51010Denzene-da	152	5 071	5 071	(1 000)	91018	40 0000	
*	, rechirens encino	136	6 620	6 620	(1 000)	368841	40.0000	
*	2 IN C. aprilin DC-GID	164	9 665	9.665	(1 000)	212670	40 0000	
*	4 Phenanthrene-dl0	188	12 982	12 982	(1 000)	404163	40 0000	
*	3 Chrysene-d12	240	19 628	19 628	(1.000)	376940	40.0000	
*	n relytene-diz	264	22.988	22 988	(1 000)	495736	40 0000	
	13 N-Nitrosodimethylamine	74	2 047	2 047	(0 404)	140465	120 000	119 47
	10 Pyridire	79	2 047	2 047	(0 404)	265319	120.000	121 40
	19 Methyl methancsulfonate	80	3 586	3 586	(0 707)	232437	120 000	119.67
	22 Aniline	93	4.772	4 772	(0 941)	654969	120 000	114 58
	23 Phenol	94	4 772	4 772	(0 941)	523648	120 000	112 82
	24 bis(2-Chloroethyl)ether	93	4.846		(0 956)	434951	120 000	117.02
	25 2-Chlorophenol	128	4 889	4 889	(0 964)	371416	120 000	116 27
	27 1,3-Dichlorobenzene	146	5 033	5 033	(0 993)	406256	120 000	114 54
	28 1,4 Dichlorobenzene	146	5 092		(1 004)	411941	120 000	114.29
	29 1,2-Dicalorobenzenc	146	5 301		(1 045)	375651	120.000	117.29
	30 Benzyl Alcohol	108	5 258	5 258	•	279444	120.000	120 59
	31 2 Methylphenol	108	5 413	5 413	-	361956	120 000	116.29
	32 2.2' oxybis(1-Chloropropane)	45	5 429	5 429	•	5 171 03	120.000	114.35
	33 N-Nitroso-di-n-propylamine	70	5.600	5 600		312034	120.000	
	35 4-Metnylphenol	108	S 573	5.573		376899	120 000	119 85
	38 Hexach oroethane	117	5 642	5 642		162638	120.000	115 43
	39 Witrobenzene	71	S 760	5 760		476429	120.000	116 87
	44 Isophorone	82	6.038	6.038		795086		117.85
	45 2 Nirrophenol	139	6 150	6 150		213687	120 000	118 57
	46 2,4-Directhylphenol ,	107	6 198	6 198		386804	120 000	122 76
		•	0	V 138	(0 330)	300804	120 000	119 87

							AMOUN	TS
		QUAN'I 51G					CAL-AMT	ON-COL
Сопро	ounds	MASS	RT	EXP RT	REL RT	RESPONSE	(pg)	( ng)
==-		====	₩=		=====	=====	======	======
4	bis(2-Chloroctnoxy)mothane	93	6 326	6 326	(0 956)	504646	120 000	117 20
51	7,4-Dichlorophenol	162	6.449	6 449	(0 974)	299590	120 000	118 07
2د	? Benzoic Acia	122	6 390	6 390	(0 965)	229616	120 000	147 96
53	1,2,4-Tr_chlorobeczene	380	6 561	6 561	(0 991)	310565	120 ano	115 11
54	Naphthalene	128	6 647	6 647	(1 004)	1125211	120 000	113 11
55	4-Cnlorpanil ne	127	6 754	6 754	(1 020)	474328	120 000	118 99
59	Hexachlorobutadione	225	6 903	6 903	(1 043)	156918	120 000	113 78
62	4-Chloro 3-Methylphenol	107	7 491	7 491	(1 132)	356498	120 000	121 61
65	2-Methylnaphthalenc	142	7 688	7 688	(1 161)	752903	120 000	114 75
66	1-Methylnaphthalene	142	7 870	7 870	(1 189)	708225	120 000	115 23
67	Hexachlorocyc'cpcnLadiene	237	8 078	8.078	(0 836)	198479	120 000	123 60
59	2,4,6-Trichlorophenol	196	8 228	8 228	(0 851)	208719	120 000	120 51
70	2,4,5-Irichlorophenol	196	8 303	8 303	(0 859)	225056	120 000	121 77
73	2-CF oronaphthalene	162	8.543	8 543	(0 884)	654994	120.000	113 22
77	2-Nitroaniline	65	8 810	8.810	(0 912)	255998	120 000	124.88
80	Dimethylphthalate	163	9 254	9 254	(0 957)	821238	120 000	118 47
82	2, C-Dinitrofoluene	165	9 377	9 377	(0 970)	189150	120 000	123 39
83	Acchaphthylene	152	9.350	9 350	(0 967)	1138422	120 000	116 49
85	3-Vitroaniline	138	9 649	9.649	(0 998)	237519	120 000	123 76
86	Acenaphthene	153	9 735	9 735	(1 007)	710547	120 000	115 86
87	2,4-Dinitrophonel	184	9 863	9 863	{1 0201	111904	120 000	148 15
89	4 Nitrophenoi	109	10.044	10 044	(1 039)	128908	120 000	130 59
JC	Dibenzofuran	168	10 071	10 071	(1 042)	991648	120 000	115 08
91	2,4-Dinitrorolucne	165	10 205	10 205	(1 056)	268722	120 000	174 91
95	2 3,5,5-Tetrachlorophonol	232	10 354	10 354	(1 071)	175722	120 000	124.60
92	2,3,4,6-Tetrachlorophenol	232	10.456	10 456	(1 082)	172017	120 000	120.49
96	2-Naphthy_amine	143	10 424	10 424	(1 078)	703413	120 000	112 96
97	D_ethylphthalate	149	10 787	10.787	(1 116)	835129	120 000	118 38
98	∍luorene	166	10 824	10 824	(1 120)	813051	120 000	115 21
99	4-Caloropaenyi-phenylether	204	10 856	10 856	(1 123)	364585	120 000	115 64
100	4 N_troaniline	138	11 001	11 001	(1 138)	253630	120 000	124 82
	4,6-D.nitro-2-methylphenol	198	11 102	11 102	(0 855)	155763	120 200	135 05
- 03	N-Nitrosodiphenylamine (1)	169	11 155	11 155	(0 859)	620978	120 000	115 96
	1,2-Diphenylhydrazine	77	11 214	11 214	(0 864)	1049055	120 000	116 85
112	4-Bromophenyl-phenylethor	248	11 967	11 967	(0 922)	208810	120 000	115 71
113	Hexachlorobenzeno	284	12 272	12.272	(0 945)	223372	120 000	115 46
117	[cntach]orophenol	266	12 721	12.721	(0 980)	147159	120 000	137 39
122	Phonan .hrene	1.78	13 036	13 036	(1 004)	1193957	120 000	113.37
12.5	Anthracene	178	13 148	13 148	(1 013)	1231447	120 000	113.84
125	Carbazole	167	13 586	13.586	(1 046)	1218038	120 000	116 54
130	Di n-Butylphthalate	149	14 724	14.724	(1 134)	1507329	120 000	118 38
135	Fluoranthene	202	15 979	15.979 (	(1 231)	1295419	120 000	115 88
136	Benzidire	184	16 401	16 401 (	(0 836)	626658	120 000	130 40
137	Pyrene	202	16 524	16 524 (	(0 842)	1338555	120 000	119 54
	Butylbenzylphthalate	149	18 469	18.469 (	(0 941)	744600	120 000	125 31
149	3,3' Dichlorobenzidine	252	19 639	19.639 (	1 001)	497656	120 000	127 42
150	Benzo (a) Anthracene	228	19 585	19 585 (	0 998)	1278544	120 000	120 64

	0573377 077				AMOUN	TS.
Compounds	QUANT SIG				CAIAMT	ON-COL
* ***** '= *** '- *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *** - *	MAS5	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
15° Chrysene	# 1 12 2	==	****== =====	=======	*****	221:222
7.00	228	19 698	19 698 (1 004)	1215822	120 000	118 63
153 bis(2-ethy_hexyl)Fhthalate	149	20 039	20 039 (1 021)	1039131	120 000	124.45
155 Di-n-octylphthalate	149	21.514	21 514 (0 936)	1898233	120 000	118 04
157 Benzo(s)[luoranthene	252	22 187	22 187 (0 965)	1539369	120 000	118 38
158 Ber.zo(k)11uoranthene	252	22 251	22 251 (0 968)	1379772	120 000	106 08
159 7,12-dimethylbenz[a]ant.wacon	256	22 246	22 246 (0 968)	649032	120 000	-
167 Renzo (a pyrene	252	22 881	22 881 (0 995)	1485055	120 COO	112 63
169 ludeno(1,2,3-cd)pyzene	276	25 296	25 296 (1 100)	2051034	120 000	115 /2
170 Dibenz(a,h)anthracene	278	25 344	25 344 (1 102)	1761245		116.58
171 Renzo(g,n,i)perylone	276	25 910	25 910 (1 127)		120 000	117 16
\$ 1/2 Nitrobenzene-d5	82	5 739	5 739 (0 867)	1846045	120 000	118 10
\$ 173 2 Fluorobiphenyl	172	8 372	•	445/1R	120 000	119 36
\$ 174 Terphonyl-d14	244		8 372 (0.866)	751389	120 000	115 42
5 175 Phenol-d5	99	17 053	17 053 (0 869)	1015026	120 000	120 73
\$ 1/6 2 Fluorophenol		4 761	4 761 (0 939)	476073	120 000	116 01
\$ 177 2,4,5 lribromophenol	112	3 746	3 746 (0 739)	362026	120 000	118 71
	330	11 417	11 417 (0 879)	122400	120.000	122 54
TOPIN NOT GI	132	4 873	4 873 (0 961)	334707	120 000	116 74
\$ 179 %,2-DichTorobenzene-d4	152	5 285	5 285 (1 042)	229760	120 000	113 37



Data File: \\QPITPA02\D\chem\71.i\s051900.b\S0519CC6.D Report Date: 22-May-2000 08:01

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file: \\QPITPA02\D\chem\71.i\s051900.b\S0519CC6.D Lab Smp Id: sstd160 Client Smp ID:

Client Smp ID: SSTD160

Inj Date : 19-MAY-2000 15:44

Operator: 45183 Inst ID: 71.i Smp Info : sstd160(80.0ug/ml) 194-175-14 8270/clp Misc Info: sstd160,s051900.b,8270clp.m,1-82701.sub,1,5

Comment

: \\QPITPA02\D\chem\71.i\s051900.b\8270clp.m Method Meth Date : 22-May-2000 07:54 bachas Quant Type: ISTD

Cal File: S0519CC6.D

Calibration Sample, Level: 5

Cal Date: 19-MAY-2000 15:44
Als bottle: 7
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: 1-82701.sub

Target Version: 4.03 Processing Host: PITPC050

						ris /s/m/w
					NUOMA	ITS /
	QUANT STG				CAL-AMT	ON-COT
Compounds	MASS	RT	EXP RT RFL RT	RESPONSE	( ng)	( ng)
	#= -=	==	=======================================	=======	=======	프루=크 링스스
* 1 1,4-Bichlorobenzone-d4	152	5 085	5 085 (1 000)	84332	40 0000	
* 2 Naphthalene-d8	136	5 629	6 629 (1 000)	342337	40 0000	
* 3 Acemaphthene-d13	164	9.679	9 679 (1 000)	196197	40.0000	
* 4 Phenanthrene-d'0	188	12,991	12 991 (1 000)	372498	40 0000	
* 5 Chrysche-d12	240	19 637	19 637 (1.000)	342183	40 0000	
* 6 Perylene d12	264	23 003	23 003 (1 000)	164496	40.0000	
13 N-Nitresodimethylamine	74	2 061	2 061 (0 405)	136800	160 000	140 99
10 Pyridine	79	2 056	2 056 (0 404)	258214	160 000	141,84 (M)
19 Methyl mernanesulfonate	80	3 595	3 595 (0 707)	286226	160.000	159.57
22 Aniline	93	4 781	4 781 (0 940)	791706	10 000	154 55
23 Phenol	94	4 781	4 781 (0 940)	637759	160 000	153 52
24 pis(z-Chloroethyl)ether	93	4 855	4 855 (0.955)	535994	160 000	157.67
25 2-Chlorophenol	128	4 898	4.898 (0 963)	454329	160 000	155.77
27 1,3-Dichlorobenzene	146	5 042	5.042 (0 992)	496176	160 000	155 02
29 1,4 Jichlorobenzenc	146	5 101	5 101 (1 003)	50:146	160 000	154 61
29 1,2-Dichlorobenzene	146	5 309	5 309 (1 044)	453120	160 000	153 03
30 Benzyl Alconol	108	5 272	5.272 (1 037)	343171	160 000	158.73
31 2-McLhylpheno_	108	5 422	5 422 (1 066)	444907	160 000	156 00
32 2,2'-oxybis("-Chloropropane)	45	5 443	5 443 (1 070)	652440	160 000	154 68
33 N-Nitroso-di-n-propylamine	70	5 609	5 609 (1 103)	384578	160 000	158 94
33 4-Me nylpachol	108	5.582	5.582 (1 098)	457296	160 000	154 51
38 Hexachloroethane	117	5 651	5.651 (1 111)	200617	160 000	157 04
39 Nitrobenzene	77	5 769	5.769 (0 870)	590453	160 000	357 67
44 Tsopherone	82	€ 052	6 052 (0 913)	986481	160 000	158 50
45 2-Nitropienol	139	6 159	6 159 (0 929)	266358	160 000	159 57
46 2,4-Dimotrylphenol	107	6 212	6 212 (0 937)	478492	160 000	158 06
			•	=		

					AMOUN	ITS
Compounds	QUANT SIG				CAL-AM1	ON COL
-	MASS	RT	EXP RT REL RT	RESPONSE	( ng)	( ng)
	#22L	₩ 2	Z====# ===#==	<b>→ ← = □ = 1</b>	===-===	~##E#==
47 bis(2-Coloroethoxy) methane	93	6 335	6 335 (0 956)	623042	160 000	156 88
51 2,4-Dichloropheno?	162	6 463	6 463 (0 975)	369166	160 000	156 92
52 Berzold Acid	122	6 415	6 415 (0 968)	292754	160 000	170 24(A)
53 L,2,4-Trichloropenzene	180	6 575	6 575 (0 992)	380572	160 000	155 35
54 Naphthalere	128	6 656	6 656 (1.004)	1373563	160 000	153 71
F5 4 Chloroaniline	127	6 763	6 763 (1 020)	578694	160 000	157 02
59 Hexachicroputadiere	225	6 912	6 912 (1 043)	192320	160 000	153 92
62 4-Chloro-3-Mathyiphenol	107	7 505	7 505 (1 132)	442396	160 000	160 02(A)
65 2 Methylnaphthalone	142	7 703	7 703 (1.162)	918080	160 000	154 79
66 1-Methy_naphthalene	142	7 884	7.884 (1 189)	860697	160 000	155 84
67 Hexachlorocyclopentadiene	237	8 087	8 087 (0 836)	245794	160 000	159 29
69 2,4,6-1richlorophenol	196	8 242	8 242 (0 852)	256192	160 000	158 99
70 2,4,5 Trichlorophenol	196	8 317	8 317 (0 859)	272800	160 000	158 04
/3 2 Chloronaphthalene	162	8 557	8.557 (0 884)	790346	160 000	153 61
/7 2 Nitroaniline	65	8 825	8 825 (0 912)	317545	160 000	161 23(A)
80 Dimothylubthalate	163	9 263	9 263 (0 957)	1006062	160 000	157 65
82 2,6 Jinitrotoluene	165	9 391	9.391 (0 970)	234010	160 000	160 Ol(A)
83 Adenaphthylend	152	9 364	9 364 (0 967)	1379297	160 000	155 42
85 3-Nitroan_line	138	9 663	9 663 (0 998)	293952	160 000	160 21 (A)
66 Acchaphthone	153	9 749	9.749 (1 007)	865826	160 000	156 09
8/ 2,4 Dinitropheno	184	9 882	9 882 (1 021)	146949	160 000	173 62 (A)
89 1-Nitrophenol	109	10 064	10 064 (1 040)	163553	160 200	165 31(A)
∃0 Dibenzofuran	168	10.085	10 085 (1 042)	1197125	160 000	154 72
91 2,4-Dimitrofoluone	165	10.219	10 219 (1 056)	329538	160 000	161 11(A)
95 2,3,5,6-Totrachlorophenol	232	10 363	10 363 (1 071)	213723	160 000	160 34 (A)
92 2,3,4,6 Tetrachlorophenol	232	10 470	10 470 (1 082)	204574	160 000	156 54
96 2-Naphthylamine	143	10.443	10 443 (1 079)	838532	160 000	152 93
9/ Diethylphthalale	149	10 801	10 801 (1 116)	1025661	160 000	157 99
98 Fluorene	166	10 839	10 839 (1 120)	978880	160 000	154 79
99 4-Chlorophelyl-phenylethor	204	10.871	10 871 (1 123)	439916	160 000	154 78
100 4-Nitrosniline	118	11.020	11 020 (1,139)	313278	160 000	161 63 (A)
102 4,6 Dinitro-2-methylphenol	198	11 116	11 116 (0 856)	196078	160 000	165 64 (A)
103 N-Nitrosodiphonylamine (1)	169	11 175	11 175 (0 860)	738125	160 000	156 82
104 1,2 Diphenylhydrazine	17	11.228	11 228 (0 864)	1294965	160 000	156.89
312 4-Bromophenyl-phenylether	248	11 976	11 976 (0 922)	254204	160 000	155 84
113 Hexachlorobenzone	284	12 286	12 286 (0 946)	272560	160 000	155 69
117 Pentachlorophenol	266	12 730	12 730 (0 980)	183247	160 000	
122 Phonant hireno	178		13 050 (1 005)	1453970	160 000	165 23(A) 154.52
123 Autoracene	178		13 162 (1 013)	1496932	160 000	154.52
126 Carbazole	167	13 600	13.600 (1 047)	1483447	150 000	154 O1 156 10
ido Dn-ButylphuhaTato	149		14 733 (1 134)	1832482	160 000	155 94
135 Filiprantiene	202	15 994	15 994 (1 231)	1565039	160.000	155 94
136 benzidine	184	16 416	16 416 (0 836)	764670	160 000	163 15 (A)
137 Pyrene	202	16 539	16 539 (0 842)	1620456	160 000	
144 Butylberzylphihalato	149	18 478	18 478 (0 941)	913162	160 000	158 82
149 3,3'-Dicalorobenzidine	252		19 653 (1 001)	606162	160 000	162 32(A)
150 Benzo(a)Anthracene	228		19 605 (0 998)	1567751	160 000	163 67(A)
		. =	,,,,		200 000	160 72(A)

Page 3

Data File: \\QPITPA02\D\chem\71.i\s051900.b\S0519CC6.D Report Date: 22-May-2000 08:01

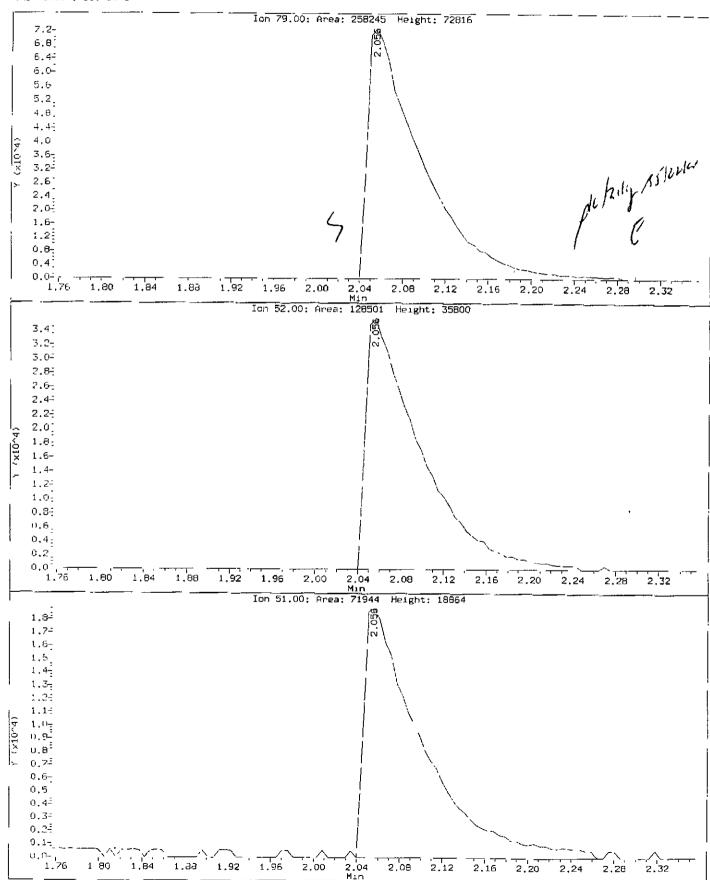
AMOUNTE OHANT SIG CAL AMT ON-COL Compounds MASS RT EXP RT REL R1 RESPONSE ( ng) (ng) **电影:10000 10000 "00000 "0000** # F 7 # == ----- \*\*=== ====== == **= = = = =** ===;=== 151 Chrysene 228 19 712 19 712 (1 004) 1478824 160 000 158 97 153 bis (2 ctay)hexy\_)Phthalate 149 20 048 20 048 (1 021) 1275518 160 000 161.98(A) 155 Di-n-octylphthalate 149 21.528 21.528 (0 936) 2328181 160 000 154 81 157 Benzo(b) f Loranthene 252 22 207 22 207 (N 965) 2058443 160 000 165 28 (A) 158 benzo(k)fluoranthone 252 22 276 22 276 (0 968) 1779718 160 000 145 83 (M) 159 7,12-dimethylbonz[a]anthracen 256 27 265 22 265 (0 958) 789180 160 000 152.68 167 Berzo(a)pyrano 22 901 22 961 (0 996) 252 1840120 160 000 155 81 169 Indenc(1,2,3-cd)pyrone 276 25 316 25 316 (1 101) 2590118 160 000 157 30 170 Disenz(a, h) anthracene 2/8 25 374 25 374 (1 103) 2199257 160 000 157 05 17t Benzo(q,h,1)perylene 276 25 935 25 935 (1 127) 2325283 160 000 157 66 \$ 172 Nitroberzene-d5 82 5 747 5 747 (0 867) 551411 160 000 158 11 \$ 173 2-Fluorotiphenyl 172 8 381 8 381 (0.866) 917510 160 000 155 33 \$ 174 Terphonyi-d14 214 17 067 17 067 (0 869) 1236326 160 000 160 21 (A) \$ 1/5 Pheno1-d5 99 4 770 4 770 (0 938) 576070 160 000 155.55 \$ 176 2-Fluorophonol 112 3 755 3 755 (0 738) 143659 160 000 158.36 \$ 177 2,4,6 Iribromopheno 330 11 426 11 426 (0 880) 150658 160 000 159 25 \$ 178 2 Chlorophenol-d4 132 4 882 4 882 (0 960) 408261 160 000 156 92 5 1/9 1,2-Dichloropenzere-d4 5 293 5 293 (1 041) 152 279585 160 000 153.50

## QC Flag Legend

- A Target compound detected but, quantitated amount exceeded maximum amount.
- M Compound response manually integrated.

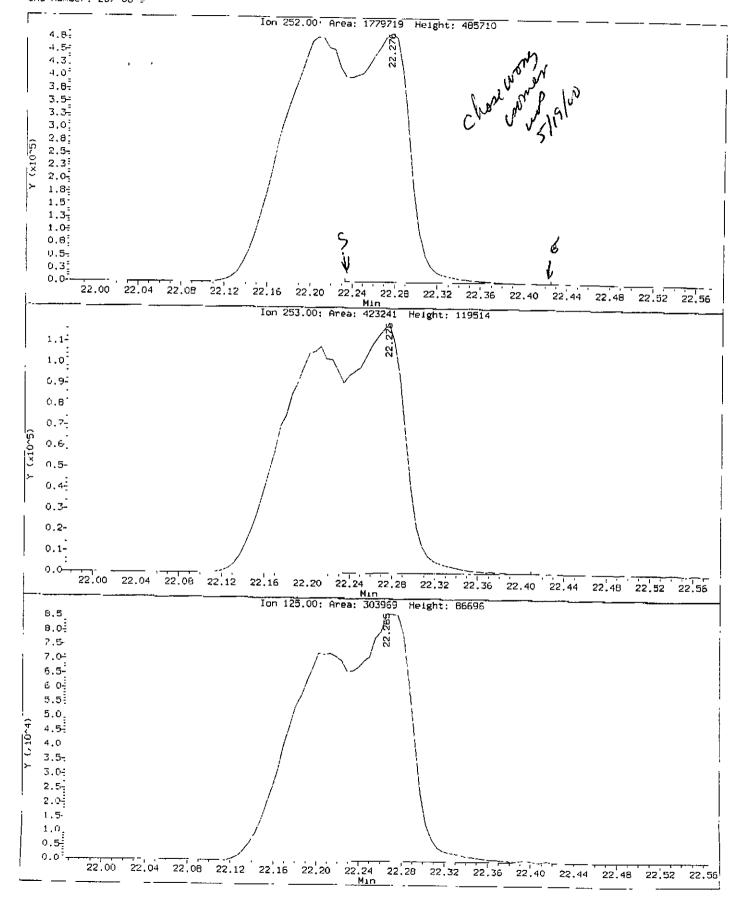
657 268 Data File: \QPITPRO2\B\chem\71.1\s051900.b\S0519006.D Injection Date: 19-MHY-2000 15:44 Instrument: 71.1 Client Sample ID. SSTD160

Compound: Pyridine CAS Mumber: 110–86–1



657 269

Compound: Benzo(k)fluoranthene CAS Number: 207-08-9



FORM 7

657 270 SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name:

Contract:

Lab Code: Case No.:

SAS No.: SDG No.: C0E130194

Instrument ID: 71 Calibration Date: 05/22/00 Time: 1127

Lab File ID: S0522CCC Init. Calib. Date(s): 05/19/00 05/19/00

Init. Calib. Times: 1320 1544

GC Column: HP5-MS ID: 0.25 (mm)

			MIN	2.5	MAX
COMPOUND	RRF	RRF50	RRF	%D	%D
######################################	2 010	1.933			20.0
Phenol	2.010	1.567			50.0
bis(2-Chloroethyl)ether	1.392	1.398			50.0
2-Chlorophenol	1.542	1.590			50.0
1,3-Dichlorobenzene		1.608			20.0
1,4-Dichlorobenzene	1.564				50.0
1,2-Dichlorobenzene	1.441				50.0
2-Methylphenol	2.038				50.0
2,2'-oxybis(1-Chloropropane)	1.419	1.319			50.0
4-Methylphenol Hexachloroethane	0.608	0.599		1 5	50.0
Hexachioroethane	0.608				50.0
Nitrobenzene	0.437				
Isophorone					20.0
2-Nitrophenol	0.190				50.0
2,4-Dimethylphenol	0.350			•	50.0
bis (2-Chloroethoxy) methane_	0.465				50.0
N-Nitroso-di-n-propylamine_	1.143	1	L .		20.0
2,4-Dichlorophenol	0.274				50.0
1,2,4-Trichlorobenzene	- 1.064				50.0
Naphthalene		i .	1		50.0
4-Chloroaniline	0.431 0.148				20.0
Hexachlorobutadiene	- 0.148				20.0
4-Chloro-3-Methylphenol	0.703				50.0
2-Methylnaphthalene	0.703				50.0
Hexachlorocyclopentadiene	0.304				20.0
2,4,6-Trichlorophenol	0.328				50.0
2,4,5-Trichlorophenol	1.072				50.0
2-Chioronaphtharene	0.389			11.0	50.0
Dimethylphthalate	1.299	Į.			50.0
	1.822				50.0
Acenaphthylene	0.290				
2,6-Dinitrotoluene 3-Nitroaniline	0.364			12.1	50.0
Acenaphthene	- 1.143		1		
2,4-Dinitrophenol	0.151				50.0
I	- 0.190				50.0
4-Nitrophenol Dibenzofuran	1.602				
2,4-Dinitrotoluene	- 0.408	1			
2,4-Diffictocordene	-  0.400	0.515		]	

page 1 of 3

FORM VII SV

# FORM 7 SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: Contract:

Lab Code: Case No.: SAS No.: SDG No.: C0E130194

Instrument ID: 71 Calibration Date: 05/22/00 Time: 1127

Lab File ID: S0522CCC Init. Calib. Date(s): 05/19/00 05/19/00

Init. Calib. Times: 1320 1544

GC Column: HP5-MS ID: 0.25 (mm)

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
Diethylphthalate	1.324	1.176	0.01	11.2	
4-Chlorophenyl-phenylether_	0.586	0.589	0.01		50.0
Fluorene	1.311	1.275	0.01		50.0
4-Nitroaniline	0.386	0.316	0.01		50.0
4,6-Dinitro-2-methylphenol	0.118	0.103	0.01		50.0
N-Nitrosodiphenylamine (1)	0.526	0.544	0.01		20.0
4-Bromophenyl-phenylether	0.177	0.194			50.0
Hexachlorobenzene	0.190	0.205			50.0
Pentachlorophenol	0.109				20.0
Phenanthrene	1.029	1.059	0.01		50.0
Anthracene	1.058	1.077	0.01		50.0
Carbazole	1.027	1.001	0.01		50.0
Di-n-Butylphthalate	1.254	1.195			50.0
Fluoranthene	1.095	1.090		0.4	
Pyrene	1.187	. 1.086		8.5	50.0
Butylbenzylphthalate	0.638	0.581			50.0
3,3'-Dichlorobenzidine	0.420	0.453			50.0
Benzo(a) Anthracene	1.129	1.152			50.0
Chrysene	1.086	1.111			
bis(2-ethylhexyl)Phthalate	0.895	0.834			
Di-n-octylphthalate	1.289	1.217	0.01	5.6	
Benzo (b) fluoranthene	1.061	1.082			50.0
Benzo(k) fluoranthene	1.088	1.101		1.2	
Benzo (a) pyrene	1.026				
Indeno(1,2,3-cd)pyrene	1.414		0.01		50.0
Dibenz (a, h) anthracene	1.207	1.296	0.01	7.4	50.0
Benzo(g,h,i)perylene	1.259	1.339	0.01	6.4	50.0
Pyridine	0.924	1.539	0.01		50.0
N-Nitrosodimethylamine	0.494	0.877	0.01		50.0
Aniline	2.479	2.304	0.01		50.0
Benzyl Alcohol	1.018	0.934	0.01		50.0
Benzoic Acid	0.177		0.01	43.5	50.0
1-Methylnaphthalene	0.659	0.627			50.0
2,3,4,6-Tetrachlorophenol	0.267		0.01		50.0
2,3,5,6-Tetrachlorophenol	0.266	0.252	0.01	5.3	50.0
1,2-Diphenylhydrazine	0.885	0.901	0.01		50.0
Benzidine	0.520		0.01		50.0
		<u> </u>			

page 2 of 3

FORM VII SV

657 272

#### FORM 7 SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name:

Contract:

Lab Code: Case No.:

SAS No.: SDG No.: COE130194

Instrument ID: 71

Calibration Date: 05/22/00 Time: 1127

Lab File ID: S0522CCC

Init. Calib. Date(s): 05/19/00 05/19/00

Init. Calib. Times: 1320 1544

GC Column: HP5-MS ID: 0.25 (mm)

COMPOUND	RRF	RRF50	MIN RRF	%D	MAX %D
**************************************	0.852	0.759	0.01	10.9	50.0
Methyl methanesulfonate 2-Naphthylamine	1.150	0.945	0.01	17.8	50.0
7,12-dimethylbenz[a]anthrace	0.457	0.444	0.01	2.8	50.0
Nitrobenzene-d5	0.405	0.394	0.01	2.7	50.0
2-Fluorobiphenyl	1.213	1.340 0.837	0.01		50.0
Terphenyl-d14 Phenol-d5	0.894 1.785	1.688	0.01	5.4	
2-Fluorophenol	1.335	1.355	0.01	1.5	50.0
2,4,6-Tribromophenol	0.099 1.250	0.100 1.235	0.01		
1,2-Dichlorobenzene-d4	0.878	0.902	0.01	2.7	50.0
		l <u></u> -	l	I	l

page 3 of 3

FORM VII SV

#### CONTINUING CALIBRATION COMPOUNDS PERCENT DRIFT REPORT

657 273

Instrument ID: 71.i Lab File ID: S0522CCC.D Analysis Type: NONE Injection Date: 22-MAY-2000 11:27

Lab Sample ID: sstd50

Method File: \\QPITPA02\D\chem\71.i\s052200.b\8270clp

	EXPECTED	MEASURED	}	MAX
COMPOUND	CONC	CONC	%D	₹D
1 Phenol	======= :   50.0000	48.8894	2.2	20
2 bis(2-Chloroethyl)ether	50.0000	48.8254	2 3	50.
3 2-Chlorophenol	50.0000	50.4420	0.9	50
4 1,3-Dichlorobenzene	50.0000	51.6951	3.4	50
5 1,4-Dichlorobenzene	50.0000	51 5161	3.0	20
6 1.2-Dichlorobenzene	50.0000	51 4915	3 0	50
7 2-Methylphenol	50.0000	47.9111	4 2	50.
8 2,2'-oxybis(1-Chloropropane)	50.0000	49.1843	1 6	50
9 4-Methylphenol	50.0000	47.4454	5.1	50.
11 1,4-Dichlorobenzene-d4	40.0000	40.0000	0.0	50
12 Hexachloroethane	50.0000	49.6800	0 6	50
13 Nitrobenzene	50.0000	49.0017	2.0	50.
14 Isophorone	50.0000	46 3839	7.2	50
15 2-Nitrophenol	50 0000	50.5185	1.0	20
16 2,4-Dimethylphenol	50.0000	48 8957	2.2	50
17 bis (2-Chloroethoxy) methane	50.0000	48.8324	2.3	50
18 N-Nitroso-di-n-propylamine	50 0000	44.8180	10.4	50.
18 2,4-Dichlorophenol	50.0000	50 8884	1 8	20.
19 1,2,4-Trichlorobenzene	50.0000	52 7620	5.5	50
20 Naphthalene	50.0000	51.4723	2.9	50.
21 4-Chloroaniline	50.0000	48 3655	3.3	50.
22 Hexachlorobutadiene	50.0000	54.3722	8 7	20
23 4-Chloro-3-Methylphenol	50.0000	43.7665	12.5	20.
24 2-Methylnaphthalene	50.0000	49.4468	1.1	50.
25 Hexachlorocyclopentadiene	50.0000	32 8036	34 4	50.
26 2,4,6-Trichlorophenol	50 0000	52.1118	4 2	20.
27 2,4,5-Trichlorophenol	50 0000	50.4151	0 8	50.
28 2-Chloronaphthalene	50.0000	53.6168	7.2	50.
29 2-Nitroaniline	50 0000	45.4736	9 1	50
30 Dimethylphthalate	50.0000	47 7428	4 5	50
31 Acenaphthylene	50 0000	50.8376	1 7	50
32 2,6-Dinitrotoluene	50.0000	46.2888	7.4	50
32 Naphthalene-d8	40 0000	40.0000	0 0	50
33 3-Nitroaniline	50 0000	45 2084	9.61	50
34 Acenaphthene	50.0000	50 7712	1.5	20
35 2,4-Dinitrophenol	50.0000	36.5147	27.0	50
36 4-Nitrophenol	50.0000	37.7721	24.5	50
37 Dibenzofuran	50.0000	50.7751	16	50
38 2,4-Dinitrotoluene	50 0000	44 1237	11.8	50

Data File: \\QPITPA02\D\chem\71.i\s052200.b/S0522CCC.D
Report Date: 05/22/2000

657 274

CONTINUING CALIBRATION COMPOUNDS PERCENT DRIFT REPORT

Instrument ID: 71.i Lab File ID: S0522CCC.D Analysis Type: NONE Injection Date: 22-MAY-2000 11:27

Lab Sample ID: sstd50

Method File: \\QPITPA02\D\chem\71.i\s052200.b\

	EXPECTED	MEASURED	- 1	XAM
COMPOUND	CONC	CONC.	₹D	*D
39 Diethylphthalate	50 0000	45.6782		50 0
40 4-Chlorophenyl-phenylether	50 0000	50 4630		50.0
41 Fluorene	50.0000	49 2542	1 5	50.0
42 4-Nitroaniline	50.0000	42.5780		50 0
43 4,6-Dinitro-2-methylphenol	50.0000	44.8612	10 3	50 0
44 N-Nitrosodiphenylamine (1)	50.0000	51 4102	2.8	20 0
45 4-Bromophenyl-phenylether	50 0000	54.1355	8.3	50.0
46 Hexachlorobenzene	50.0000	53.4779	7 0	50 0
47 Pentachlorophenol	50.0000	49.6633	0.7	20 0
48 Phenanthrene	50 0000	51 5293	3.1	50.0
49 Anthracene	50 0000	51.1209	2 2	50 0
50 Carbazole	50.0000	49 3025	1.4	50.0
51 Di-n-Butylphthalate	50.0000	48.4588	3 1	50.0
52 Fluoranthene	50.0000	50 2521	0 5	20 0
52 Acenaphthene-d10	40.0000	40 0000	0.0	50.0
53 Pyrene	50.0000	46 5459	6 9	50 0
54 Butylbenzylphthalate	50 0000	46 3629	7.3	50 0
55 3,3'-Dichlorobenzidine	50.0000	53.1632	6 3	50.0
56 Benzo (a) Anthracene	50.0000	50 9682	1.9	50.0
57 Chrysene	50 0000	51.0834	2 2	50.0
58 bis (2-ethylhexyl) Phthalate	50 0000	47.2990	5.4	50.0
59 Di-n-octylphthalate	50.0000	48.2218	3.6	20 (
60 Benzo(b) fluoranthene	[ 50 0000	51 0118	2.0	50.0
61 Benzo(k) fluoranthene	50 0000	51 1174	2.2	50 (
62 Benzo(a)pyrene	50 0000	51 9372	3 9	20 (
63 Indeno(1,2,3-cd)pyrene	50.0000	52 9159	58	50.6
64 Dibenz(a,h)anthracene	50 0000	53 0955	6 2	50.
65 Benzo(g,h,1)perylene	50.0000	52 6871	5 4	50
66 Pyridine	50.0000	73 9924	48.0	50
67 N-Nitrosodimethylamine	50.0000	77.8410	55.7	50
68 Anıline	50.0000	47 3099	5.4	50
69 Benzyl Alcohol	50.0000	46 6409	6.7	50
71 Benzoic Acid	50.0000	30.5833	38.8	50.
72 1-Methylnaphthalene	50.0000	48.2430	3.5	50.
76 2,3,4,6-Tetrachlorophenol	50.0000	48.5322	2.9	50.
77 2,3,5,6-Tetrachlorophenol	50.0000	47 6907	7 4.6	50.
78 1,2-Diphenylhydrazine	50 0000	50 8954	1.8	50.
78 Phenanthrene-d10	40.0000	40 0000	0 0	50
79 Benzidine	50.0000	1 60.0099	el 20 0	50.

Data File: \\QPITPA02\D\chem\71.i\s052200.b/S0522CCC.D Report Date: 05/22/2000

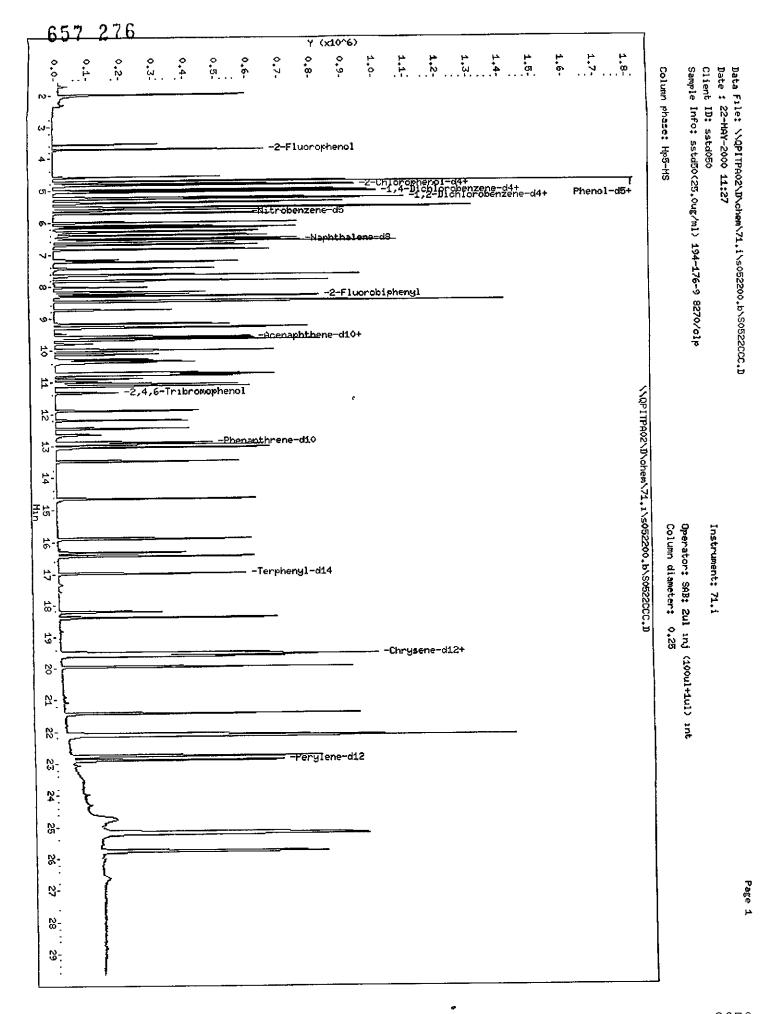
#### CONTINUING CALIBRATION COMPOUNDS PERCENT DRIFT REPORT

Instrument ID: 71.i Lab File ID: S0522CCC.D Analysis Type: NONE

Injection Date: 22-MAY-2000 11:27

Lab Sample ID: sstd50
Method File: \\QPITPA02\D\chem\71.i\s052200.b\

		E	XPECTED	MEASURED	1	MAX
CO	MPOUND	I	CONC.	CONC	*D	*D
	3世纪日本中国的西班牙里的中国的中国的中国的古代日代记记记	-				
87	Methyl methanesulfonate	1	50 0000	45 4607	9.1	50.0
90	Chrysene-dl2	1	40 0000	40 0000	0.0	50.0
101	Perylene-d12	1	40 0000	40 0000	0.0	50.0
114	2-Naphthylamine	1	50.0000	42 6206	14.8	50 (
143	7,12-dimethylbenz(a)anthracen	1	50.0000	49.0668	1 9	50 (
181	Nitrobenzene-d5	1	50 0000	49 2210	1.6	50.0
182	2-Fluorobiphenyl	1	50.00001	54 4698	8.9	50.0
183	Terphenyl-d14	1	50.0000	47.4612	5.1	50.0
184	Phenol-d5	1	50.0000	47.9495	4.1	50.0
185	2-Fluorophenol	1	50 0000	50.6878	1 4	50 (
186	2,4,6-Tribromophenol	1	50 0000	50.0751	0.2	50.
187	2-Chlorophenol-d4	1	50.00001	49 6478	0 7	50.
188	1,2-Dichlorobenzene-d4	1	50 0000	51.5543	3.1	50



1/1/1

Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522CCC.D Report Date: 22-May-2000 12:04

### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file : \\QPITPA02\D\chem\71.i\s052200.b\S0522CCC.D

Lab Smp Id: sstd50 Client Smp ID: sstd050

Inj Date : 22-MAY-2000 11:27 Operator : SAB: 2ul inj (100ul+1ul) int Inst ID: 71.i

Smp Info : sstd50(25.0ug/ml) 194-176-9 8270/clp

Misc Info : sstd50, s052200.b, 8270clp.m, 1-82701.sub, 2, 2

Comment

Method : \\QPITPA02\D\chem\71.i\s052200.b\8270clp.m Meth\_Date : 22-May-2000 11:59 bachas Quant Type: I Quant Type: ISTD Cal Date : 19-MAY-2000 15:44

Cal File: S0519CC6.D Als bottle: 2

Continuing Calibration Sample Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: 1-82701.sub Target Version: 4.03

Processing Host: PITPC050

							1551707
						AMOUN	,
	QUANT SIG					CAL-AMT	ON-COL
Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	(ng)	(ng)
美国克莱亚亚国际 医水水 医多耳节 化二氯苯酚 医克耳耳科	====	==	<b>40222</b> 0	*****			*****
* 1 1,4-Dichlorobenzene-d4	152	5.029	5.029	(1.000)	145518	40 0000	
* 2 Naphthalene-d8	136	6 562	6.562	(1.000)	535078	40 0000	
* 3 Acenaphthene-d10	164	9.596	9.596	(1.000)	265064	40.0000	
* 4 Phenanthrene-d10	188	12.898	12 898	(1.000)	427212	40.0000	
* 5 Chrysene-d12	240	19.533	19 53 <u>3</u>	(1.000)	446724	40.0000	
* 6 Perylene-d12	264	22.893	22. 893	(1.000)	599590	40.0000	
13 N-Nitrosodimethylamine	74	1 989	1.989	(0.396)	159465	50 0000	88 650
10 Pyridine	79	2.000	2.000	(0.398)	279889	50.0000	83 243 (M)
19 Methyl methanesulfonate	80	3 544	3 544	(0 705)	138055	50.0000	44.512
22 Aniline	93	4.730	4.730	(0.941)	419170	50.0000	46.475
23 Phenol	94	4.730	4 730	(0.941)	351579	50.0000	48 082
24 bis(2-Chloroethy1)ether	93	4 799	4 799	(0.954)	285018	50.0000	48.226
25 2-Chlorophenol	128	4.847	4.847	(0 964)	254360	50.0000	50.213
27 1,3-Dichlorobenzene	146	4.992	4.992	(0 993)	289190	50.0000	51.566
28 1,4-Dichlorobenzene	146	5 050	5.050	(1 004)	292500	50.0000	51.397
29 1,2-Dichlorobenzene	146	5 253	5 253	(1.045)	268741	50 0000	51.250
30 Benzyl Alcohol	108	5 216	5.216	(1.037)	169832	50.0000	45 850
<pre>31 2-Methylphenol</pre>	108	5 365	5 365	(1.067)	232682	50 0000	47 094
<pre>32 2.2'-oxybis(1-Chloropropane)</pre>	45	5.387	5.387	(1.071)	359586	50.0000	48.496
33 N-Nitroso-di-n-propylamine	70	5.542	5.542	(1.102)	181164	50 0000	43 555
35 4-Methylphenol	108	\$.526	5.526	(1.099)	239853	50 0000	46 461
38 Hexachloroethane	117	5.601	5.601	(1.114)	109043	50 0000	49.283
39 Nitrobenzene	77	5.707	5.707	(0 870)	283286	50.0000	48 463
44 Isophorone	82	5.985	5.985	(0.912)	440040	50.0000	45.320
45 2-Nitrophenol	139	6.097		(0.929)	128641	50.0000	50.634
46 2,4-Dimethylphenol	107	6.151	6.151	(0.937)	226731	50.0000	48 448

Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522CCC.D
Report Date: 22-May-2000 12:04

A7 bis (2-Chloroethoxy)methane   33   6.268   6.268   (0.955)   300352   50.0000   44 332								AMOUN	TS
47 bis (2-Chloroethoxy)methane	_		QUANT SIG					CAL-AMT	ON-COL
17	-		MASS	RT	EXP RT	REL RT	RESPONSE	( ng)	( ng)
Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary   Seminary									
Seminary Act   122   6 306 (6 951)   67015   50.0000   20 328   53 1,24-Trichlorobenses   100   6 509   6.509 (0.932)   200526   50.0000   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070   53.070				6.268	6.268	(0.955)	300352	50.0000	48 332
\$ 1,2,4-Trichlorobensene			162	6.396	6 396	(0 975)	185915	50.0000	50 713
Section   128			122	6 306	6 306	(0 961)	67015	50.0000	28 328
55 4-Chloroaniline			180	6 509	6.509	(0.992)	205626	50.0000	53,070
Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   Second   S		-	128	6 594	6.594	(1.005)	729762	50.0000	51 287
62 4-Chloro-3-Mechylphenol 107 7.433 7.433 (1.133) 180380 50 0000 42 279 65 2-Methylnaphthalene 142 7.630 7.630 (1.163) 459590 50.0000 48.849 65 1-Methylnaphthalene 142 7.630 7.630 (1.163) 459590 50.0000 47.546 67 Mexachlorocyclopentadiene 237 8.021 8.021 (0.836) 60733 50.0000 37.546 67 Mexachlorocyclopentadiene 237 8.021 8.021 (0.836) 60733 50.0000 52.521 00.24,6-Trichlorophenol 196 8.139 6.239 (0.859) 11603 50.0000 52.521 00.24,5-Trichlorophenol 196 8.239 6.239 (0.859) 11603 50.0000 50.0374 00.24,5-Trichlorophenol 196 8.239 6.859) 11603 50.0000 54.010 00.24,5-Trichlorophenol 196 8.239 6.859 11603 50.0000 54.010 00.24,5-Trichlorophenol 196 8.239 6.859 11603 50.0000 54.010 00.24 00.24 00.24 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00.25 00		•	127	6.696	6.696	(1.020)	274481	50.0000	47.677
65 2-Methylnaphthalene	59	Hexachlorobutadiene	225	6 845	6 845	(1.043)	108754	50 0000	55.030
Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Section   Sect		- ·	107	7.433	7.433	(1.133)	180380	50 0000	42 279
67 Rexachlorocyclopentadiene 237 8.021 8.021 (0.836) 60733 50.0000 30.123 69 2.4,6-Trichlorophenol 196 8.170 (0.851) 113420 50.0000 52.521 70 2.4,5-Trichlorophenol 196 8.229 8.239 (0.859) 116033 50.0000 52.521 73 2-Chloronaphthalene 162 8.475 8.475 (0.863) 33633 50.0000 54.010 77 2-Nitroaniline 65 8.736 8.736 (0.910) 114707 50.0000 44.456 80 Dimethylphthalate 163 9.180 9.180 (0.957) 404884 50.0000 46.985 82 2.6-Dinitrotoluene 165 9.303 9.303 (0.969) 87163 50.0000 45.310 83 Acenaphthylene 152 9.281 9.281 (0.967) 611562 50.0000 50.564 85 3-Nitroaniline 138 9.570 9.570 (0.977) 611585 50.0000 44.050 86 Acenaphthene 153 9.661 9.661 (1.007) 383137 50.0000 44.050 86 Acenaphthene 153 9.661 9.661 (1.007) 383137 50.0000 50.564 87 2.4-Dinitrophenol 184 9.789 9.789 (1.020) 34651 50.0000 50.564 87 2.4-Dinitrophenol 199 9.965 9.965 (1.038) 44938 50.0000 34.866 34-Nitrophenol 109 9.965 9.965 (1.038) 44938 50.0000 34.866 94-Nitrophenol 168 9.997 9.997 (1.042) 536417 50.0000 34.866 94-Nitrophenol 168 9.997 9.997 (1.042) 536417 50.0000 34.806 95 2.3.5,6-Tetrachlorophenol 232 10.280 (10.280 (10.77) 83428 50.0000 47.211 92 2.3.4,6-Tetrachlorophenol 232 10.280 (10.280 (10.77) 83428 50.0000 47.211 92 2.3.4,6-Tetrachlorophenol 232 10.280 (10.280 (10.77) 83428 50.0000 47.211 92 2.3.4,6-Tetrachlorophenol 232 10.280 (10.781 (1.071) 83595 50.0000 48.004 97 92-tethylphthalate 149 10.708 (1.078) 10.765 (1.081) 85197 50.0000 48.004 97 94-Chlorophenyl-phenylether 204 10.782 (1.745 (1.120) 422350 50.0000 41.084 97 Diethylphthalate 149 10.708 (1.078) 10.708 (1.124) 135077 50.0000 50.141 10.4-Stroaniline 138 10.000 (1.070) 853 55193 50.0000 43.929 10.380 (1.071) 8340 50.0000 43.929 10.380 (1.078) 8340 50.0000 43.929 10.380 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.782 (1.084) 10.7		= "	142	7.630	7 630	(1.163)	459580	50.0000	48.849
69 2,4,6-Trichlorophenol 196 8.170 8.170 (0.851) 113420 50.0000 52.531 70 2,4,5-Trichlorophenol 196 8.239 8.239 (0.859) 116033 50.0000 50.374 73 2-chloronaphthalene 162 8.475 8.475 (0.883) 383639 50.0000 54.010 77 2-Nitroaniline 65 8.736 8.736 (0.910) 114707 50.0000 44.456 80 Dimethylphthalate 163 9.180 9.180 (0.957) 404584 50.0000 46.985 2.6-Dinitrotoluene 165 9.303 9.03 (0.969) 87163 50.0000 50.654 82 2.6-Dinitrotoluene 155 9.281 9.281 (0.967) 611862 50.0000 45.310 83 Acenaphthylene 152 9.281 9.281 (0.967) 611862 50.0000 44.060 85 3-Nitroaniline 138 9.570 9.570 (0.977) 106185 50.0000 44.060 86 Acenaphthene 153 9.661 9.661 (1.007) 383137 50.0000 50.564 87 2,4-Dinitrophenol 184 9.789 9.789 (1.020) 34851 50.0000 34.866 89 4-Nitrophenol 109 9.365 9.365 (1.038) 44338 50.0000 35.652 90.000 Dihenzofuran 168 9.997 9.997 (1.042) 536417 50.0000 50.541 91.2-4-Dinitrotoluene 165 10.115 (0.151 (1.054) 115637 50.0000 42.803 95 2,3,5,6-Tetrachlorophenol 232 10.280 10.280 (0.071) 83428 50.0000 47.211 92.3,4,6-Tetrachlorophenol 232 10.376 (1.038) 44.008 81.358 95.0000 44.388 96 2-Naphthylamine 143 10.344 (1.078) 313265 50.0000 44.390 97 Fluorene 166 10.745 (1.078) 131365 50.0000 44.804 99 Hillorophenol 198 10.708 (1.116) 389505 50.0000 44.804 99 Hillorophenol 198 10.708 (1.116) 389505 50.0000 44.804 99 4-Chlorophenyl-phenylether 204 10.782 (1.784) (1.120) 422350 50.0000 44.804 99 4-Chlorophenyl-phenylether 204 10.782 (1.784) (1.120) 422350 50.0000 49.188 10.904 (1.2-Diphenylhydrazine 17 11.35 11.35 (0.863) 481308 50.0000 50.942 11.2 4-Bromophenyl-phenylether 248 11.888 11.889 (1.922) 10.380 50.0000 49.782 (1.124) 195077 50.0000 50.942 11.2 4-Bromophenyl-phenylether 248 11.889 11.889 (1.922) 10.380 50.0000 49.782 (1.124) 195077 50.0000 50.942 11.2 4-Bromophenyl-phenylether 248 11.889 11.889 (1.922) 10.380 50.0000 50.493 11.3 Harkachlorophenol 256 12.647 12.647 (0.981) 58168 50.0000 50.493 11.3 Harkachlorophenol 266 12.647 12.647 (0.981) 58168 50.0000 49.782 11.3 Harkachlorophenol 266 12.647 12.647 (0.981) 58168 50.0000 49.782			142	7.812	7.812	(1.190)	419108	50 0000	47.546
70 2,4,5-Trichlorophenol   196			237	8.021	8.021	(0 836)	60733	50.0000	30.123
73 2-Chloronaphthalene 162 8 475 (0 883) 3363639 50 0000 51.010 77 2-Nitroaniline 65 8.736 8.736 (0.910) 114707 50.0000 44.456 80 Dimethylphthalate 163 9.180 9 180 (0 957) 404584 50.0000 46.985 82 2.6-Dunitrotoluene 165 9.303 9 303 (0 969) 87163 50.0000 45.310 83 Acenaphthylene 152 9.281 9.281 (0 967) 611562 50 0000 50.654 85 3-Nitroaniline 138 9.570 9.570 (0 997) 106185 50 0000 44.060 86 Acenaphthene 153 9.661 9.661 (1.007) 338137 50 0000 50.564 87 2.4-Dunitrophenol 184 9.789 9.789 (1.020) 34851 50 0000 34.806 89 4-Nitrophenol 109 9.965 9.965 (1.038) 4493 50 0000 35.652 90 Dibenzofuran 168 9.997 9.997 (1.042) 536417 50 0000 55.61 91 2.4-Dunitrotoluene 165 10.115 (1.015 (1.054) 115637 50 0000 42.803 95 2.3.5.6-Tetrachlorophenol 232 10.280 (1.071) 83428 50.0000 47.211 92 2.3.4.6-Tetrachlorophenol 232 10.76 (1.038) 61.071 83428 50.0000 47.211 93 2.3.5.6-Tetrachlorophenol 232 10.76 (1.038) 61.091 8195 50.0000 48.188 96 2-Naphthylamine 143 10.344 10.344 (1.078) 313265 50.0000 48.188 97 Diethylphthalate 149 10.708 (1.16) 389505 50.0000 44.804 97 Diethylphthalate 149 10.708 10.708 (1.16) 389505 50.0000 44.804 98 Fluorene 166 10.745 10.745 (1.120) 422350 50.0000 40.918 100 4-Nitroaniline 138 10.900 10.900 (1.136) 104552 50.0000 40.918 101 4-Nitroaniline 138 10.900 10.900 (1.136) 104552 50.0000 40.918 102 4.6-chinitro-2-methylphenol 198 11.007 11.007 (0.858) 290432 50.0000 50.942 103 N-Nitrosodiphenylamine (1) 169 11.071 11.071 (0.858) 290432 50.0000 50.942 104 4.9-chinitro-2-methylphenol 226 12.647 12.647 (0.981) 5818 50.0000 50.942 113 Hexachlorophenol 256 12.647 12.647 (0.981) 5818 50.0000 50.943 114 Pentachlorophenol 256 12.647 12.647 (0.981) 5818 50.0000 50.935 115 Pentachlerone 178 12.951 12.951 11.004 5578 50.0000 51.491 129 Di-n-Butylphthalate 149 12.958 13.058 (1.012) 57522 50.0000 51.491 130 Di-n-Butylphthalate 149 13.058 13.058 (1.012) 57522 50.0000 51.481 130 Di-n-Butylphthalate 149 13.979 13.496 (1.046) 534645 50.0000 49.782 130 Di-n-Butylphthalate 149 18.379 18.379 (0.941) 324288 50.0000 45.522 149 3.3°-D	69	2,4,6-Trichlorophenol	196	8.170	8.170	(0 851)	113420	50.0000	52.521
77 2-Nitroaniline 65 8.736 (0.910) 114707 50.000 44.456 80 Dimethylphthalate 163 9.180 9 180 (0.957) 404584 50.0000 46.985 82 2.6-Dinitrotoluene 165 9.303 9.303 (0.969) 87163 50.0000 45.310 83 Acenaphthylene 152 9.281 9.281 (0.967) 611562 50.0000 50.554 85 3-Mitroaniline 138 9.570 9.570 (0.997) 106185 50.0000 44.060 86 Acenaphthene 153 9.661 9.661 (1.007) 383137 50.0000 50.564 87 2.4-Dinitrophenol 184 9.789 1.020 34851 50.0000 50.564 88 3-Nitrophenol 109 9.965 9.965 (1.038) 44938 50.0000 35.652 90 Dibenzofuran 168 9.997 9.997 (1.042) 536417 50.0000 55.541 91 2.4-Dinitrotoluene 165 10.115 (1.054) 115637 50.0000 42.803 95 2.3.5,6-Tetrachlorophenol 232 10.280 (1.071) 83428 50.0000 42.803 95 2.3.4,6-Tetrachlorophenol 232 10.280 (1.071) 83428 50.0000 47.211 92 2.3.4,6-Tetrachlorophenol 232 10.376 (1.081) 85197 50.0000 48.158 96 2-Naphthylamine 143 10.344 (1.034 (1.038) 131265 50.0000 44.390 97 Dischylphthalate 149 10.708 10.708 (1.16) 389505 50.0000 48.604 99 4-Chlorophenyl-phenylether 204 10.782 10.782 (1.124) 195077 50.0000 50.194 104 4-Kitroaniline 138 10.900 (1.036) 104552 50.0000 41.929 103 N-Nitrosodiphenylamine (1) 169 11.071 1.007 (0.853) 55193 50.0000 43.929 103 N-Nitrosodiphenylamine (1) 169 11.071 1.007 (0.853) 55193 50.0000 43.929 103 M-Ritroaniline 226 12 647 (1.64) 5863 481308 50.0000 50.942 114 4-Bromophenyl-phenylether 248 11.888 11.888 (0.922) 103801 50.0000 51.721 104 1.2-Diphenylhydrazine 77 11.135 (1.135) (1.063) 50.0000 51.721 105 1.2-Dhenylhydrazine 178 12.551 (1.2-S)1 (1.046) 534645 50.0000 49.782 113 Hexachlorophenol 266 12 647 (1.64) 534645 50.0000 49.782 114 Hexachlorophenol 278 12.551 (1.2-S)1 (1.04) 5352 50.0000 49.782 115 Pentachlorophenol 266 12 647 (1.64) 534645 50.0000 50.935 115 Pentachlorophenol 266 12 647 (1.64) 534645 50.0000 49.782 116 Benzaline 184 16.317 (1.63) (1.04) 53520 50.0000 49.782 117 Pentachlorophenol 266 12 647 (1.64) 534645 50.0000 49.782 118 Benzaline 184 16.317 (1.63) (1.04) 53520 50.0000 49.782 119 3.3*-Dichlorobenzaline 22 16.434 (1.64) (1.64) 534645 50.0000 45.5	70	2,4,5-Trichlorophenol	196	8.239	8.239	(0.859)	116033	50.0000	50.374
80 Dimethylphthalate 163 9.180 9.180 (0.957) 404584 50.0000 46.985 82 2.6-Dnittrotoluene 165 9.303 9.303 (0.959) 87163 50.0000 45.310 83 Acenaphthylene 152 9.281 9.281 (0.967) 611562 50.0000 50.654 85 3-Mitroanilline 138 9.570 9.570 (0.997) 106185 50.0000 50.654 85 3-Mitroanilline 138 9.570 9.570 (0.997) 106185 50.0000 44.060 86 Acenaphthene 153 9.661 9.661 (1.007) 383137 50.0000 50.564 87 2.4-Dnittrophenol 184 9.789 9.789 (1.020) 34651 50.0000 33.8066 89 4-Mitrophenol 109 9.9565 9.965 (1.038) 44938 50.0000 35.652 90 Dibenzofuran 168 9.997 9.997 (1.042) 536417 50.0000 35.652 90 Dibenzofuran 168 9.997 9.997 (1.042) 536417 50.0000 50.541 91 2.4-Dnittrotoluene 165 10.115 10.115 (1.054) 115637 50.0000 42.803 92 2.3.4,6-Tetrachlorophenol 232 10.280 10.280 (1.071) 83428 50.0000 47.211 92 2.3.4,6-Tetrachlorophenol 232 10.376 10.376 (1.061) 85197 50.0000 48.158 96 2-Maphthylamine 143 10.344 10.344 (1.038) 313265 50.0000 48.158 96 2-Maphthylamine 143 10.344 10.344 (1.038) 313265 50.0000 44 390 98 Fluorene 166 10.745 10.745 (1.120) 422350 50.0000 44 390 98 Fluorene 166 10.745 10.745 (1.124) 195077 50.0000 50.194 100 4-Nitroanilline 138 10.900 10.900 (1.136) 104552 50.0000 40.918 102 4.6-Dinittro-2-methylphenol 198 11.071 11.071 (0.858) 290432 50.0000 50.934 10.24 (4.6-Dinittro-2-methylphenol 198 11.071 11.071 (0.858) 290432 50.0000 50.942 12.4-Diphenylhydrazzne 77 11.135 11.135 (0.663) 481308 50.0000 50.942 12.4-Diphenylhydrazzne 77 11.138 11.135 (0.663) 481308 50.0000 50.942 12.4-Diphenylhydrazzne 77 11.138 11.135 (0.663) 481308 50.0000 50.942 12.4-Diphenylhydrazzne 77 11.138 11.135 (0.663) 481308 50.0000 50.942 12.4-Diphenylhydrazzne 178 12.951 12.951 11.004 53465 50.0000 49.782 13.496 13.496 (1.046) 534645 50.0000 49.782 13.496 13.496 (1.046) 534645 50.0000 49.782 13.496 13.496 (1.046) 534645 50.0000 49.782 13.496 13.496 (1.046) 534645 50.0000 49.782 13.496 13.496 (1.046) 534645 50.0000 49.782 13.496 13.496 (1.046) 534645 50.0000 49.782 13.496 10.446 40.446 40.446 40.446 40.446 40.446 40.446 40.446 40.446 40.446 40.44	73	2-Chloronaphthalene	162	8 475	8 475	(0 883)	383639	50 0000	54.010
82 2,6-Dinitrotoluene 165 9.303 9.303 (0.969) 87163 50.0000 45.310 83 Acenaphthylene 152 9.281 9.281 (0.967) 611562 50.0000 50.654 85 3-Natroaniline 138 9.570 (0.997) 106185 50.0000 44.060 86 Acenaphthene 153 9.661 9.661 (1.007) 383137 50.0000 50.564 87 2,4-Dinitrophenol 184 9.789 9.789 (1.020) 34851 50.0000 34.806 89 4-Nitrophenol 109 9.965 9.965 (1.038) 44938 50.0000 35.652 90.0000 34.806 89 4-Nitrophenol 109 9.965 9.965 (1.038) 44938 50.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.652 90.0000 35.0000			65	8.736	8.736	(0.910)	114707	50.0000	44.456
83 Acenaphthylene	80	Dimethylphthalate	163	9.180	9 180	(0 957)	404584	50.0000	46.985
85 3-Nitroaniline	82	2,6-Dinitrotoluene	165	9.303	9 303	(0 969)	87163	50.0000	45.310
86 Acenaphthene	83	Acenaphthylene	152	9 281	9 281	(0 967)	611562	50 0000	50.654
87 2,4-Dinitrophenol 184 9 789 9.789 (1.020) 34851 50 0000 34.806 89 4-Nitrophenol 109 9.965 9.965 (1.038) 44938 50 0000 35.652 90 Dibenzofuran 168 9 997 9 997 (1.042) 536417 50 0000 50 541 91 2,4-Dinitrotoluene 165 10 115 10.115 (1.054) 115637 50 0000 42.803 95 2,3,5,6-Tetrachlorophenol 232 10 280 10.280 (1.071) 83428 50.0000 47.211 92 2,3,4,6-Tetrachlorophenol 232 10 376 10.376 (1.081) 85197 50.0000 48.158 96 2-Naphthylamine 143 10.344 10.344 (1.078) 313265 50.0000 41 084 97 Diethylphthalate 149 10.708 10.708 (1.16) 389505 50.0000 44 390 98 Fluorene 166 10.745 10.745 (1.120) 422350 50.0000 44 390 99 4-Chlorophenyl-phenylether 204 10.782 10 782 (1.124) 195077 50.0000 40.918 102 4,6-Dinitro-2-methylphenol 198 11.007 11.007 (0.853) 55193 50.0000 40.918 102 4,6-Dinitro-2-methylphenol 198 11.007 11.007 (0.853) 55193 50.0000 40.918 104 1,2-Diphenylhydrazine 77 11.135 (1.863) 481308 50.0000 51.721 104 1,2-Diphenylhydrazine 77 11.135 (0.863) 481308 50.0000 51.721 104 2-Bromophenyl-phenylether 248 11.888 11.880 (0.922) 103801 50.0000 54.909 113 Hexachlorobenzene 284 12.193 12.193 (0.945) 109639 50.0000 54.096 117 Pentachlorophenol 266 12.647 12.647 (0.981) 58168 50.0000 49.783 (M) 122 Phenanthrene 178 12.951 (1.046) 534645 50.0000 49.783 (M) 122 Phenanthrene 178 12.951 (1.046) 534645 50.0000 49.783 (M) 123 Anthracene 178 12.951 (1.046) 534645 50.0000 49.783 (M) 124 Carbazole 167 13.496 13.496 (1.046) 534645 50.0000 49.782 (M) 125 Phenanthrene 202 15.889 15.889 (1.232) 582296 50.0000 49.782 (M) 136 Phenanthrene 202 15.889 15.889 (1.232) 582296 50.0000 49.782 (M) 137 Pyrene 202 16.434 16.434 (0.841) 606633 50.0000 45.746 (M) 148 Butylbenzylphthalate 149 18.379 18.379 (0.941) 324288 50.0000 53.964	85	3-Nitroaniline	138	9 570	9 570	(0 997)	106185	50 0000	44.060
89 4-Nitrophenol 109 9.965 9.965 (1.038) 44938 50 0000 35.652 90 Dibenzofuran 168 9 997 9 997 (1.042) 536417 50 0000 50 541 91 2.4-Dinitrotoluene 165 10 115 10.115 (1.054) 115637 50 0000 42.803 95 2,3.5,6-Tetrachlorophenol 232 10 280 10.280 (1.071) 83428 50.0000 47.211 92 2,3.4,6-Tetrachlorophenol 232 10 376 10.376 (1.081) 85197 50.0000 48.158 96 2-Naphthylamine 143 10.344 10.344 (1.078) 313265 50.0000 41 084 97 Diethylphthalate 149 10.708 10.708 (1.116) 389505 50.0000 41 390 98 Fluorene 166 10.745 10.745 (1.120) 422350 50.0000 48.604 99 4-Chlorophenyl-phenylether 204 10.782 10 782 (1.124) 195077 50.0000 40.918 102 4.6-Dinitro-2-methylphenol 198 11 007 11.007 (0.853) 55193 50.0000 40.918 102 4.6-Dinitro-2-methylphenol 198 11 007 11.007 (0.853) 55193 50.0000 40.918 104 1.2-Diphenylhydrazine 77 11 135 11 135 (0.863) 481308 50.0000 51.721 104 1.2-Diphenylhydrazine 77 11 135 11 135 (0.863) 481308 50.0000 54.096 117 Pentachlorophenol 256 12 647 12 647 (0.981) 58168 50.0000 54.096 117 Pentachlorophenol 256 12 647 12 647 (0.981) 58168 50.0000 54.096 117 Pentachlorophenol 256 12 647 12 647 (0.981) 58168 50.0000 51.481 123 Anthracene 178 12.951 12.951 (1.004) 557738 50.0000 48.761 130 Di-n-Butylphthalate 149 14 639 14.639 (1.135) 638330 50.0000 48.761 130 Di-n-Butylphthalate 149 14 639 14.639 (1.135) 638330 50.0000 49.782 136 Puoranthene 202 15 889 15.889 (1.232) 582296 50.0000 49.782 136 Benzialne 184 16.317 (6.317 (0.851) 53222 50.0000 45.746 144 Butylbenzylphthalate 149 18 379 18.379 (0.941) 324288 50.0000 45.746 144 Butylbenzylphthalate 149 18 379 18.379 (0.941) 324288 50.0000 53.964	86	Acenaphthene	153	9 661	9 661	(1.007)	383137	50 0000	50.564
90 Dibenzofuran 168 9 997 9 997 (1.042) 536417 50 0000 50 541 91 2.4-Dinitrotoluene 165 10 115 10.115 (1.054) 115637 50 0000 42.803 95 2.3.5,6-Tetrachlorophenol 232 10 280 10.280 (1.071) 83428 50.0000 47.211 92 2.3.4,6-Tetrachlorophenol 232 10 376 10.376 (1.081) 85197 50.0000 48.158 96 2-Naphthylamine 143 10.344 (1.078) 313265 50.0000 41 084 97 Diethylphthalate 149 10.708 10.708 (1.16) 389505 50.0000 44 390 98 Fluorene 166 10.745 10.745 (1.120) 422350 50.0000 48.604 99 4-Chlorophenyl-phenylether 204 10.782 10 782 (1.124) 195077 50.0000 50 194 100 4-Natroaniline 138 10.900 10 900 (1.136) 104552 50.0000 40.918 102 4.6-Dinitro-2-methylphenol 198 11 007 11.007 (0.853) 55193 50.0000 51.721 104 1,2-Diphenylhydrazine 77 11 135 11 135 (0.863) 481308 50.0000 50 942 12 4-Bromophenyl-phenylether 248 11.888 11 888 (0.922) 103801 50.0000 54.909 113 Hexachlorobenzene 284 12 193 12 193 (0.945) 109639 50.0000 54.909 113 Hexachlorophenol 266 12 647 12 647 (0.981) 58158 50.0000 54.909 113 Accachlorophenol 266 12 647 12 647 (0.981) 58158 50.0000 54.909 113 Anthracene 178 12.951 12.951 (1.004) 565778 50.0000 50.935 126 Carbazole 167 13.496 13.496 (1.046) 534645 50.0000 47.657 135 Pluoranthene 202 15 889 15.889 (1.325) 582296 50.0000 49.782 136 Benzialne 184 16.317 (0.835) 13.532 50.0000 45.574 137 Pyrene 202 16.434 16.337 (0.841) 324288 50.0000 45.522 149 3,3"-Dichlorobenzialne 252 19.543 19.543 (1.001) 253202 50.0000 53.9964	87	2,4-Dinitrophenol	184	9 789	9.789	(1.020)	34851	50 0000	34.806
91 2.4-Dinitrotoluene	89	4-Nitrophenol	109	9.965	9.965	(1.038)	44938	50 0000	35.652
91 2.4-Dinitrotoluene 165 10 115 10.115 (1.054) 115637 50 0000 42.803 95 2,3,5,6-Tetrachlorophenol 232 10 280 10.280 (1.071) 83428 50.0000 47.211 92 2,3,4,6-Tetrachlorophenol 232 10 376 10.376 (1.081) 85197 50.0000 48.158 96 2-Naphthylamine 143 10.344 10.344 (1.078) 313265 50.0000 41 084 97 Dischylphthalate 149 10.708 10.708 (1.166) 389505 50.0000 44 390 98 Fluorene 166 10.745 10.745 (1.120) 422350 50.0000 48.604 99 4-Chlorophenyl-phenylether 204 10.782 10.782 (1.124) 195077 50.0000 50 194 100 4-Nitroaniline 138 10.900 10.900 (1.136) 104552 50.0000 40.918 102 4.6-Dinitro-2-methylphenol 198 11.007 11.007 (0.853) 55193 50.0000 43.929 103 N-Nitrosodiphenylamine (1) 169 11.071 11.071 (0.858) 290432 50.0000 51.721 104 1,2-Diphenylhydrazine 77 11.135 11.135 (0.863) 481308 50.0000 50.942 112 4-Bromophenyl-phenylether 248 11.888 11.888 (0.922) 103801 50.0000 54.096 117 Pentachlorophenol 266 12.647 (0.981) 58168 50.0000 54.096 117 Pentachlorophenol 266 12.647 (0.981) 58168 50.0000 54.096 117 Pentachlorophenol 266 12.647 (0.981) 58168 50.0000 51.481 123 Anthracene 178 13.058 13.058 (1.012) 575232 50.0000 47.657 135 Pluoranthene 202 15.889 15.889 (1.012) 575232 50.0000 47.657 135 Pluoranthene 202 15.889 15.889 (1.012) 575232 50.0000 49.782 136 Benzidine 184 16.317 16.317 (0.835) 363229 50.0000 62.577 137 Pyrene 202 16.434 16.434 (0.841) 606633 50.0000 45.746 144 Butylbenzylphthalate 149 18.379 18.379 (0.941) 324288 50.0000 45.522 149 3,3"-Dichlorobenzidine 252 19.543 19.543 (1.001) 253202 50.0000 53.964	90	Dibenzofuran	168	9 997	9 997	(1.042)	536417	50 0000	50 541
95 2,3,5,6-Tetrachlorophenol 232 10 280 10.280 (1.071) 83428 50.0000 47.211 92 2,3,4,6-Tetrachlorophenol 232 10 376 10.376 (1.081) 85197 50.0000 48.158 96 2-Naphthylamine 143 10.344 10.344 (1.078) 313265 50.0000 41 084 97 Diethylphthalate 149 10.708 10.708 (1.116) 389505 50.0000 44 390 98 Fluorene 166 10.745 10.745 (1.120) 422350 50.0000 48.604 99 4-Chlorophenyl-phenylether 204 10.782 10 782 (1.124) 195077 50.0000 50 194 100 4-Nitroaniline 138 10.900 10 900 (1.136) 104552 50.0000 40.918 102 4,6-Dinitro-2-methylphenol 198 11 007 11.007 (0.853) 55193 50 0000 43.929 103 N-Nitrosodiphenylamine (1) 169 11.071 11 071 (0.858) 290432 50.0000 51.721 104 1,2-Diphenylhydrazine 77 11 135 11 135 (0.863) 481308 50.0000 50 942 112 4-Bromophenyl-phenylether 248 11.888 11.888 (0.922) 103801 50.0000 54.099 113 Hexachlorobenzene 284 12 193 12 193 (0.945) 109639 50.0000 54.096 117 Pentachlorophenol 266 12 647 (2.647 (0.981) 58168 50.0000 54.096 117 Pentachlorophenol 266 12 647 (0.981) 58168 50.0000 54.096 117 Pentachlorophenol 266 12 647 (0.981) 58168 50.0000 50.935 126 Carbazole 167 13.496 (1.046) 534645 50.0000 48.761 130 Di-n-Butylphthalate 149 14 639 14.639 (1.012) 575232 50.0000 47.657 137 Pyrene 202 16.434 16.434 (0.841) 606633 50.0000 47.657 137 Pyrene 202 16.434 16.434 (0.841) 606633 50.0000 45.522 149 3,3'-Dichlorobenzidine 252 19.543 19.543 (1.001) 253202 50.0000 53.964	91	2,4-Dinitrotoluene	165	10 115	10.115	(1.054)	115637		
92 2,3,4,6-Tetrachlorophenol 232 10 376 10.376 (1.081) 85197 50.0000 48.158 96 2-Naphthylamine 143 10.344 10.344 (1.078) 313265 50.0000 41 084 97 Diethylphthalate 149 10.708 10.708 (1.116) 389505 50.0000 44 390 98 Fluorene 166 10.745 10.745 (1.120) 422350 50.0000 48.604 99 4-Chlorophenyl-phenylether 204 10.782 10.782 (1.124) 195077 50.0000 50 194 100 4-Mitroaniline 138 10.900 10 900 (1.136) 104552 50.0000 40.918 102 4.6-Dinitro-2-methylphenol 198 11 007 11.007 (0.853) 55193 50 0000 43.929 103 N-Nitrosodiphenylamine (1) 169 11.071 11.071 (0.858) 290432 50.0000 51.721 104 1,2-Diphenylhydrazine 77 11 135 11 135 (0.863) 481308 50.0000 50.942 112 4-Bromophenyl-phenylether 248 11.888 11.888 (0.922) 103801 50.0000 54.909 113 Hexachlorophenol 266 12 647 (1.981) 58168 50.0000 54.909 117 Pentachlorophenol 266 12 647 (1.981) 58168 50.0000 49 783 (M) 122 Phenanthrene 178 12.951 (1.004) 565778 50.0000 51.481 123 Anthracene 178 13.058 13.058 (1.012) 575232 50.0000 50.935 126 Carbazole 167 13.496 13.496 (1.046) 534645 50.0000 49.782 136 Benzidine 184 16.317 16.317 (0.835) 363229 50.0000 62.577 137 Pyrene 202 16.434 16.434 (0.841) 606633 50.0000 49.782 149 3,3'-Dichlorobenzidine 252 19.543 19.543 (1.001) 253202 50.0000 53.964	95	2,3,5,6-Tetrachlorophenol	232	10 280	10.280	(1.071)	83428	50.0000	
96 2-Naphthylamine	92	2,3,4,6-Tetrachlorophenol	232	10 376	10.376	(1.081)	85197		
97 Diethylphthalate 149 10.708 10.708 (1 116) 389505 50.0000 44 390 98 Fluorene 166 10.745 10.745 (1.120) 422350 50.0000 48.604 99 4-Chlorophenyl-phenylether 204 10.782 10 782 (1.124) 195077 50.0000 50 194 100 4-Nitroaniline 138 10.900 10 900 (1.136) 104552 50.0000 40.918 102 4,6-Dinitro-2-methylphenol 198 11 007 11.007 (0 853) 55193 50 0000 43.929 103 N-Nitrosodiphenylamine (1) 169 11.071 11 071 (0 858) 290432 50.0000 51.721 104 1,2-Diphenylhydrazine 77 11 135 11 135 (0 863) 481308 50.0000 50 942 112 4-Bromophenyl-phenylether 248 11.888 11 888 (0.922) 103801 50.0000 54.909 113 Hexachlorobenzene 284 12 193 12 193 (0 945) 109639 50.0000 54.096 117 Pentachlorophenol 266 12 647 12 647 (0.981) 58168 50.0000 49 783 (M) 122 Phenanthrene 178 12.951 12.951 (1 004) 565778 50.0000 51.481 123 Anthracene 178 13.058 13.058 (1.012) 575232 50 0000 50.935 126 Carbazole 167 13.496 13.496 (1 046) 534645 50.0000 49.782 130 Di-n-Butylphthalate 149 14 639 14.639 (1 135) 638330 50.0000 47.657 135 Pluoranthene 202 15 889 15.889 (1 232) 582296 50 0000 49.782 136 Benzidine 184 16.317 16.317 (0 835) 363229 50.0000 42.577 137 Pyrene 202 16.434 16.434 (0 841) 606633 50.0000 45.746 144 Butylbenzylphthalate 149 18 379 18.379 (0 941) 324288 50.0000 45.522 149 3,3'-Dichlorobenzidine 252 19.543 19.543 (1 001) 253202 50.0000 53.964	96	2-Naphthylamine	143	10.344	10.344	(1.078)	313265	50.0000	
98 Fluorene 166 10.745 10.745 (1.120) 422350 50.0000 48.604 99 4-Chlorophenyl-phenylether 204 10.782 10 782 (1.124) 195077 50.0000 50 194 100 4-Nitroaniline 138 10.900 10 900 (1.136) 104552 50.0000 40.918 102 4,6-Dinitro-2-methylphenol 198 11 007 11.007 (0 853) 55193 50 0000 43.929 103 N-Nitrosodiphenylamine (1) 169 11.071 11 071 (0 858) 290432 50.0000 51.721 104 1,2-Diphenylhydrazine 77 11 135 11 135 (0 863) 481308 50.0000 50 942 112 4-Bromophenyl-phenylether 248 11.888 11 888 (0.922) 103801 50.0000 54.996 113 Hexachlorobenzene 284 12 193 12 193 (0 945) 109639 50.0000 54.096 117 Pentachlorophenol 266 12 647 12 647 (0.981) 58168 50.0000 49 783 (M) 122 Phenanthrene 178 12.951 (1 004) 565778 50.0000 51.481 123 Anthracene 178 13.058 13.058 (1.012) 575232 50 0000 50.935 126 Carbazole 167 13.496 13.496 (1 046) 534645 50.0000 49.782 136 Benzidine 184 16.317 16.317 (0 835) 363229 50.0000 62.577 137 Pyrene 202 16.434 16.434 (0 841) 606633 50.0000 45.746 144 Butylbenzylphthalate 149 18 379 18.379 (0 941) 324288 50.0000 45.522 149 3,3'-Dichlorobenzidine 252 19.543 19.543 (1 001) 253202 50.0000 53.964	97	Diethylphthalate	149	10.708	10.708	(1 116)	389505		
99 4-Chlorophenyl-phenylether 204 10.782 10 782 (1.124) 195077 50.0000 50 194 100 4-Nitroaniline 138 10.900 10 900 (1.136) 104552 50.0000 40.918 102 4.6-Dinitro-2-methylphenol 198 11 007 11.007 (0 853) 55193 50 0000 43.929 103 N-Nitrosodiphenylamine (1) 169 11.071 11 071 (0 858) 290432 50.0000 51.721 104 1,2-Diphenylhydrazine 77 11 135 11 135 (0 863) 481308 50.0000 50 942 112 4-Bromophenyl-phenylether 248 11.888 11 888 (0.922) 103801 50.0000 54.909 113 Hexachlorophenol 266 12 647 12 647 (0.981) 58168 50.0000 54.096 117 Pentachlorophenol 266 12 647 12 647 (0.981) 58168 50.0000 49 783 (M) 122 Phenanthrene 178 12.951 12.951 (1 004) 565778 50.0000 51.481 123 Anthracene 178 13.058 13.058 (1.012) 575232 50 0000 50.935 126 Carbazole 167 13.496 13.496 (1 046) 534645 50.0000 48.761 130 Di-n-Butylphthalate 149 14 639 14.639 (1 135) 638330 50.0000 47.657 135 Pluoranthene 202 15 889 15.889 (1 232) 582296 50 0000 49.782 136 Benzidine 184 16.317 16.317 (0 835) 363229 50.0000 42.776 137 Pyrene 202 16.434 16.434 (0 841) 606633 50.0000 45.746 144 Butylbenzylphthalate 149 18 379 18.379 (0 941) 324288 50.0000 45.522 149 3.37-Dichlorobenzidine 252 19.543 19.543 (1 001) 253202 50.0000 53.964	98	Fluorene	166	10.745	10.745	(1.120)	422350		
100 4-Nitroaniline	99	4-Chlorophenyl-phenylether	204	10.782	10 782	(1.124)	195077	50.0000	
102 4,6-Dinitro-2-methylphenol 198 11 007 11.007 (0 853) 55193 50 0000 43.929  103 N-Nitrosodiphenylamine (1) 169 11.071 11 071 (0 858) 290432 50.0000 51.721  104 1,2-Diphenylhydrazine 77 11 135 11 135 (0 863) 481308 50.0000 50 942  112 4-Bromophenyl-phenylether 248 11.888 11 888 (0.922) 103801 50.0000 54.909  113 Hexachlorobenzene 284 12 193 12 193 (0 945) 109639 50.0000 54.096  117 Pentachlorophenol 266 12 647 12 647 (0.981) 58168 50.0000 49 783 (M)  122 Phenanthrene 178 12.951 12.951 (1 004) 565778 50.0000 51.481  123 Anthracene 178 13.058 13.058 (1.012) 575232 50 0000 50.935  126 Carbazole 167 13.496 13.496 (1 046) 534645 50.0000 48.761  130 Di-n-Butylphthalate 149 14 639 14.639 (1 135) 638330 50.0000 47.657  135 Fluoranthene 202 15 889 15.889 (1 232) 582296 50 0000 49.782  136 Benzidine 184 16.317 16.317 (0 835) 363229 50.0000 62.577  137 Pyrene 202 16.434 16.434 (0 841) 606633 50.0000 45.746  149 3,3'-Dichlorobenzidine 252 19.543 19.543 (1 001) 253202 50.0000 53.964	100	4-Natroanilane	138	10.900	10 900	(1.136)			
103 N-Nitrosodiphenylamine (1) 169 11.071 11 071 (0 858) 290432 50.0000 51.721 104 1,2-Diphenylhydrazine 77 11 135 11 135 (0 863) 481308 50.0000 50 942 112 4-Bromophenyl-phenylether 248 11.888 11 888 (0.922) 103801 50.0000 54.909 113 Hexachlorobenzene 284 12 193 12 193 (0 945) 109639 50.0000 54.096 117 Pentachlorophenol 266 12 647 12 647 (0.981) 58168 50.0000 49 783 (M) 122 Phenanthrene 178 12.951 12.951 (1 004) 565778 50.0000 51.481 123 Anthracene 178 13.058 13.058 (1.012) 575232 50 0000 50.935 126 Carbazole 167 23.496 13.496 (1 046) 534645 50.0000 48.761 130 Di-n-Butylphthalate 149 14 639 14.639 (1 135) 638330 50.0000 47.657 135 Fluoranthene 202 15 889 15.889 (1 232) 582296 50 0000 49.782 136 Benzidine 184 16.317 16.317 (0 835) 363229 50.0000 62.577 137 Pyrene 202 16.434 16.434 (0 841) 606633 50.0000 45.746 144 Butylbenzylphthalate 149 18 379 18.379 (0 941) 324288 50.0000 53.964	102	4,6-Dinitro-2-methylphenol	198	11 007	11.007	(0 853)	55193		
104 1,2-Diphenylhydrazine 77 11 135 11 135 (0 863) 481308 50.0000 50 942  112 4-Bromophenyl-phenylether 248 11.888 11 888 (0.922) 103801 50.0000 54.909  113 Hexachlorobenzene 284 12 193 12 193 (0 945) 109639 50.0000 54.096  117 Pentachlorophenol 266 12 647 12 647 (0.981) 58168 50.0000 49 783 (M)  122 Phenanthrene 178 12.951 12.951 (1 004) 565778 50.0000 51.481  123 Anthracene 178 13.058 13.058 (1.012) 575232 50 0000 50.935  126 Carbazole 167 13.496 13.496 (1 046) 534645 50.0000 48.761  130 Di-n-Butylphthalate 149 14 639 14.639 (1 135) 638330 50.0000 47.657  135 Fluoranthene 202 15 889 15.889 (1 232) 582296 50 0000 49.782  136 Benzidine 184 16.317 16.317 (0 835) 363229 50.0000 62.577  137 Pyrene 202 16.434 16.434 (0 841) 606633 50.0000 45.746  149 3.3'-Dichlorobenzidine 252 19.543 19.543 (1 001) 253202 50.0000 53.964	103	N-Nitrosodiphenylamine (1)	169	11.071	11 071	(0 858)	290432	50.0000	
112 4-Bromophenyl-phenylether       248       11.888       11.888       0.922       103801       50.0000       54.909         113 Hexachlorobenzene       284       12 193       12 193       0.945       109639       50.0000       54.096         117 Pentachlorophenol       266       12 647       12 647       0.981       58168       50.0000       49 783 (M)         122 Phenanthrene       178       12.951       12.951       1 004       565778       50.0000       51.481         123 Anthracene       178       13.058       13.058       1.012       575232       50.0000       50.935         126 Carbazole       167       13.496       13.496       (1.012)       575232       50.0000       48.761         130 Di-n-Butylphthalate       149       14 639       14.639       (1.135)       638330       50.0000       47.657         135 Fluoranthene       202       15 889       15.889       (1.232)       582296       50.0000       49.782         136 Benzidine       184       16.317       16.317       (0.835)       363229       50.0000       45.746         144 Butylbenzylphthalate       149       18 379       18.379       (0.941)       324288       50.0000	104	1,2-Diphenylhydrazine	77	11 135	11 135	(0 863)	481308	50.0000	
117 Pentachlorophenol 266 12 647 12 647 (0.981) 58168 50.0000 49 783 (M)  122 Phenanthrene 178 12.951 12.951 (1.004) 565778 50.0000 51.481  123 Anthracene 178 13.058 13.058 (1.012) 575232 50 0000 50.935  126 Carbazole 167 13.496 13.496 (1.046) 534645 50.0000 48.761  130 Di-n-Butylphthalate 149 14 639 14.639 (1.135) 638330 50.0000 47.657  135 Pluoranthene 202 15 889 15.889 (1.232) 582296 50 0000 49.782  136 Benzidine 184 16.317 16.317 (0.835) 363229 50.0000 62.577  137 Pyrene 202 16.434 16.434 (0.841) 606633 50.0000 45.746  144 Butylbenzylphthalate 149 18 379 18.379 (0.941) 324288 50.0000 45.522  149 3,3'-Dichlorobenzidine 252 19.543 19.543 (1.001) 253202 50.0000 53.964	112	4-Bromophenyl-phenylether	248	11.888	11 888	(0,922)	103801	50.0000	
122 Phenanthrene       178       12.951       12.951       (1 004)       565778       50.0000       51.481         123 Anthracene       178       13.058       13.058       (1.012)       575232       50 0000       50.935         126 Carbazole       167       13.496       13.496       (1 046)       534645       50.0000       48.761         130 Di-n-Butylphthalate       149       14 639       14.639       (1 135)       638330       50.0000       47.657         135 Pluoranthene       202       15 889       15.889       (1 232)       582296       50 0000       49.782         136 Benzidine       184       16.317       16.317       (0 835)       363229       50.0000       62.577         137 Pyrene       202       16.434       16.434       (0 841)       606633       50.0000       45.746         144 Butylbenzylphthalate       149       18 379       18.379       (0 941)       324288       50.0000       45.522         149 3.3'-Dichlorobenzidine       252       19.543       (1 001)       253202       50.0000       53.964	113	Hexachlorobenzene	284	12 193	12 193	(0 945)	109639	50.0000	
122 Phenanthrene       178       12.951       12.951       (1 004)       565778       50.0000       51.481         123 Anthracene       178       13.058       13.058       (1.012)       575232       50 0000       50.935         126 Carbazole       167       13.496       13.496       (1 046)       534645       50.0000       48.761         130 Di-n-Butylphthalate       149       14 639       14.639       (1 135)       638330       50.0000       47.657         135 Pluoranthene       202       15 889       15.889       (1 232)       582296       50 0000       49.782         136 Benzidine       184       16.317       16.317       (0 835)       363229       50.0000       62.577         137 Pyrene       202       16.434       16.434       (0 841)       606633       50.0000       45.746         144 Butylbenzylphthalate       149       18 379       18.379       (0 941)       324288       50.0000       45.522         149 3.3'-Dichlorobenzidine       252       19.543       (1 001)       253202       50.0000       53.964	117	Pentachlorophenol	266	12 647	12 647	(0.981)	58168	50.0000	49 783 (M)
123 Anthracene       178       13.058       13.058       13.058       1.012)       575232       50 0000       50.935         126 Carbazole       167       13.496       13.496       10.46)       534645       50.0000       48.761         130 Di-n-Butylphthalate       149       14 639       14.639       1 135)       638330       50.0000       47.657         135 Fluoranthene       202       15 889       15.889       1 232)       582296       50.0000       49.782         136 Benzidine       184       16.317       16.317       (0 835)       363229       50.0000       62.577         137 Pyrene       202       16.434       16.434       (0 841)       606633       50.0000       45.746         144 Butylbenzylphthalate       149       18 379       18.379       (0 941)       324288       50.0000       45.522         149 3.3'-Dichlorobenzidine       252       19.543       19.543       (1 001)       253202       50.0000       53.964	122	Phenanthrene	178						
126 Carbazole       167       13.496       13.496       (1 046)       534645       50.0000       48.761         130 Di-n-Butylphthalate       149       14 639       14.639       (1 135)       638330       50.0000       47.657         135 Fluoranthene       202       15 889       15.889       (1 232)       582296       50 0000       49.782         136 Benzidine       184       16.317       16.317       (0 835)       363229       50.0000       62.577         137 Pyrene       202       16.434       16.434       (0 841)       606633       50.0000       45.746         144 Butylbenzylphthalate       149       18 379       18.379       (0 941)       324288       50.0000       45.522         149 3,3'-Dichlorobenzidine       252       19.543       19.543       (1 001)       253202       50.0000       53.964	123	Anthracene	178	13.058	13.058	(1.012)	575232	50 0000	
130 Di-n-Butylphthalate       149       14 639 14.639 (1 135)       638330 50.0000 47.657         135 Fluoranthene       202       15 889 15.889 (1 232)       582296 50 0000 49.782         136 Benzidine       184       16.317 16.317 (0 835)       363229 50.0000 62.577         137 Pyrene       202       16.434 16.434 (0 841) 606633 50.0000 45.746         144 Butylbenzylphthalate       149       18 379 18.379 (0 941) 324288 50.0000 45.522         149 3,3'-Dichlorobenzidine       252       19.543 19.543 (1 001) 253202 50.0000 53.964	126	Carbazole	167	13.496	13.496	(1 046)	534645		
135 Pluoranthene       202       15 889 15.889 (1 232)       582296 50 0000 49.782         136 Benzidine       184 16.317 16.317 (0 835)       363229 50.0000 62.577         137 Pyrene       202 16.434 16.434 (0 841) 606633 50.0000 45.746         144 Butylbenzylphthalate       149 18 379 18.379 (0 941) 324288 50.0000 45.522         149 3,3'-Dichlorobenzidine       252 19.543 19.543 (1 001) 253202 50.0000 53.964	130	Di-n-Butylphthalate	149						
136 Benzidine     184     16.317 16.317 (0 835)     363229     50.0000     62.577       137 Pyrene     202     16.434 16.434 (0 841)     606633     50.0000     45.746       144 Butylbenzylphthalate     149     18 379 18.379 (0 941)     324288     50.0000     45.522       149 3,3'-Dichlorobenzidine     252     19.543 19.543 (1 001)     253202     50.0000     53.964	135	Fluoranthene	202						
137 Pyrene 202 16.434 16.434 (0 841) 606633 50.0000 45.746 144 Butylbenzylphthalate 149 18 379 18.379 (0 941) 324288 50.0000 45.522 149 3,3'-Dichlorobenzidine 252 19.543 19.543 (1 001) 253202 50.0000 53.964	136	Benzidine	184						
144 Butylbenzylphthalate     149     18 379 18.379 (0 941)     324288 50.0000 45.522       149 3,3'-Dichlorobenzidine     252     19.543 19.543 (1 001)     253202 50.0000 53.964	137	Pyrene	202						
149 3,3'-Dichlorobenzidine 252 19.543 19.543 (1 001) 253202 50.0000 53.964	144	Butylbenzylphthalate	149						
and the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of the same of th	149	3,3'-Dichlorobenzidine	252	19.543	19.543	(1 001)		50.0000	
· · · · · · · · · · · · · · · · · · ·	150	Benzo (a) Anthracene	228	19 490	19 490	(0 998)	643559		51.051

Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522CCC.D Report Date: 22-May-2000 12:04

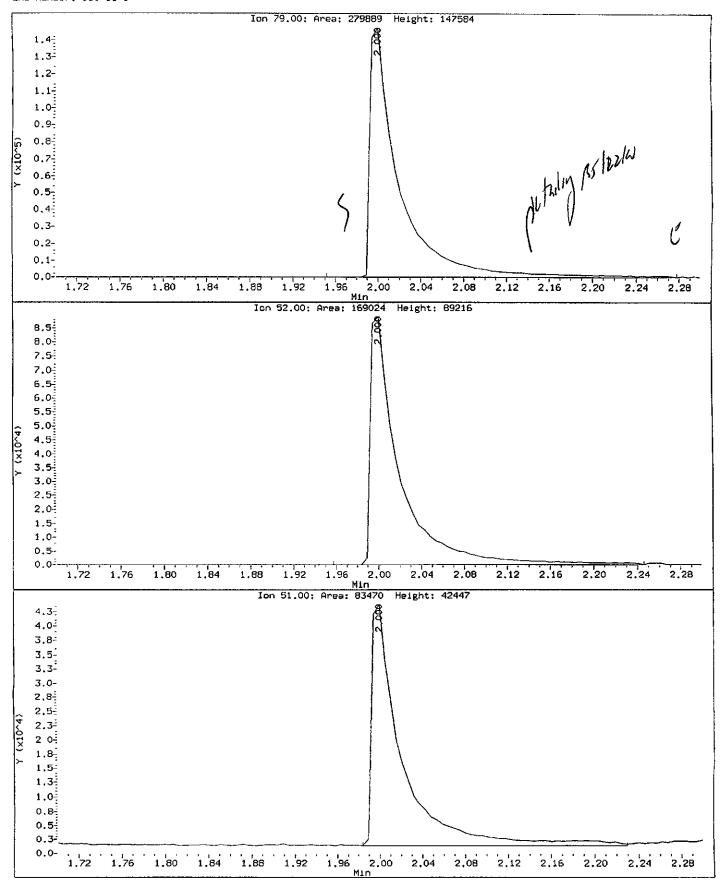
						MUOMA	TS
	QUANT SIG					CAL-AMT	ON-COL
Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	( ng)	( ng)
	5. <b>2</b> 0. \$		*****	8E===0		======	00=2055
151 Chrysene	228	19 597	19.597	(1 003)	620609	50 0000	51 161
153 bis(2-ethylhexyl)Phthalate	149	19 949	19.949	(1 021)	465605	50 0000	46.570
155 Di-n-octylphthalate	149	21 419	21.419	(0.936)	912315	50 0000	47.230
157 Benzo(b) fluoranthene	252	22.076	22.076	(0 964)	810619	50 0000	50 972
158 Benzo(k)fluoranthene	252	22 134	22 134	(0.967)	825097	50.0000	50.602
159 7,12-dimethylbenz(a)anthracen	256	22.134	22,134	(0.967)	332481	50.0000	48 544
167 Benzo(a)pyrene	252	22.770	22 770	(0.995)	798737	50 0000	51 911
169 Indeno(1,2,3-cd)pyrene	276	25 158	25 158	(1.099)	1132982	50.0000	53.436
170 Dibenz(a,h)anthracene	278	25 206	25.206	(1 101)	971394	50.0000	53 685
171 Benzo(g,h,1)perylene	276	25 746	25.746	(1 125)	1003463	50.0000	53 160
\$ 172 Nitrobenzene-d5	82	5 686	5.686	(0.866)	263493	50.0000	48.696
\$ 173 2-Fluorobiphenyl	172	8 309	B 309	(0.866)	443954	50.0000	55 215
\$ 174 Terphenyl-d14	244	16 969	16 969	(0 869)	467386	50.0000	46 786
\$ 175 Phenol-d5	99	4.719		(0 938)	306982	50.0000	47.290
\$ 176 2-Fluorophenol	112	3 709	3.709	(0 738)	246558	50 0000	50.758
\$ 177 2,4,6-Tribromophenol	330	11 338		(0.879)	53349	50 0000	50.300
\$ 178 2-Chlorophenol-d4	132	4.831		(0 961)	224630	50.0000	49 394
\$ 179 1,2-Dichlorobenzene-d4	152	5 237		(1.041)	164052	50.0000	51.344

## QC Flag Legend

M - Compound response manually integrated.

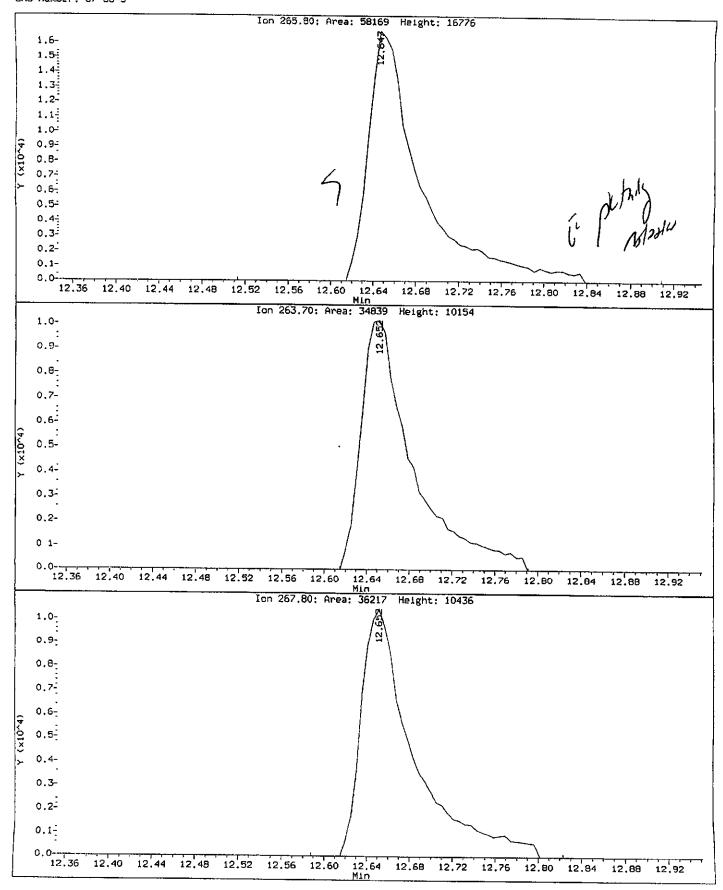
Data File: \\QPITPAOZ\D\chem\71.1\s052200.b\S0522CCC.D Injection Date: 22-MAY-2000 11:27 Instrument: 71.1 Client Sample ID: sstd050

Compound: Pyridine CAS Number: 110-86-1



Data File: \\QPITPAO2\D\chem\71.i\s052200.b\S0522CCC.D Injection Date: 22-MAY-2000 11:27 Instrument: 71.i Client Sample ID: sstd050

Compound: Pentachlorophenol CAS Number: 87-86-5



GC/MS SEMIVOLATILE QC DATA

Data File: \\OP!TPA02\D\chem\71.1\s051900.b\S0519DF2.D

Date : 19-MA/-2000 11:44

Client ID: DFTPP02

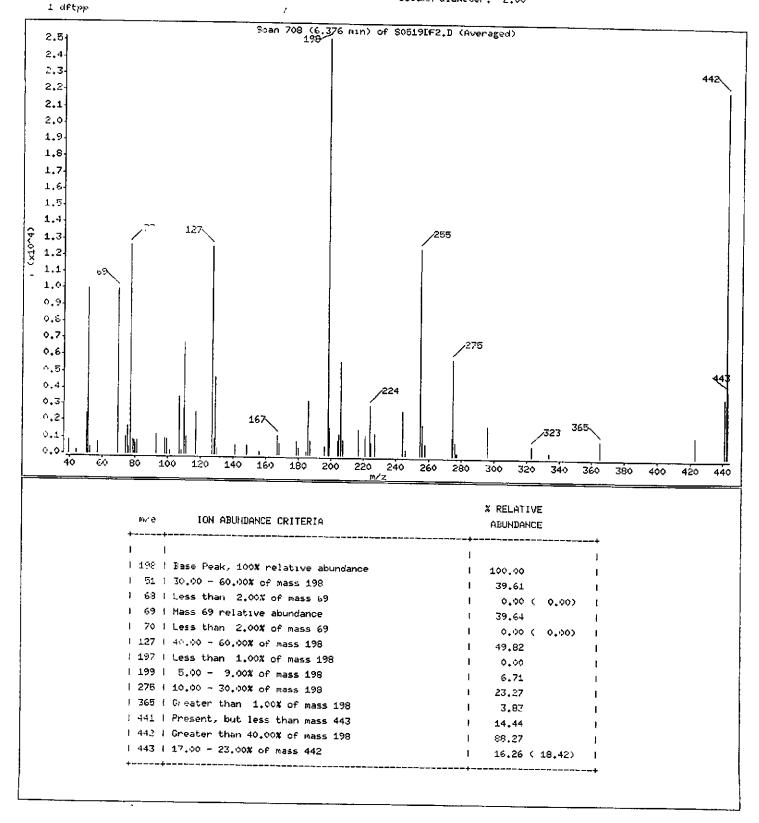
Instrument: 71.1

Sample Info: DFTPF050 (25) ph) 194-158-6

Operator: SAB: 2ul ind (100ul+ful) int

Column diameter: 2.00

Column phase:



Data File: \\QPITPA02\D\chem\71.1\s051900.b\\$0519DF2.D

Date : 19-MAY-2000 11:44

Client ID: DFTPP02

Instrument: 71.1

Sample Info: DFTPFv50 (25ppb) 194-158-6

Operator: SAB: 2ul inj (100ul+1ul) int

Column phase:

Column diameter: 2.00

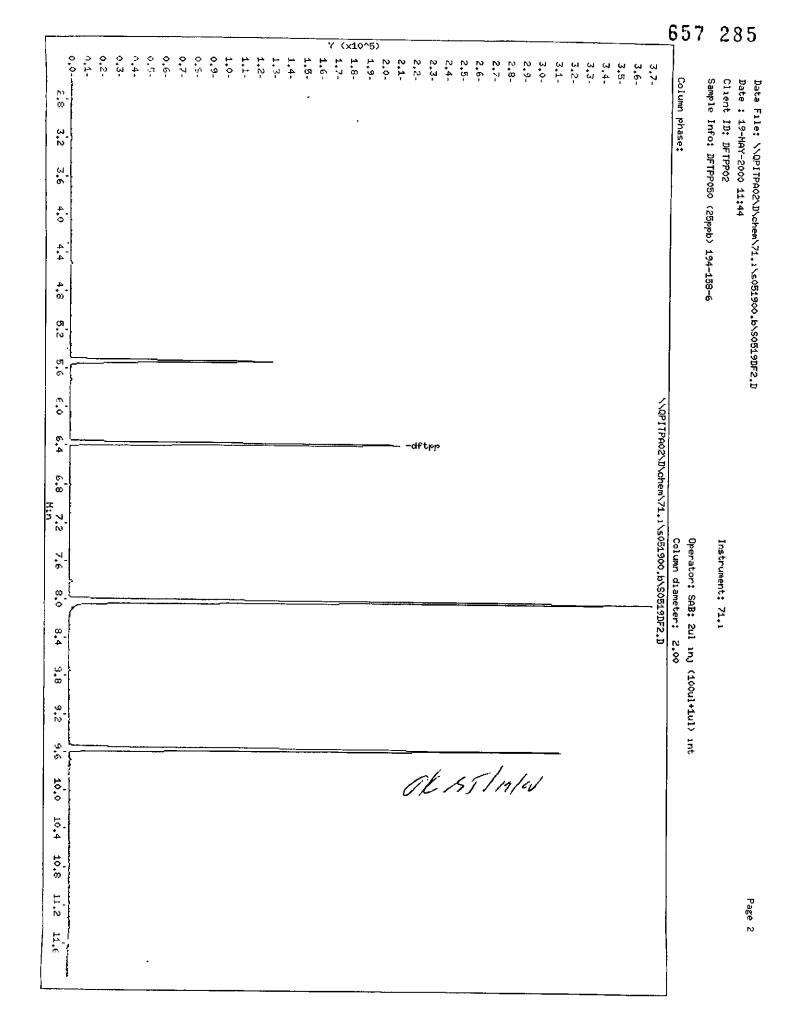
Data File: S0519DF2.D

Spectrum: Scan 709 (6.376 min) of SO519DF2.D (Averaged)

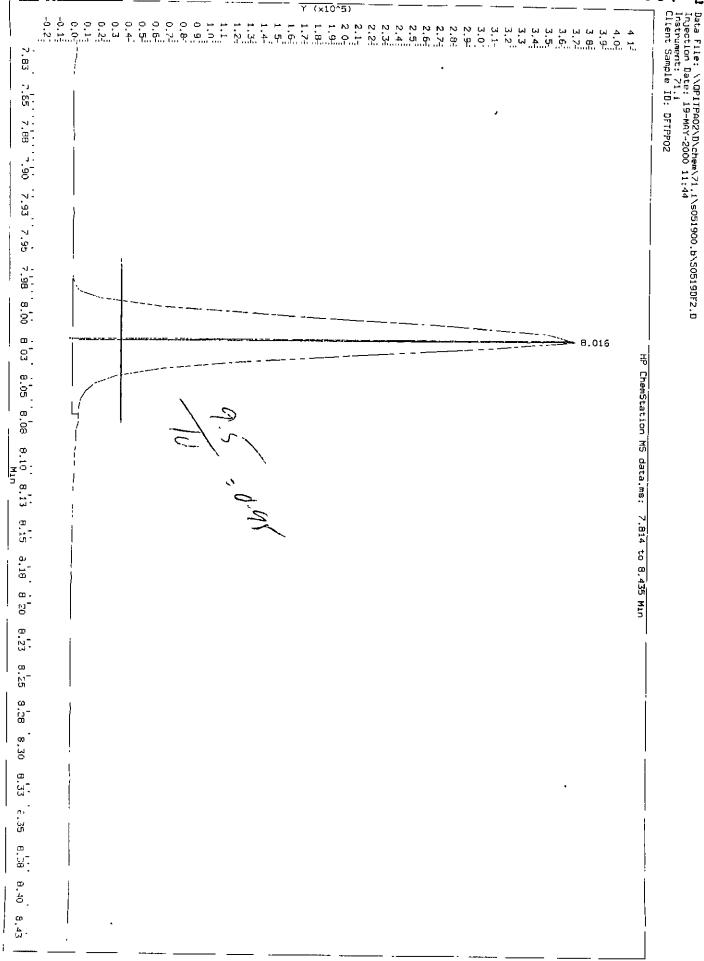
Location of Maximum: 198,00 Number of points: 67

I	m/z	Y <del></del>	MZ	Υ.	m/z	Y 	m/z 
347	246.00		180.00		99,00		39,00
12521	255.00	189 I	185.00	193 (	101,00	167	44.00
1846	256,00	3323	186.00	3462 I	107.00	2417	50.00
676	258.00	854	187.00	188 I	108.00	9981 I	51,00
	274.00	512   		6758 1	-	400	52.00
	275.00		198.00		111.00		57.00
799	276.00	1692	199,00	2547 (	117.00	9989	69,00
178	277,00	890 I	204,00	12555	127,00	1019	74,00
1789	296,00	1258	205.00	899 1	128,00	1653 I	75.00
<u>6</u> £4	323.00		206.00		129.00		76.00
208		917 I			130.00		77.00
966	365,00	1552 I	217,00	577 1	141,00	876 I	78.00
1286	423,00	1189	221.00	596	148.00	793 !	79.00
3638	441.00	3011 I	224,00	190 (	156.00	620 1	80,00
22245		779 J			•	83 <b>0</b>	
	443,00			690 I			93.00
		2715 I	244.00	836	179,00	922 i	98.00

2078



Data File: \\QPITPA02\D\chem\71.1\s051900.b\S0519DF2.D Injection Date: 19-MAY-2000 11:44 Instrument: 71.1 Client Sample ID: DFTPP02



Date : 22-MAY-2000 11:09

Client ID: DFTPP02

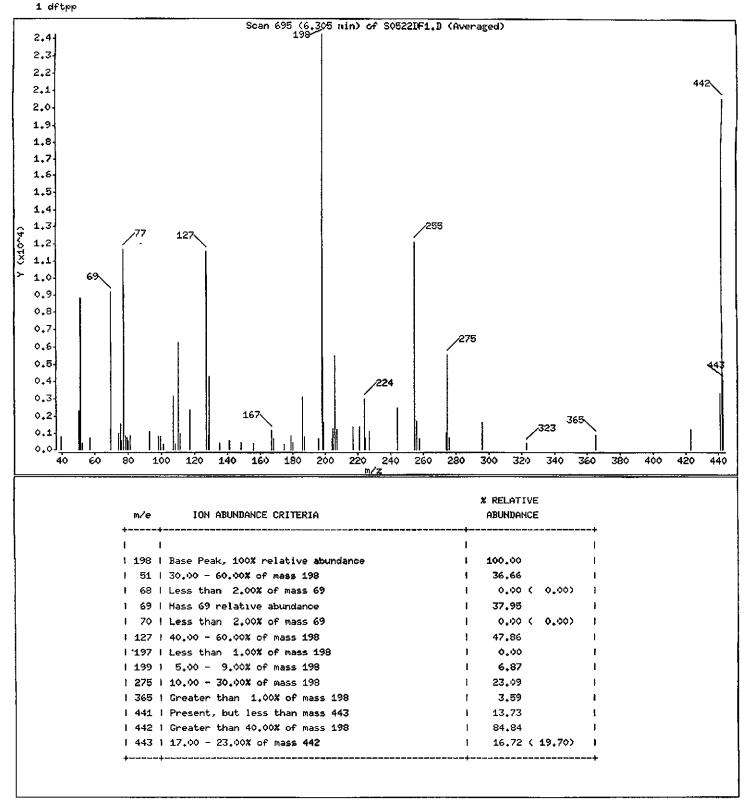
Instrument: 71.1

Sample Info: DFTPP050 (25ppb) 194-158-6

Operator: SAB: 2ul inj (100ul+1ul) int

Column phase:

Column diameter: 2.00



Data File: \\QPITPAO2\D\chem\71.1\s052200.b\S0522DF1.D

Date : 22-MAY-2000 11:09

Client ID: DFTPP02

Instrument: 71.1

Sample Info: DFTPP050 (25ppb) 194-158-6

Operator: SAB: 2ul inj (100ul+1ul) int

Column phase:

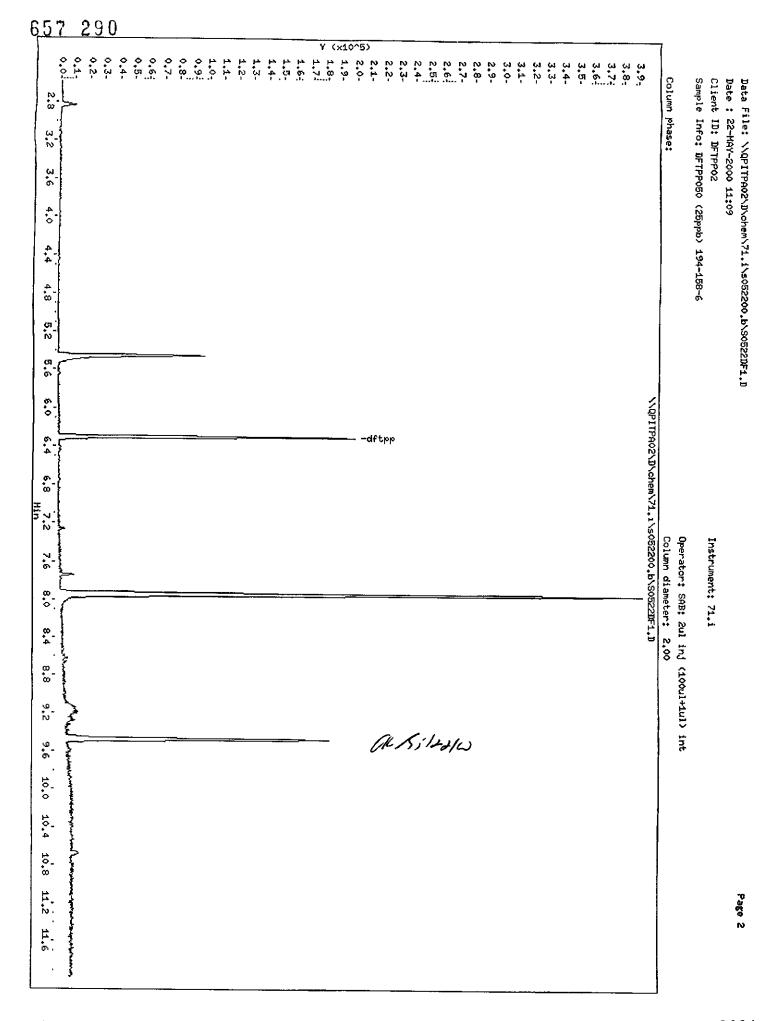
Column diameter: 2.00

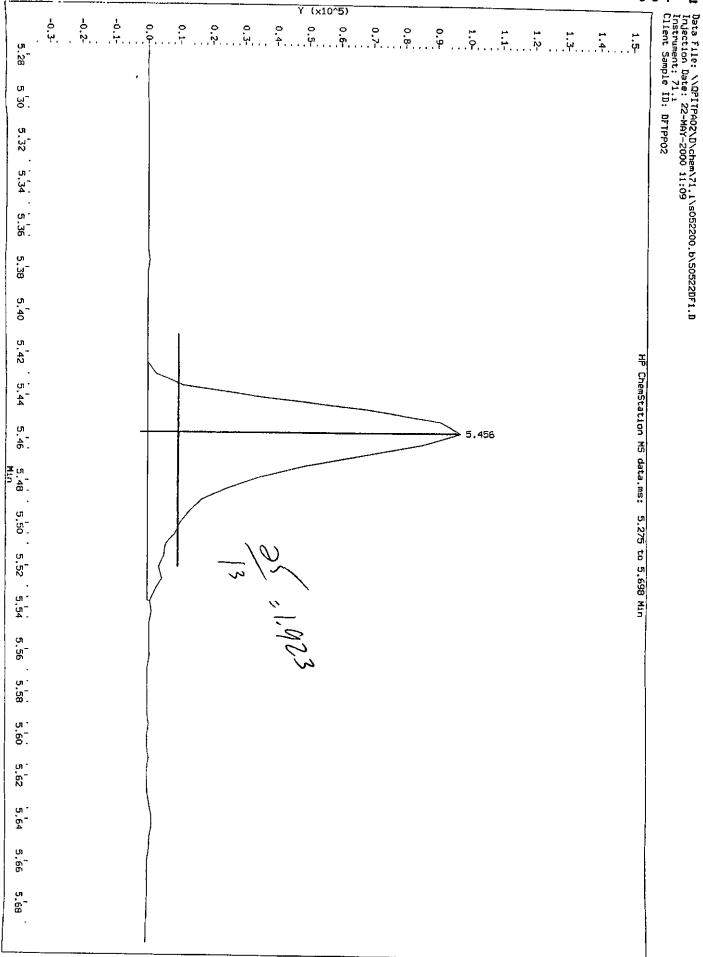
Data File: S0522DF1.D

Spectrum: Scan 696 (6,305 min) of S0522DF1.D (Averaged)

Location of Maximum: 198.00 Number of points: 63

Y	m/z	Y	m/z	Υ .	m/z		Y	m/z	_
1125	227,00	342 I	175.00				796	39,00	
2523		828	179.00	336	101.00	ţ	2334	50,00	
12128	255.00	433 1	189,00	3152 i	107,00	t	8874	51,00	
1738		3092	186,00	350	108.00	1	383	52,00	
689	258.00		187,00		110.00			57,00	
1035	274.00	678 l	196.00		111.00			69.00	•
5590	•	24208	198.00		117,00	ı	1006	74,00	
707	276.00		199.00	11585 I	127,00	ŀ	1577	75.00	
1680	296.00	692 I	204.00	892	128,00	ŀ	579	76.00	
400	323.00	1270 I	205.00		129.00			77.00	
<del></del> 868	365.00	5515 I	206.00		135,00			78,00	
1223	423.00	1250	207.00	545	141.00	ı	735	79,00	
3324	441.00		217.00	455	148,00	ı	584	80,00	
20538	442.00	1365 1	221.00	377	156,00	ı	917	81,00	
4047	443.00		224.00		167.00			93.00	
		739 1	225,00	676 1	168.00	1	845	98,00	





Data File: \\QPITPAOZ\D\chem\71.i\s052200.b\S0522DF1.D Injection Date: 22-MAY-2000 11:09 Instrument: 71.i Client Sample ID: DFTPF02

## UXB INTERNATIONAL METHOD BLANK COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E190000 269

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DDF2D101 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

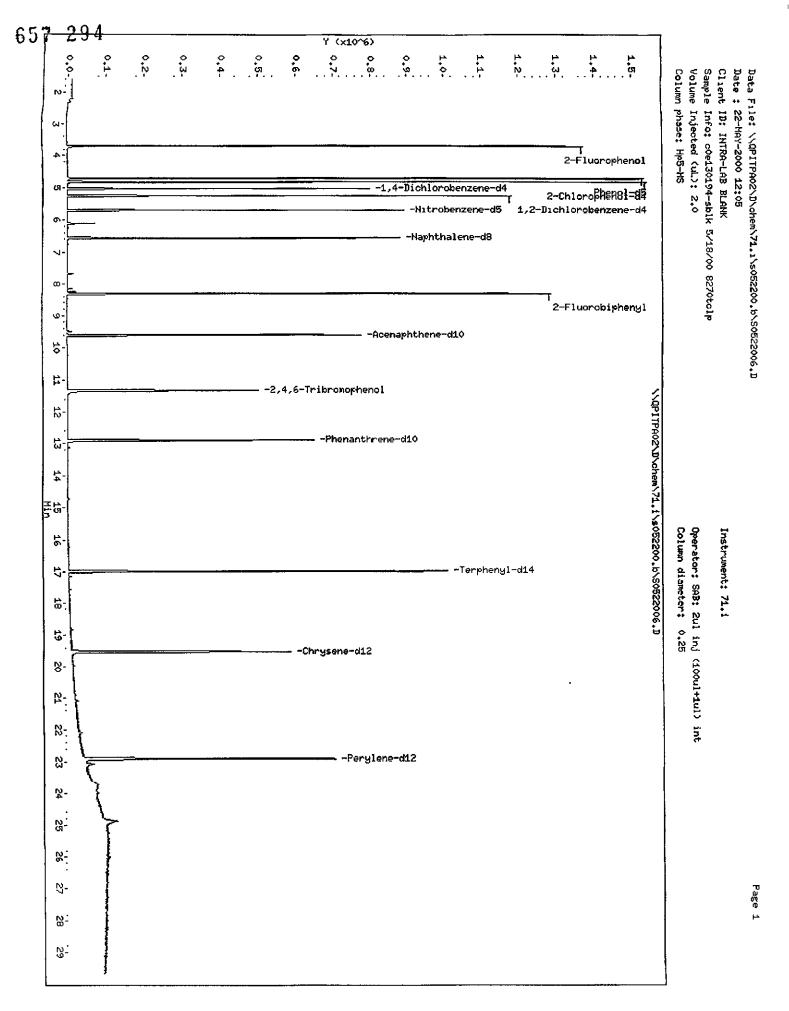
Moisture %:NA

QC Batch: 0140269

Client Sample Id: INTRA-LAB BLANK

#### CONCENTRATION UNITS:

CONCENTIALL ZON CHILD.					
COMPOUND (ug/L	or ug/kg) mg/L	Q			
1,4-Dichlorobenzene	0.050	ן ט			
2,4-Dinitrotoluene	0.050	ן ט			
Hexachlorobenzene	0.050	ן ט			
Hexachlorobutadiene	0.050	ט			
Hexachloroethane	0.050	ן ט			
Nitrobenzene	0.050	ן די			
Pentachlorophenol	0.25	ן ט			
Pyridine	0.10	U			
2,4,5-Trichlorophenol	0.050	ן ט			
2,4,6-Trichlorophenol	0.050	<u>  U</u>			
Cresols (total)	0.050	<u>                                     </u>			
	1,4-Dichlorobenzene 2,4-Dinitrotoluene Hexachlorobenzene Hexachlorobutadiene Hexachloroethane Nitrobenzene Pentachlorophenol Pyridine 2,4,5-Trichlorophenol 2,4,6-Trichlorophenol	1,4-Dichlorobenzene       0.050         2,4-Dinitrotoluene       0.050         Hexachlorobenzene       0.050         Hexachlorobutadiene       0.050         Hexachloroethane       0.050         Nitrobenzene       0.050         Pentachlorophenol       0.25         Pyridine       0.10         2,4,5-Trichlorophenol       0.050         2,4,6-Trichlorophenol       0.050			



Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522006.D Report Date: 22-May-2000 12:45

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file : \\QPITPA02\D\chem\71.i\s052200.b\S0522006.D

Lab Smp Id: DDF2D101 Client Smp ID: INTRA-LAB BLANK

Inj Date : 22-MAY-2000 12:05

Operator : SAB: 2ul inj (100ul+1ul) int Inst ID: 71.i

Smp Info : c0e130194-sblk 5/18/00 8270tclp

Misc Info : ddf2d101,s052200.b,8270clp.m,1-tclp.sub

Comment

Method : \\QPITPA02\D\chem\71.i\s052200.b\8270clp.m

Meth Date : 22-May-2000 12:37 bachas Quant Type: ISTD

Cal Date : 19-MAY-2000 15:44 Cal File: S0519CC6.D

Als bottle: 8 QC Sample: BLANK

Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: 1-tclp.sub

Target Version: 4.03
Processing Host: PITPC050

Concentration Formula: Amt \* DF \* Uf \* Vt/(Vo \* Vi)

Name	Value	Description
DF Uf Vt Vo Vi	0.001 1000.000 200.000	Dilution Factor ng unit correction factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected (uL)

							CONCENTRA	ATIONS
		QUANT SIG					ON-COLUMN	PINAL
Co	ompounds	MASS	RT	EXP RT	REL RT	RESPONSE	( ng)	(mg/L)
-			==	505000				******
*	1 1,4-Dichlorobenzene-d4	152	5.030	5 029	(1.000)	161868	40.0000	(a)
*	2 Naphthalene-d8	136	6.569	6.562	(1 000)	630828	40.0000	(a)
*	3 Acenaphthene-d10	164	9.603	9.596	(1.000)	338411	40 0000	(a)
*	4 Phenanthrene-dl0	188	12.904	12.898	(1 000)	564305	40 0000	(a)
*	5 Chrysene-d12	240	19 534	19.533	(1.000)	527745	40.0000	(a)
*	6 Perylene-d12	264	22 894	22.893	(1.000)	648566	40 0000	(a)
	10 Pyridine	79	Com	pound No	t Detecte	đ		
	28 1,4-Dichlorobenzene	146	Con	pound No	ot Detecte	đ		
М	34 Cresols, total	100	Com	pound No	ot Detecte	đ		
	31 2-Methylphenol	108	Com	pound No	ot Detecte	đ		
	35 4-Methylphenol	108	Con	pound No	ot Detecte	đ		
	38 Hexachloroethane	117	Con	pound No	ot Detecte	d		
	39 Nitrobenzene	77	Сοπ	pound No	ot Detecte	d.		
	59 Hexachlorobutadiene	224	Con	pound No	ot Detecte	đ.		

	OULDNIE OFO		CONCENTRATIONS
	QUANT SIG		ON-COLUMN FINAL
Compounds	MASS	RT EXP RT REL RT RESPONSE	( ng) ( mg/L)
****************	2289	四百 医医疗法院区 日野江江日本 计多点体出版设置	
69 2,4,6-Trichlorophenol	196	Compound Not Detected	
70 2,4,5-Trichlorophenol	196	Compound Not Detected	
91 2,4-Dinitrotoluene	165	Compound Not Detected	
113 Hexachlorobenzene	283	Compound Not Detected.	
117 Pentachlorophenol	265	Compound Not Detected.	
\$ 172 Nitrobenzene-d5	82	5 687 5.686 (0.866) 407907	63.9433 0.15986(a)
\$ 173 2-Fluorobiphenyl	172	8.310 8.309 (0.865) 699615	68 1531 0.17038(a)
\$ 174 Terphenyl-d14	244	16.975 16 969 (0 869) 796415	67.4835 0.16871(a)
\$ 175 Phenol-d5	99	4.715 4.719 (0 937) 711350	98 5136 0 24628(a)
\$ 176 2-Fluorophenol	112	3.705 3.709 (0.737) 513293	94 9950 0.23749(a)
\$ 177 2,4,6-Tribromophenol	330	11.345 11.338 (0.879) 138807	99.0802 0.24770(a)
\$ 178 2-Chlorophenol-d4	132	4 832 4 831 (0.961) 546795	108 089 0.27022(a)
\$ 179 1,2-Dichlorobenzene-d4	152	5.244 5 237 (1 042) 239430	67.3668 0 16842(a)

#### QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).

# UXB INTERNATIONAL CHECK SAMPLE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:C0E190000 269

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DDF2D102 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/22/00

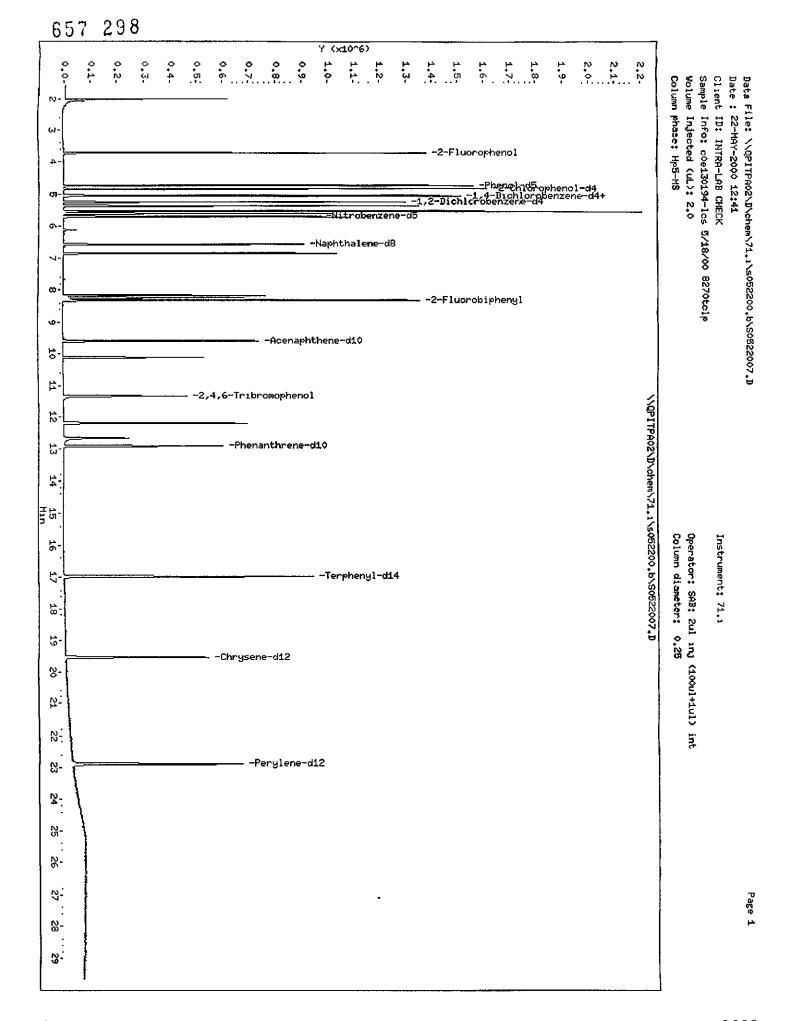
Moisture %:NA

QC Batch: 0140269

Client Sample Id: CHECK SAMPLE

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L Q
106-46-7	1,4-Dichlorobenzene	0.173
121-14-2	2,4-Dinitrotoluene	0.136
118-74-1	Hexachlorobenzene	0.191
87-68-3	Hexachlorobutadiene	0.179
67-72-1	Hexachloroethane	0.170
98-95-3	Nitrobenzene	0.168
87-86-5	Pentachlorophenol	0.162
110-86-1	Pyridine	0.272
95-95-4	2,4,5-Trichlorophenol	0.171
88-06-2	2,4,6-Trichlorophenol	0.173
1319-77-3	Cresols (total)	0.489



Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522007.D Report Date: 22-May-2000 13:36

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file : \\QPITPA02\D\chem\71.i\s052200.b\S0522007.D

Lab Smp Id: DDF2D102 Client Smp ID: INTRA-LAB CHECK

Inj Date : 22-MAY-2000 12:41

Operator : 045183 Inst ID: 71.i

Smp Info : c0e130194-lcs 5/18/00 8270tclp

Misc Info : ddf2d102,s052200.b,8270clp.m,1-tclp.sub

Comment

Method : \\QPITPA02\D\chem\71.i\s052200.b\8270clp.m
Meth Date : 22-May-2000 12:37 bachas Quant Type: I Quant Type: ISTD Cal File: S0519CC6.D Cal Date : 19-MAŶ-2000 15:44 Als bottle: 9

QC Sample: LCS Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: 1-tclp.sub

Target Version: 4.03 Processing Host: PITPC050

Concentration Formula: Amt \* DF \* Uf \* Vt/(Vo \* Vi)

Name	Value	Description
DF Uf Vt Vo Vi	0.001 1000.000 200.000	Dilution Factor  ng unit correction factor  Volume of final extract (uL)  Volume of sample extracted (mL)  Volume injected (uL)

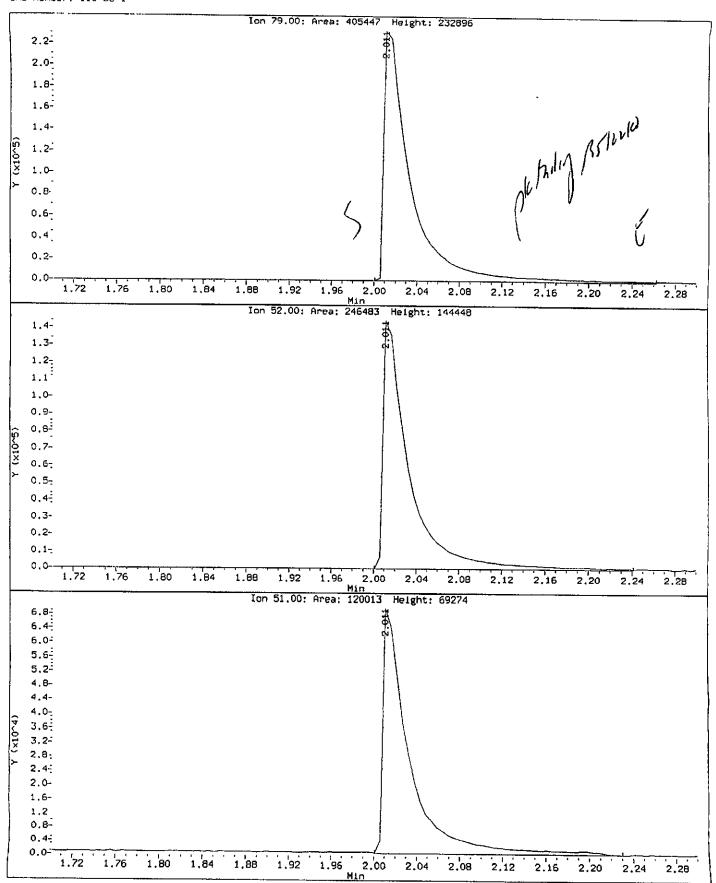
								CONCENTR	ATIONS
_		_	QUANT SIG					ON-COLUMN	FINAL
Con	oqı	unds	MASS	RT	EXP RT	REL RT	RESPONSE	( ng)	(mg/L)
	<b>#</b> #	(B) 在正正正可含含的过去式和过去分词	•===		*****			*****	
*	1	1,4-Dichlorobenzene-d4	152	5.035	5 029	(1 000)	161298	40 0000	(a)
*	2	Naphthalene-d8	136	6.563	6.562	(1 000)	635000	40.0000	(a)
*	3	Acenaphthene-d10	164	9.597	9 596	(1 000)	326610	40.0000	-
*	4	Phenanthrene-d10	188	12.899		(1.000)	529319	40.0000	(a)
*	5	Chrysene-d12	240	19 523		(1 000)	491727	40 0000	(a)
*	6	Perylene-d12	264	22 888		(1 000)	616566	40.0000	(a)
	10	Pyridine	79	2.011		(0.399)	405446	108 788	(a)
	28	1,4-Dichlorobenzene	146	5.051		(1.003)	437236		0 27197(aM)
M	34	Cresols, total	100	0.002	3 030	(1.003)		69 3127	0 17328(a)
	31	2-Methylphenol	108	F 366	- 266		1101657	195.409	0 48852(a)
		4-Methylphenol		5 366		(1 066)	368204	67.2332	0 16808(a)
			108	5.532	5.526	(1.099)	733453	128 176	0 32044(a)
	38		117	5.601	5.601	(1.112)	167245	68.1931	0 17048(a)
	39	Nitrobenzene	77	5 713	5.707	(0.871)	465607	67.1198	0.16780(a)
	59	Hexachlorobutadiene	225	6.851	6.845	(1.044)	168139	71.6908	0 17923(a)

					CONCENTRA	ATIONS
	QUANT SIG				ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT REL RT	RESPONSE	(ng)	(mg/L)
	8=28	==	9852== BE33E5	202-3685		###C===
69 2,4,6-Trichlorophenol	196	8.171	8 170 (0 851)	184505	69.3385	0.17335(a)
70 2,4,5-Trichlorophenol	196	8 235	8 239 (0.858)	194645	68.5781	0.17144(a)
91 2,4-Dinitrotoluene	165	10 115	10 115 (1.054)	181220	54.4380	0 13609(a)
113 Hexachlorobenzene	284	12.199	12 193 (0 946)	192275	76 5677	0 19142(a)
117 Pentachlorophenol	266	12.642	12 647 (0.980)	94061	64 9729	0.16243 (aM)
\$ 172 Nitrobenzene-dS	82	5.692	5.686 (0.867)	418081	65 1074	0 16277(a)
\$ 173 2-Fluorobiphenyl	172	8 310	8 309 (0.866)		72.1947	0.18049(a)
\$ 174 Terphenyl-d14	244	16 969	16 969 (0 869)	763089	69.3960	0.17349(a)
\$ 175 Phenol-d\$	99	4 720	4 719 (0.937)		94.8326	0.17349(a) 0.23708(a)
\$ 176 2-Fluorophenol	112	3.710	3 709 (0.737)	492894	91.5425	0.22886(a)
\$ 177 2,4,6-Tribromophenol	330	11.339	11.338 (0.879)	131989	100.441	0.22880(a) 0.25110(a)
\$ 178 2-Chlorophenol-d4	132	4 832	4.831 (0.960)	531455	105.429	, -
\$ 179 1,2-Dichlorobenzene-d4	152	5 243	5 237 (1 041)	246938	69.7250	0 26357(a) 0 17431(a)

#### QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).
 M - Compound response manually integrated.

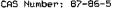
Compound: Pyridine CAS Number, 110-86-1

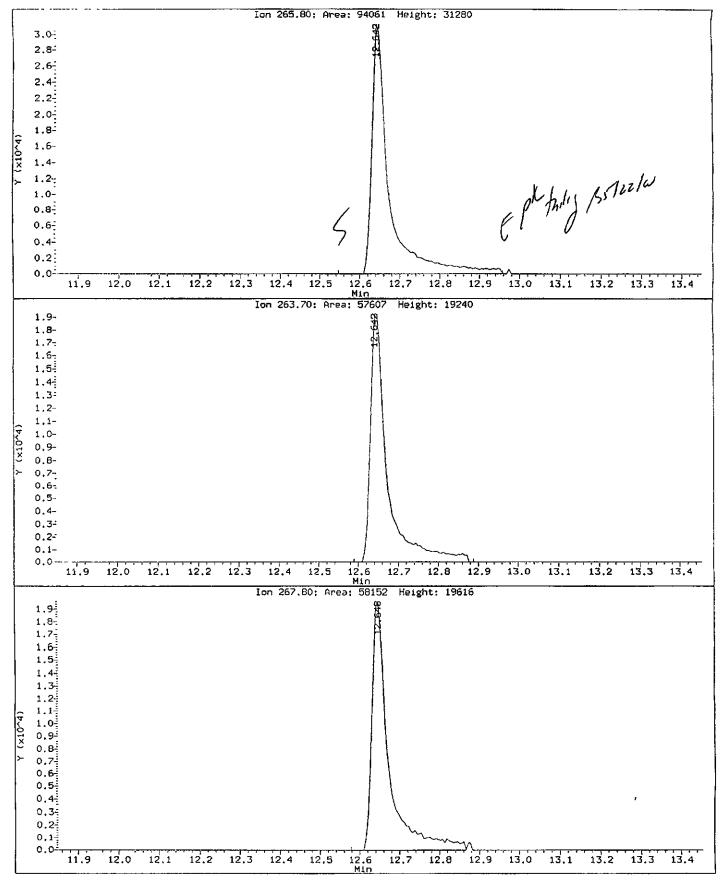


65; 302
Data File: \\QPITPA02\D\chem\71.1\s052200.b\S0522007.D
Injection Date: 22-MAY-2000 12:41

Instrument: 71.i Client Sample ID: INTRA-LAB CHECK

Compound: Pentachlorophenol CAS Number: 87-86-5





## UXB INTERNATIONAL MATRIX SPIKE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID Lab Sample ID:COE130194 001

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Date Received: 05/13/00 Work Order: DD6A411G Date Extracted:05/18/00 Dulution factor: 1 Date Analyzed: 05/22/00

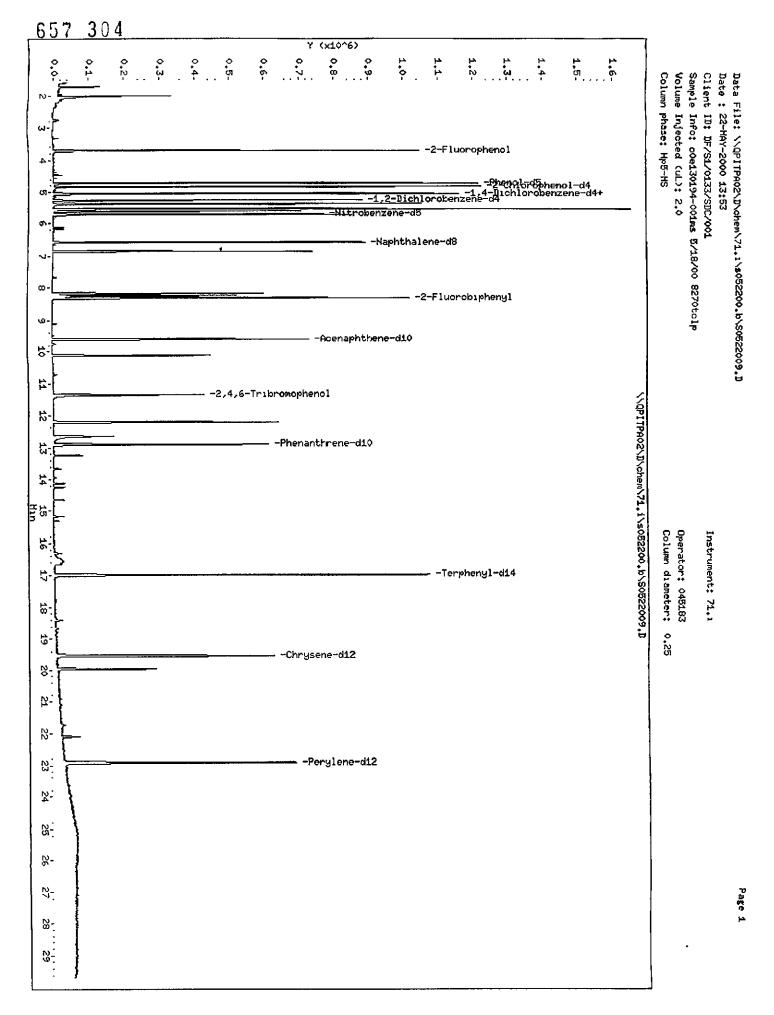
Moisture %:10

Client Sample Id: DF/S1/0133/SDC/001 A

#### CONCENTRATION UNITS:

QC Batch: 0140269

CAS NO.	COMPOUND (ug/L or	r ug/kg) mg/L Q
106-46-7	1,4-Dichlorobenzene	0.127
121-14-2	2,4-Dinitrotoluene	0.118
118-74-1	Hexachlorobenzene	0.174
87-68-3	Hexachlorobutadiene	0.130
67-72-1	Hexachloroethane	0.118
98-95-3	Nitrobenzene	0.131
87-86-5	Pentachlorophenol	0.138
110-86-1	Pyridine	0.157
95-95-4	2,4,5-Trichlorophenol	0.142
88-06-2	2,4,6-Trichlorophenol	0.140
1319-77-3	Cresols (total)	0.375



Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522009.D

Report Date: 22-May-2000 14:37

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file : \\QPITPA02\D\chem\71.i\s052200.b\S0522009.D

Lab Smp Id: DD6A411G Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 22-MAY-2000 13:53

Operator : 045183 Inst ID: 71.i

Smp Info : c0e130194-001ms 5/18/00 8270tclp

Misc Info : dd6a411g, s052200.b, 8270clp.m, 1-tclp.sub

Comment

Method : \\QPITPA02\D\chem\71.i\s052200.b\8270clp.m Meth Date : 22-May-2000 12:37 bachas Quant Type: Is Quant Type: ISTD Cal Date : 19-MAY-2000 15:44 Cal File: S0519CC6.D

Als bottle: 11 QC Sample: MS

Dil Factor: 1.00000

38 Hexachloroethane

39 Nitrobenzene 59 Hexachlorobutadiene

Compound Sublist: 1-tclp.sub Integrator: HP RTE

Target Version: 4.03 Processing Host: PITPC050

Concentration Formula: Amt \* DF \* Uf \* Vt/(Vo \* Vi)

117

77

225

Name	Value	Description	July
DF Uf		Dilution Factor ng unit correction factor	55/WW
Vt		Volume of final extract (uL)	1
Vo		Volume of sample extracted (mL)	
Vi	2.000	Volume injected (uL)	

						CONCENTRA	ATIONS
	QUANT SIG					ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT	REL RT	response	( ng)	( mg/L)
		25	*****	=====	***		****
* 1 1,4-Dichlorobenzene-d4	152	5 034	5.029	(1.000)	165477	40.0000	(a)
* 2 Naphthalene-d8	136	6 572	6.562	(1.000)	633807	40.0000	(a)
<ul> <li>3 Acenaphthene-d10</li> </ul>	164	9 601	9.596	(1.000)	317069	40.0000	(a)
* 4 Phenanthrene-d10	188	12.902	12 898	(1 000)	528651	40 0000	(a)
* 5 Chrysene-d12	240	19 532	19 533	(1 000)	549885	40 0000	(a)
* 6 Perylene-d12	264	22 898	22.893	(1.000)	656369	40 0000	(a)
10 Pyridine	79	2.015	2 000	(0 400)	240197	62 8217	0 15705 (aM)
28 1,4-Dichlorobenzene	146	5.055	5.050	(1 004)	329410	50.9008	0.12725(a)
M 34 Cresols, total	100				868515	150.162	0 37540(a)
31 2-Methylphenol	108	5.365	5 365	(1.066)	289982	51.6128	0.12903(a)
35 4-Methylphenol	108	5.536	5.526	(1 100)	578533	98 5493	0.24637(a)

5 605 5.601 (1.114)

5.712 5 707 (0 869)

6 855 6.845 (1.043)

119108

362252

121827

47.3391

52.3189

52.0422

0.11835(a)

0.13080(a)

0 13010(a)

Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522009.D

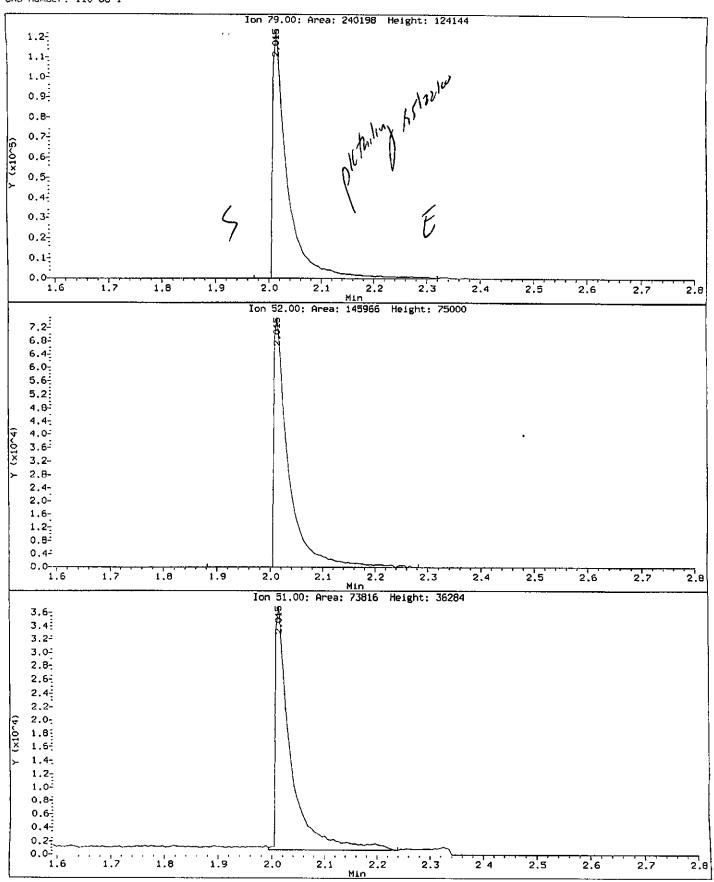
Report Date: 22-May-2000 14:37

							CONCENTRA	TIONS
		QUANT SIG					on-column	FINAL
Compo	unds	MASS	RT	EXP RT	REL RT	RESPONSE	( ng)	( mg/L)
====	프로보보다로리를등등등등록	₹##	==		*****	********	******	*******
69	2,4,6-Trichlorophenol	196	8 175	8 170	(0.851)	145057	56.1540	0.1403B(a)
70	2,4,5-Trichlorophenol	196	8.239	8 239	(0.858)	156600	56 8342	0 14208(a)
91	2,4-Dinitrotoluene	165	10 125	10 115	(1 055)	152278	47.1204	0 11780(a)
113	Hexachlorobenzene	284	12 203	12.193	(0.946)	174333	69.5105	0.17378(a)
117	Pentachlorophenol	266	12.657	12 647	(0.981)	79905	55 2649	0.13816 (aM)
\$ 172	Nitrobenzene-d5	82	5 696	5 686	(0.867)	332132	51.8200	0.12955(a)
\$ 173	2-Fluorobiphenyl	172	8 314	8.309	(0.866)	544048	56.5659	0.14141 (a)
\$ 174	Terphenyl-d14	244	16 979	16 969	(0 869)	871902	70 9053	0.17726(a)
\$ 175	Phenol-d5	99	4 718	4 719	(0 937)	540144	73 1725	0.18293(a)
\$ 176	2-Fluorophenol	112	3.709	3 709	(0.737)	380537	68 8902	0.17222(a)
\$ 177	2,4,6-Tribromophenol	330	11,348	11 338	(0.880)	118400	90 2135	0.22553(a)
\$ 178	2-Chlorophenol-d4	132	4.836	4 831	(0 961)	433921	83 9063	0.20976(a)
\$ 179	1,2-Dichlorobenzene-d4	152	5.247	5.237	(1 042)	176201	48,4954	0.12124(a)

#### QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation (BLOQ).
- M Compound response manually integrated.

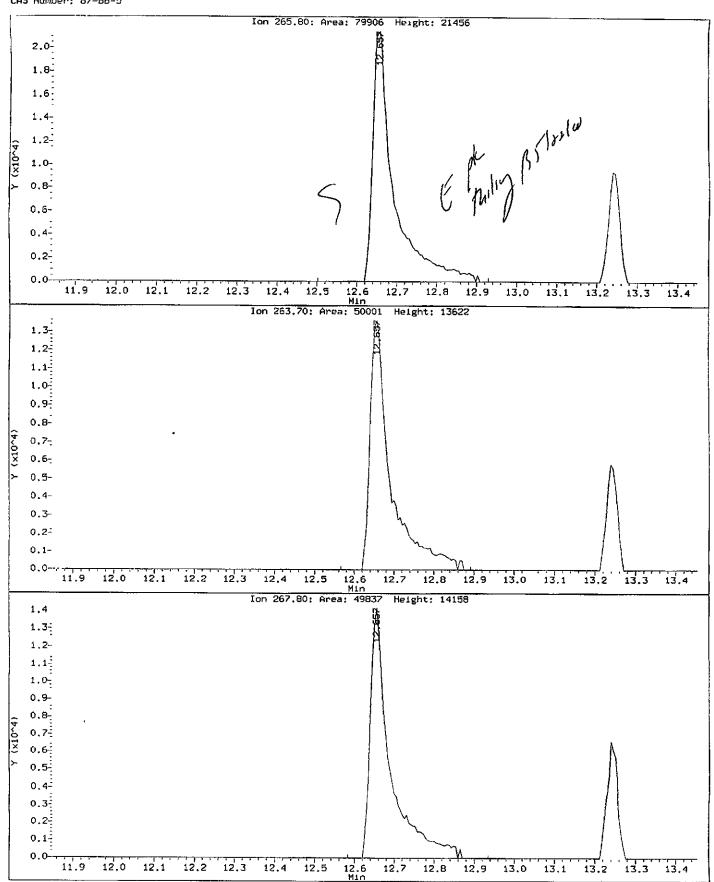
Compound Pyridine CAS Number: 110-86-1



ģ

Data File: \\QPITPAO2\D\chem\71.i\\s052200.b\\S0522009.D Injection Date: 22-MAY-2000 13:53 Instrument: 71.i Client Sample ID: DF/S1/0133/SDC/001

Compound: Pentachlorophenol CAS Number: 87-86-5



#### UXB INTERNATIONAL MATRIX SPIKE DUPLICATE COMPOUNDS

Lab Name · Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID

Lab Sample ID:COE130194 001

Method: SW846 8270C

Base/Neutrals and Acids (8270C)

Sample WT/Vol: 200 / mL Work Order: DD6A411H Dilution factor: 1

Date Received: 05/13/00 Date Extracted:05/18/00 Date Analyzed: 05/22/00

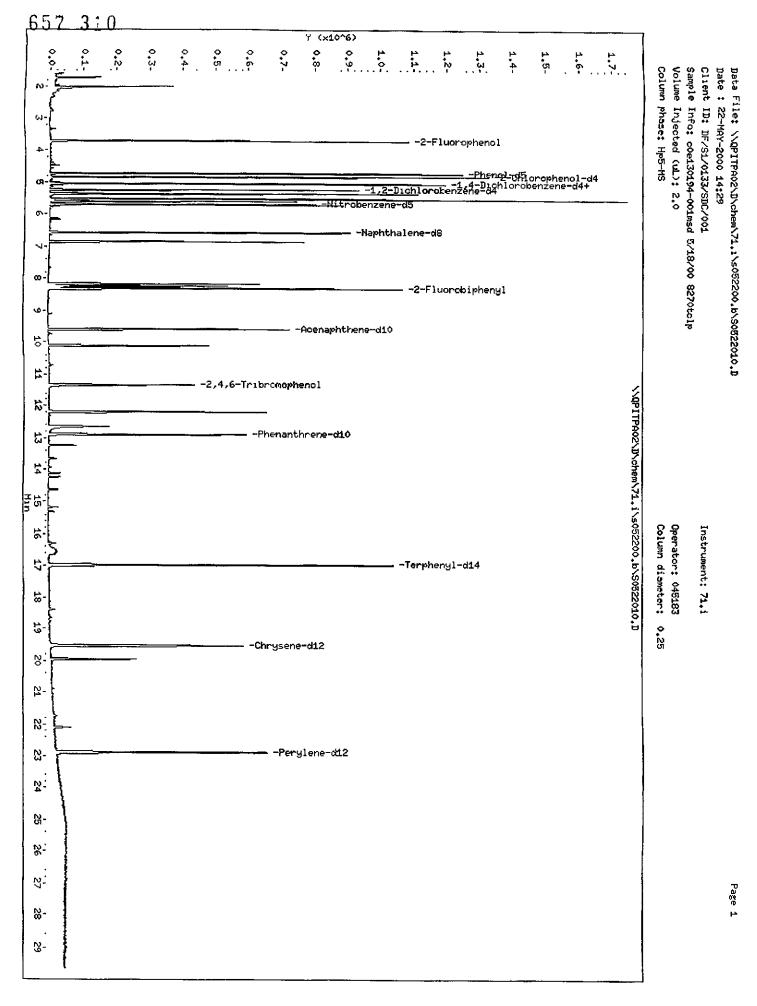
Moisture %:10

QC Batch: 0140269

Client Sample Id: DF/S1/0133/SDC/001 A

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L Q
106-46-7	1,4-Dichlorobenzene	0.131
121-14-2	2,4-Dinitrotoluene	0.120
118-74-1	Hexachlorobenzene	0.181
87-68-3	Hexachlorobutadiene	0.135
67-72-1	Hexachloroethane	0.122
98-95-3	Nitrobenzene	0.136
87-86-5	Pentachlorophenol	0.143
110-86-1	Pyridine	0.164
95-95-4	2,4,5-Trichlorophenol	0.146
88-06-2	2,4,6-Trichlorophenol	0.143
1319-77-3	Cresols (total)	0.391



Data File: \\QPITPA02\D\chem\71.i\s052200.b\S0522010.D Report Date: 22-May-2000 15:14

Page 1

#### STL Pittsburgh

Semivolatile REPORT SW-846 Method 8270

Data file : \\QPITPA02\D\chem\71.i\s052200.b\S0522010.D

Lab Smp Id: DD6A411H Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 22-MAY-2000 14:29

Operator : 045183 Inst Smp Info : c0e130194-001msd 5/18/00 8270tclp Inst ID: 71.i

Misc Info : dd6a411h, s052200.b, 8270clp.m, 1-tclp.sub

Comment

: \\QPITPA02\D\chem\71.i\s052200.b\8270clp.m Method Meth Date : 22-May-2000 12:37 bachas Quant Type: ISTD Cal Date :  $19-MA\bar{Y}-2000\ 15:44$ Cal File: S0519CC6.D Als bottle: 12

QC Sample: MSD Dil Factor: 1.00000

Integrator: HP RTE Compound Sublist: 1-tclp.sub

Target Version: 4.03 Processing Host: PITPC050

Concentration Formula: Amt \* DF \* Uf \* Vt/(Vo \* Vi)

Name	Value	Description	15/201W
DF Uf Vt Vo Vi	$0.001 \\ 1000.000$	Dilution Factor ng unit correction factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected (uL)	<i>(</i> ))**

						CONCENTRA	ATIONS
	QUANT SIG					ON-COLUMN	FINAL,
Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	( ng)	( mg/L)
19 年间 19 月 15 日 15 日 15 日 15 日 15 日 15 日 15 日 15	=400		=====			065===	======
* 1 1,4-Dichlorobenzene-d4	152	5.035	5 029	(1 000)	165720	40.0000	(a)
* 2 Naphthalene-d8	136	6.568	6 562	(1 000)	635621	40.0000	(a)
* 3 Acenaphthene-d10	164	9.602	9 596	(1 000)	322693	40 0000	(a)
* 4 Phenanthrene-d10	188	12.904	12 898	(1 000)	522771	40.0000	(a)
* 5 Chrysene-d12	240	19 533	19 533	(1 000)	519636	40 0000	(a)
* 6 Perylene-d12	264	22 899	22.893	(1 000)	634789	40.0000	(a)
10 Pyridine	79	2 016	2.000	(0.400)	251726	65.7404	0.16435 (aM)
28 1,4-Dichlorobenzene	146	5.056	5.050	(1 004)	340560	52.5466	0 13137(a)
M 34 Cresols, total	100				907164	156.588	0.39147(a)
31 2-Methylphenol	108	5.366	5.365	(1.066)	299454	53.2206	0.13305(a)
35 4-Methylphenol	108	5 531	5.526	(1.099)	607710	103.368	0.25842(a)
38 Hexachloroethane	117	5.606	5.601	(1.114)	123443	48 9901	0.12248(a)
39 Nitrobenzene	77	5.713			377412	54 3528	0.13588(a)
59 Hexachlorobutadiene	225	6 856		(1 044)	126637	53.9425	0.13388 (a) 0.13486 (a)

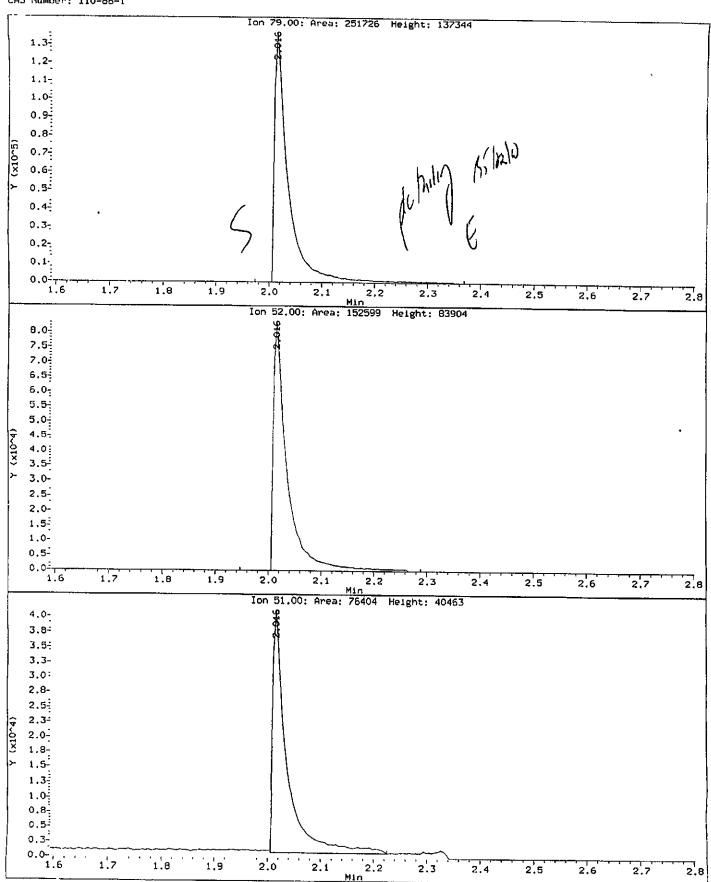
657 312
Data File: \QPITPA02\D\chem\71.i\s052200.b\S0522010.D
Report Date: 22-May-2000 15:14

						CONCENTRA	ATIONS
	QUANT SIG					ON-COLUMN	FINAL
Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	(ng)	( mg/L)
, 자료 및 도보으로 등록 및 영향 및 등록 및 등로 표표 표표 보건 는	#===	==	=====		*=====	=====	
69 2,4,6-Trichlorophenol	196	8.176	8 170	(0 851)	150179	57 1236	0.14281(a)
70 2,4,5-Trichlorophenol	196	8.240	8 239	(0.858)	163265	58 2204	0 14555(a)
91 2,4-Dinitrotoluene	165	10.120	10.115	(1.054)	157233	47 8057	0 11951(a)
113 Hexachlorobenzene	284	12.204	12.193	(0 946)	179626	72.4266	0.18107(a)
117 Pentachlorophenol	266	12 652	12.647	(0 981)	81897	57 2795	0.14320 (aM)
\$ 172 Nitrobenzene-d5	82	5.692	5 686	(0 867)	348101	54 1565	0.13539(a)
\$ 173 2-Fluorobiphenyl	172	8.315	8.309	(0.866)	565813	57 8035	0.14451(a)
\$ 174 Terphenyl-d14	244	16.980	16 969	(0.869)	851843	73 3066	0.18327(a)
\$ 175 Phenol-d5	99	4.719	4.719	(0.937)	562512	76.0909	0.19023(a)
\$ 176 2-Fluorophenol	112	3 710	3 709	(0.737)	393249	71 0872	0.17772(a)
\$ 177 2,4,6-Tribromophenol	330	11 344	11 338	(0 879)	122166	94,1300	0.23532(a)
\$ 178 2-Chlorophenol-d4	132	4.832	4 831	(0 960)	449750	86 8396	0.21710(a)
\$ 179 1,2-Dichlorobenzene-d4	152	5.243	5.237	(1.041)	182631	50 1914	0.12548(a)

#### QC Flag Legend

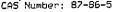
- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).
- M Compound response manually integrated.

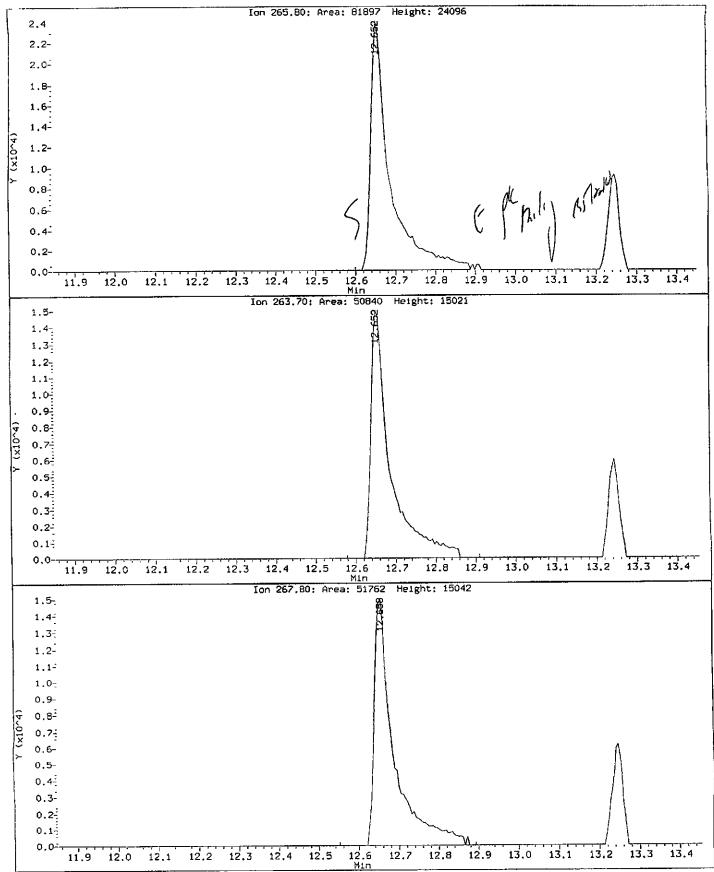
Compound: Pyridine CAS Number: 110-86-1



Data File: \\QPITPA02\D\chem\71.1\s052200.b\90522010.D Injection Date: 22-MAY-2000 14:29 Instrument: 71.1 Client Sample ID: DF/S1/0133/SDC/001

Compound: Pentachlorophenol CAS Number: 87-86-5





# GC/MS SEMIVOLATILE MISCELLANEOUS

CO  Quanterra N:\QA\LOGBOOKS_PDF\C_L L_EXT.DOC May-00 LO GO	Batch Number: (	rom above)	Prep Analyst: WT WT	22.	20.	18.	16.	14.	12.	11.	4	6. 003		NS.	1. LOE/30194 BIK	Number Samp	over 5/19/2000 Time	
	1 1	Date Time And 20.00 (6:10)	Extract(s) Received				15 2 JM	0 1000			4				N/H &00	Chent ID Ex	7 Prep Meth # 35,20C H2SO4 [86 - Analysis 82,70 NaOH. 9,2)	Tolo
Log Book Number OP-00-0037	Review	I Location Day  1. ORG Rad S-20  1. ORG RAD S 724	KG/SM   WT								7211	7/2/11	r	8/2/11		Volu	-12 Solvent MeCla	· •
	$\prod_{i \in \mathcal{I}} H_i$	or USA Constant	Extract(s) F								*		NA.	8-651-h67	194-17×1531210	Surr Vol	\$ 5 × × × × × × × × × × × × × × × × × ×	d To Your Showers
Page 1 of 100 STL Pittsbu	Date 5-20 - Sign	Che REDITORNA											NA	2.5	0.5	+	Clean up Method	STL Pittsburgh 450 William Pitt Way Pittsburgh, PA 15238 412-820-8380

TCLP (Method 1311)

Quanterra Incorporated 450 William Pitt Way Pittsburgh, Pennsylvania 15238 412/826-5477 FAX: 412/826-5571

*Muanterra* 

Sample ID Cot Number | 194 Wgt/Vol 5-16-00 Fluid Determination Init pH Final pH Checked by Calent Codes 39411 Extract Fluid #1, or #2 Conc : HNO<sub>3</sub> to Acidity (for Metals) Filter lot # 8320701W 5000 Sample ID ACCHINE ARDS SNAC SOLUTION NUMBER 1-LOS BOOK NUMBER 945404-12 pH instrument Wgt (g)
Wet/Dry Fluid Vol. TCLP Extraction SN9087030 038 Solution Number 2 - Log Book Number C94817 96-MT-526 Final pH After Tumbling Environmental
Services

317

* = Sample determined to have free-liquid, % solid determination was performe <5 = Extraction Fluid No. 1 5.7 mL Glacial Acetic Acid dil 500 mL + 64.3 mL of : to 1L (pH 4.93 ± 0.05)  >5 = Extraction Fluid No. 2 5.7 mL Glacial Acetic Acid dil to 1L (pH 2.88 ± 0.05)  \$300 M	(Record line number from above)					00647	DOGAL	1	7 Sample ID
* = Sample determined to have free-liquid, % solid determination was performed.  <5 = Extraction Fluid No. 1 5.7 mL Glacial Acetic Acid dil 500 mL + 64.3 mL of 1N Na OH dil to 1L (pH 4.93 + 0.05)  >5 = Extraction Fluid No. 2 5 7 mL Glacial Acetic Acid dil to 1L (pH 2 88 + 0.05)	Date 5-/6-0		5,0	10.7		₩ / ₩ 1.5C	9.30	5,00, 18.5.19.6	WgtVol Init pH
nination was performed. 500 mL + 64.3 mL of 1N to 1L (pH 2 88 ± 0 05)	Time Apalyst			8\		1.70	1.79	1 1.83 50	Final pH
Na OH dil	Analyst					1		# #	Extract Fluid to the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the first that the
Date/Time On  \$-\ \( \\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \	Location Protifices 5	13 12.	10	, 9	7. ED	(2) 4.0	2. C		Conc : HNO3 to Acidity (for Metals)
15:42 pH	Date Time		8	G	6DD6A4	5DOLAC 5DOLAC	2. 00CM 3. DOGAS	1. PB-7C19 5-16-00	Sample ID
Calibration 4:00 4:00	Expacts) Helinquisted		76.00		4		(0).003	20002	Wgt (g) Wet/Dry Fluid Vol.
30 ± 2   7.0    7.0					7 4	6		Jal Glass	
STL Pittsburgh	Location				.30	(2.37	36	4.93	Final pH After Tumbling

318 Comment: STL PITTSBURGH-RUN LOGBOOK hp597371-A
Operator: SAB: 2ul inj (100ul+1ul) int
Data Path: D:\HPCHEM\1\DATA\s051900.b\
Pre-Seq Cmd:
Post-Seq Cmd:

Method Sections To Run On A Barcode Mismatch

(X) Inject Anyway
( ) Don't Inject

(X) Full Method

( ) Reprocessing Only

Line Type Vial DataFile Method Sample Name 1 Sample C [144] 1 S0519DF2 DFTPP1 DFTPP050 (25ppb) 194-158-6 2 Sample\_ - 2 S0519CC1 EARLY sstd50(25.0ug/ml) 194-176-9 8270 3 Sample AQC 4 S0519CC3 EARLY sstd50(25.0ug/ml) 194-176-9 8270 sstd20(10.0ug/ml) 194-175-10 8270 4 Sample AOC 3 S0519CC2 EARLY 5 Sample AQC 5 S0519CC4 EARLY sstd80(40.0ug/ml) 194-175-6 8270 6 Sample ACC 6 S0519CC5 EARLY sstd120(60.0ug/ml) 194-175-13 e270 7 Sample ACC 7 S0519CC6 EARLY sstd160(80.0ug/ml) 194-175-14 5270 8 Sample ACC 8 S0519VER EARLY ver050(25.0ug/ml) 194-175-7 8270 9 Sample -1 S0519DF3 DFTPP1 dftpp050(25.0ug/ml) 194-158-6 10 Sample 1708 1 S0519DF4 DFTPP1 dftpp050(25.0ug/ml) 194-158-6 11 Sample 2 S0519CC7 EARLY sstd50(25.0ug/ml) 194-176-9 8270 Acc 12 Sample ARC 9 S0519001 LATE c0e120206-sblk 5/17/00 8270w AUC c0e120206-lcs 5/17/00 8270w 13 Sample 10 S0519002 LATE AC 11 S0519003 LATE 14 Sample c0e120206-lcsd 5/17/00 8270w c0e120206-001 5/17/00 8270w 20 12 S0519004 LATE 15 Sample 16 Sample AULS 13 S0519005 LATE c0e120206-002 5/17/00 8270w 17 Sample 065 14 S0519006 LATE c0e120206-003 5/17/00 8270w Mar 18 Sample c0e120206-004 5/17/00 8270w Am last Sca 1.5 S0519007 LATEON c0e120206-005 5/17/00 8270w 19 Sample | Nother 16 S0519008 LATE 20 Sample 17 S0519009 LATE(7) c0e120206-006 5/17/00 8270w 21 Sample 18 S0519010 LATE c0e120206-007 5/17/00 8270w MAN 19 S0519011 LATE (1) c0e120206-008 5/17/00 8270w Schan 22 Sample 20 S0519012 LATE 23 Sample c0e120206-009 5/17/00 8270w 21 S0519013 LATE // c0e120236-001 x50 5/17/00 8270 √24 Sample| LAR 22 S0519014 LATE 25 Sample c0e120236-002 x50 5/17/00 827**0** 26 Sample 23 S0519015 LATE 27 Sample 24 S0519016 LATE 28 Sample 25 S0519017 LATE 29 Sample 26 S0519018 LATE 29 Sample 26 S0519018 LATE 29 Sample 26 S0519018 LATE 29 Sample 26 S0519018 LATE 29 Sample 26 S0519018 LATE 29 Sample 26 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 Sample 27 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519018 LATE 29 S0519 c0e120236-003 x50 5/17/00 8270 c0e120236-004 x50 5/17/00 8270 c0e120236-005 5/17/00 8270w c0e120236-006 5/17/00 8270w 🕻 29 Sample 🗸

Last Modified: Fri May 19 13:50:50 2000 Page: 1

bequence אame: ב: אברהבאי/ ב / שבעים ארם / 50 ב 200. 5

Comment: STL PITT HP597371A LOG 2ul inj (100ul+1ul IS)

Operator: SAB: 2ul inj (100ul+1ul) int Data Path: D:\HPCHEM\1\DATA\s052200.b\

657 319

Pre-Seq Cmd: Post-Seq Cmd:

Method Sections To Run

(X) Full Method

( ) Reprocessing Only

On A Barcode Mismatch FS-/44-159-7
(X) Inject Anyway
( ) Don't Inject

	Line	Type .	Vial	DataFile	Method	Sample Name
/1/2	1/W 1 2	Sample (10)			EARLY	DFTPP050 (25ppb) 194-158-6 sstd50(25.0ug/ml) 194-176-9 8
(	3	SampleAcc		S0522006		c0e130194-sblk 5/18/00 8270tc
	4	Sample M/	9	S0522007		c0e130194-lcs 5/18/00 8270tcl
	5	Sample	10	S0522008	EARLY	c0e130194-001 5/18/00 8270tcl
- D	6	Sample 477	11	S0522009	EARLY	c0e130194-001ms 5/18/00 8270t
TCU	7	Sample	12	S0522010	EARLY	c0e130194-001msd 5/18/00 8270
	8	Sample Mic	13	S0522011	EARLY	c0e130194-002 5/18/00 8270tcl
	9	Sample ACC	14	S0522012	EARLY	c0e130194-003 5/18/00 8270tcl
~	10	Sample ASC	15	S0522013	EARLY	c0e130194-004 5/18/00 8270tcl
	11	Sample	3	S0522001	LATE	c0e170161-sblk 5/19/00 hlev82
	12	Sample	4	S0522002	LATE	c0e170161-lcs 5/19/00 hlev827
	13	Sample	5	S0522003	LATE	c0e170161-001 5/19/00 hlev827
	14	Sample	6	S0522004	LATE	c0e170161-001ms 5/19/00 hlev8
	15	Sample	7	S0522005	LATE	c0e170161-001msd 5/19/00 hlev
	16	Sample	16	S0522014	LATE	c0e120236-001 x250 5/17/00 82
	17	Sample	17	S0522015	LATE	c0e120236-002 x1500 5/17/00 8
	18	Sample	18	S0522016	LATE	c0e120236-003 x500 5/17/00 82
	19	Sample	19	S0522017	LATE	c0e120236-004 x1000 5/17/00 8
hui	20	Sample	20	S0522018	LATE	c0e120236-005 x10 5/17/00 827
15/12/4	21	Sample	21	S0522019	LATE	c0e120236-006 x5 5/17/00 8270

Last Modified: Mon May 22 11:24:21 2000 Page: 1

PSR024 5/18/00

1:46:35 MT

SAMPLE CUSTODIAN REMOVAL REQUEST

PAGE 001

REQUESTED BY: GEEHRINK

METHOD: QL

Base/Neutrals and Acids (82700)

STORAGE LOCATION WORK ORDER #	PICKED CNTR# CONT	ROL # CLIENT #	ANALYSIS LOTID	SMP# SFX	MATRIX DESCRIPTION	Q1 <u>R0</u>	TY QT	
2F CLP1 DD6A4-1-02		34412 399411	A-59-QL COE130194	001	SOLID	0	3	1
2F CLP1 DD6A5-1-02		34413 399411	A-59-QL COE130194	002	SOLID	0	3	1
2F CLP1 DD6A6-1-02		34414 399411	A-59-QL COE130194	003	SOLID	0	3	1
2F CLP1 DD6A7-1-02	a	34415 399411	A-59-QL COE130194	004	SOLID	0	3	1

B Front	RECEIVED BY  Lea Deep	DATE/TIME - 5/18/2000 19:45 5/18/2000 15:10

\*, 1 \*

•

### PESTICIDE DATA

## PESTICIDE QC SUMMARY

#### SW846 8081A SURROGATE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT QESSDG:

Lot #: C0E130194

-	CLIENT ID.	SRG01	SRG02	TOT OUT
			======	
	DF/S1/0133/SDC/001 A	92	80	00
02	DF/S1/0133/SDC/001 B	93	82	00
03	DF/S1/0133/SDC/001 C	93	82	00
04	DF/S1/0133/SDC/001 D	97	80	00
05	METHOD BLK DDE2E101	95	58	00
06	LCS DDE2E102	97	91	00
	DF/S1/0133/SDC/001 A D	95	70	00
80	DF/S1/0133/SDC/001 A S	97	69	00

SURROGA	ATES	QC LIMITS
SRG01	<pre>= Decachlorobiphenyl</pre>	( 10-147)
SRG02	<pre>= Tetrachloro-m-xylene</pre>	( 39-130)

<sup>#</sup> Column to be used to flag recovery values

<sup>\*</sup> Values outside of required QC Limits

D System monitoring Compound diluted out

657 324 SW846 8081A CHECK SAMPLE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Lot #: COE180000

WO #: DDE2E102 BATCH: 0139460

	SPIKE	SAMPLE		QC	
	ADDED	CONCENT.	%	LIMITS	ļ
COMPOUND	(mg/L)	(mg/L )	REC	REC	QUAL
	=		<b>=</b>   =====	=======================================	========
Lindane	0.00250	0.00189	75	49- 137	
Heptachlor	0.00250	0.00212	85	<u> 57 - 124</u>	
Heptachlor epoxide	0.00250	0.00219	8.8	53- 135	<u> </u>
Endrin	0.00250	0 00219	88	46- 137	<u> </u>
Methoxychlor	0.00250	0.00277	111	12- 154	

NOTES(S):		

* Values outside o	f QC	limits	
Spike Recovery: _	0	out <b>of</b>	5 outside limits
COMMENTS:			

# ŚW846 8081A MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A

Level: (low/med) LOW

Lot #: C0E130194 `

WO #: DD6A411A BATCH: 0139460

   COMPOUND	SPIKE ADDED (mg/L )	SAMPLE CONCENT. (mg/L)	MS CONCENT. (mg/L )	MS % REC	LIMITS REC	QUAL
	=======	========		======	========	=========
Lindane	0.00250	ND	0.00172	69	30~ 148	Ì
Heptachlor	0.00250	ND	0.00236	94	25- 135	
Heptachlor epoxide	0.00250	ND	0.00201	80	38- 138	
Endrin	0.00250	ND	0.00212	85	28- 148	
Methoxychlor	0.00250	ND	0.00289	116	13- 154	

NOTES(S):	

# Column to be used to flag recovery and RPD values with an asterisk \* Values outside of QC limits RPD: \_\_\_0 out of \_\_\_0 outside limits Spike Recovery. 0 out of 5 outside limits COMMENTS:

SW846 8081A MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY 657 326

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID · DF/S1/0133/SDC/001 A

Level: (low/med) LOW

Lot #: C0E130194

WO #: DD6A411C BATCH: 0139460

   COMPOUND	SPIKE ADDED (mg/L )	MSD CONCENT. (mg/L)	MSD % REC	% RPD	QC I RPD	LIMITS REC	QUAL
======================================	=======	=======	=====	======	====		=======
Lindane	0.00250	0.00172	69	0.40	22	30- 148	<u> </u>
Heptachlor	0.00250	0.00238	95	1.0	32	<u>25- 135</u>	]
Heptachlor epoxide	0.00250	0.00199	80	0.77_	31	38- 138	l [
Endrin	0.00250	0.00213	85	0.29	40	28- 148	
Methoxychlor	0.00250	0.00291	117	0.78	29	13- 154	ll

		ed to flag r of QC limits	ecovery a	ind RPD	values	with an	n aster:	isk
RPD: Spike Rec	<u>0</u> out of	5 ou 0 out of	tside lim <u>5</u>	nits outside	e limits	3		
COMMENTS:	:							
						· · · · · · · · · · · · · · · · · · ·		

NOTES(S):

# 4C PESTICIDE METHOD BLANK SUMMARY

PBLK	ļ

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325

SDG No.: C0E130194

Lab Sample ID: DDE2E101

Lab File ID: C-A2372

Matrix (soil/water) WATER

Extraction: (SepF/Cont/Sonc) SW1311

Sulfur Cleanup (Y/N) N

Date Extracted: 05/18/00

Date Analyzed (1): 05/19/00

Date Analyzed (2): 05/19/00

Time Analyzed (1): 1318

Time Analyzed (2): 1318

Instrument ID (1): GC3

Instrument ID (2): GC3

GC Column (1): RTX-CLP ID: 0.53 (mm) GC Column (2): RTX-CLP2 ID: 0.53 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	EPA	LAB	DATE	DATE
	SAMPLE NO.	SAMPLE ID	ANALYZED 1	ANALYZED 2
	=========	==========	========	~=======
01	DF/S1/0133/S	DD6A4103	05/19/00	05/19/00
02	DF/S1/0133/S	DD6A411A	05/19/00	05/19/00
03	DF/S1/0133/S	DD6A411C	05/19/00	05/19/00
04	DF/S1/0133/S		05/19/00	05/19/00
05	DF/S1/0133/S		05/19/00	05/19/00
06	DF/S1/0133/S		05/19/00	05/19/00
07	LCS	DDE2E102	05/19/00	05/19/00
08 09				<del></del> ,
10				
11			<del></del>	
12				
13			<del></del>	
14				
15				
16				
17				
18				
19				
20				
21				
22				
23 24				
25				
26	<u> </u>			
20	l	l	l	l

COMMENTS:			

page 1 of 1

FORM IV PEST

OLM03.0

# PESTICIDE SAMPLE DATA

#### UXB INTERNATIONAL

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix. (soil/water) SOLID WATEL Lab Sample ID: COE130194 001

Method: SW846 8081A LM5/23 100

Pesticides (8081A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A4103 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/19/00

QC Batch: 0139460

Client Sample Id: DF/S1/0133/SDC/001 A

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L	Q
57-74-9	Chlordane (technic	al) 0.0050	ן
72-20-8	Endrin	0.00050	ן ס
76-44-8	Heptachlor	0.00050	ן ט
1024-57-3	Heptachlor epoxide	0.00050	U
58-89-9	Lindane	0.00050	ן ט
72-43-5	Methoxychlor	0.0010	וט
8001-35-2	Toxaphene	0.020	ַן

Data File: /var/chem/gc3.i/2170-E.b/c-a2366.d

Report Date: 19-May-2000 11:16

657 330

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2366.d

Lab Smp Id: DD6A4103 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 19-MAY-2000 10:42

Operator : 1891 Inst ID: gc3.i

Smp Info : DD6A4103,2170-E.b,,PEST.sub,,,
Misc Info : 130194001

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 19-May-2000 11:01 eppinged Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-a2300 Cal File: c-a2300.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF Vt		Dilution Factor Volume of final extract (uL)
٧o	100.000	Volume of sample extracted (mL)
<b>V</b> i	1.000	Volume injected

			CONCENTRA	TIONS
			ON-COLUMN	FINAL
Compounds	RT EXP RT D	LT RT RESPONSE	( ng)	( mg/L)
	<b>38 88</b> 0022 2:	****		======
5 alpha-BHC	Compound Not	Detected		
4 HEXACHLOROBENZENE	Compound Not	Detected		
6 gamma-BHC (Lindane)	Compound Not	Detected		
7 beta-BHC	Compound Not	Detected.		
9 Chlordane	Compound Not	Detected.		
10 Heptachlor	7.667 7.675	-0.008 32294	0.00674/0	.000673765(a)
8 delta-BHC	Compound Not	Detected.	_	
11 Aldrin	Compound Not	Detected.		
12 Heptachlor epoxide	Compound Not	Detected.		
13 gamma-Chlordane	Compound Not	Detected.		
14 alpha-Chlordane	Compound Not	Detected.		
15 Endosulfan I	Compound Not	Detected.		
16 4,4'-DDE	Compound Not	Detected.		
17 Dieldrin	10.942 10.942	0 000 7486	0 00185 0	000184645(a)
20 Endrin	Compound Not	Detected.		
18 Toxaphene	Compound Not	Detected.		

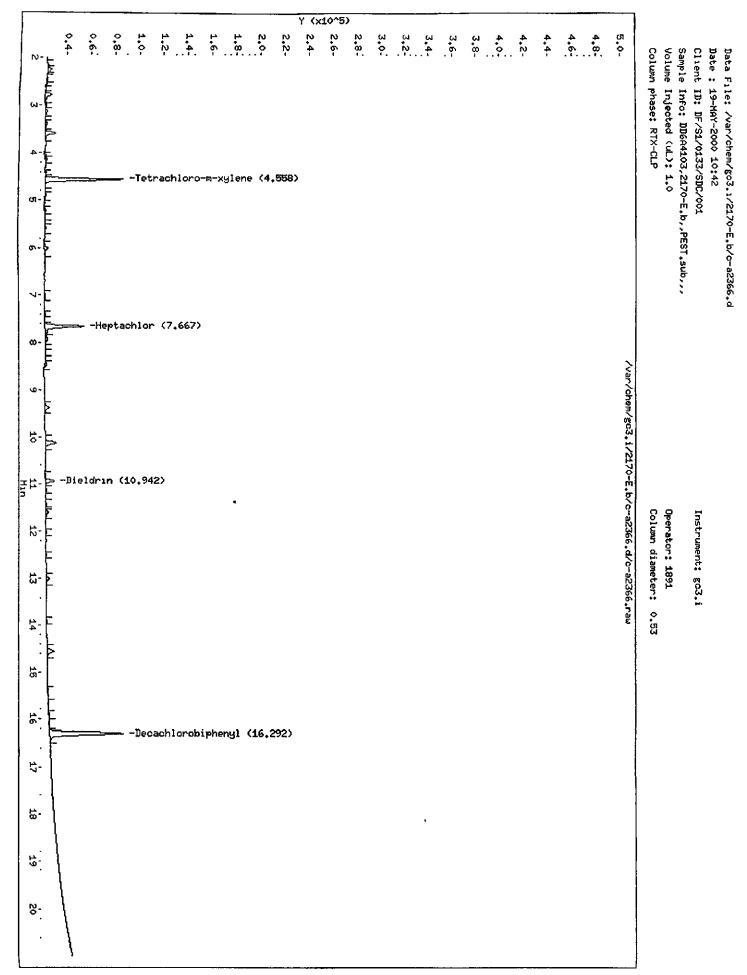
Data File: /var/chem/gc3.i/2170-E.b/c-a2366.d Report Date: 19-May-2000 11:16

					CONCENTR ON-COLUMN	ATIONS FINAL
Comp	ounds	RT	EXP RT DLT RT	RESPONSE	( ng)	( mg/L)
	· ·		*****	# # # # # # # # # # # # # # # # # # #		******
2	1 4,4'-DDD	Co	mpound Not Detecte	ed.		
2:	2 Endosulfan II	Co	mpound Not Detect	ed.		
2	3 4,4'-DDT	Co	mpound Not Detect	ed.		
2	4 Endrin aldehyde	Co	mpound Not Detect	ed.		
2	6 Endosulfan sulfate	Co	mpound Not Detecte	ed		
2	5 Methoxychlor	Co	mpound Not Detecte	ed		
` 2	7 Endrin ketone	Co	mpound Not Detect	ed.		
\$	1 Tetrachloro-m-xylene	4.558	4.558 0.000	64522	0.0159	0.00159352 (aR)
\$ 3	0 Decachlorobiphenyl	16.292	16.300 -0.008	60661	0.01841	0 00184138(aR)

# QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).
   R Spike/Surrogate failed recovery limits.

3011



Data File: /var/chem/gc3.i/2170-E.b/c-b2366.d

Report Date: 19-May-2000 12:46

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-b2366.d

Lab Smp Id: DD6A4103 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 19-MAY-2000 10:42 Operator : 1891 Inst ID: gc3.i

Smp Info : DD6A4103,2170-E.b,,PEST.sub,,,

Misc Info : 130194001

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date: 19-May-2000 11:24 eppinged Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-b2300.d

Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: PEST.sub Integrator: Falcon

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF Vt	10.000	Dilution Factor Volume of final extract (uL)
Vo		Volume of sample extracted (mL)
٧i	1.000	Volume injected

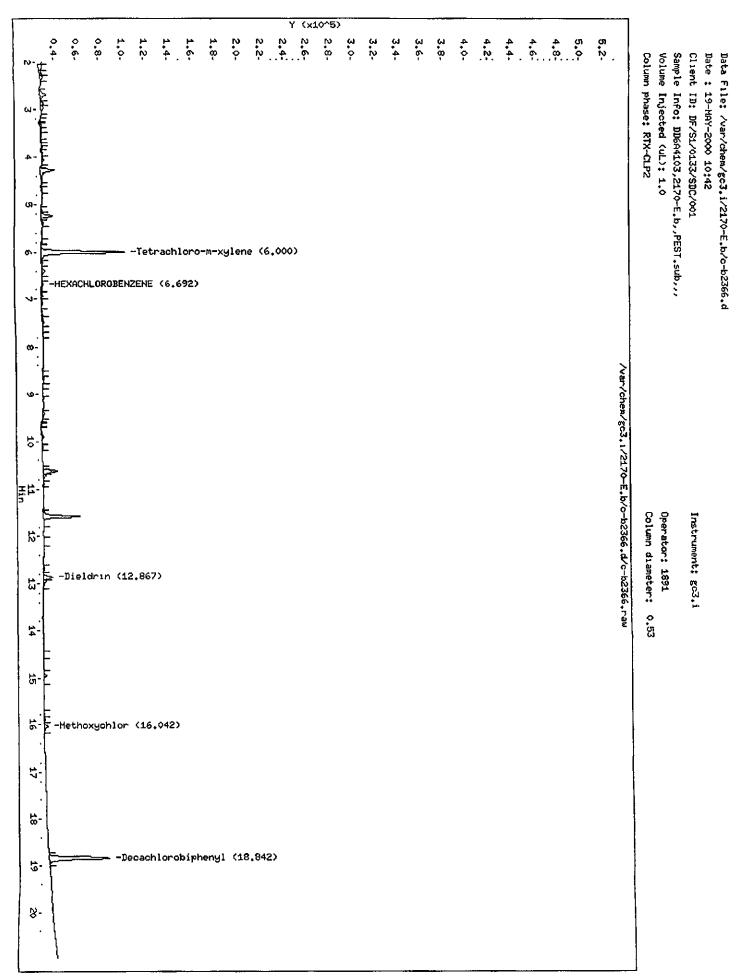
		CONCENTRATIONS
		on-column final
Compounds	RT EXP RT DLT RT RESPONSE	( ng) ( mg/L)
计计3位数据数据 网络红色对有色色色 医脊髓石 经 计 计 计		
5 alpha-BHC	Compound Not Detected.	
4 HEXACHLOROBENZENE	6 692 6 650 0.042 900	(a)
6 gamma-BHC (Lindane)	Compound Not Detected.	
7 beta-BHC	Compound Not Detected.	
9 Chlordane	Compound Not Detected	
10 Heptachlor	Compound Not Detected.	
8 delta-BHC	Compound Not Detected.	
11 Aldrin	Compound Not Detected.	
12 Heptachlor epoxide	Compound Not Detected.	
13 gamma-Chlordane	Compound Not Detected.	
14 alpha-Chlordane	Compound Not Detected.	
15 Endosulfan I	Compound Not Detected.	
16 4,4'-DDB	Compound Not Detected.	
17 Dieldrin	12 867 12.867 0.000 7830	0.00189 0.000188602(a)
20 Endrin	Compound Not Detected.	
18 Toxaphene	Compound Not Detected	

Data File: /var/chem/gc3.i/2170-E.b/c-b2366.d Report Date: 19-May-2000 12:46

		CONCENTRATIONS
		ON-COLUMN FINAL
Compounds	RT EXP RT DLT RT RESPONSE	( ng) ( mg/L)
=======================================		
21 4,4'-DDD	Compound Not Detected.	
22 Endosulfan II	Compound Not Detected.	
23 4,4'-DDT	Compound Not Detected	
24 Bndrin aldehyde	Compound Not Detected	
26 Bndosulfan sulfate	Compound Not Detected	
25 Methoxychlor	16.042 16.017 0.025 3144	0.00252 0.000251517(a)
27 Endrin ketone	Compound Not Detected.	
\$ 1 Tetrachloro-m-xylene	6.000 6.000 0.000 72974	0 01776 0.00177571(aR)
\$ 30 Decachlorobiphenyl	18.842 18.842 0.000 53208	0 01967 0.00196700 (aR)

## QC Flag Legend

- a Target compound detected but, quantitated amount
   Below Limit Of Quantitation(BLOQ).
   R Spike/Surrogate failed recovery limits.



#### UXB INTERNATIONAL

Lab Name Severn Trent Laboratories, Inc SDG Number:

Matrix: (soil/water) SOLID WATEL Lab Sample ID:COE130194 002

Method: SW846 8081A LMS23 00

Pesticides (8081A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A5103 Date Extracted 05/18/00 Dilution factor: 1 Date Analyzed: 05/19/00

QC Batch: 0139460

Client Sample Id: DF/S1/0133/SDC/001 B

#### CONCENTRATION UNITS.

CAS NO.	COMPOUND (ug/L o	or ug/kg) mg/L (	<u> </u>
57-74-9	Chlordane (technical)	0.0050	<u>U</u>
72-20-8	Endrin	0.00050	<u>u</u>
76-44-8	Heptachlor	0 00050	<u> </u>
1024-57-3	Heptachlor epoxide	0.00050	<u>U</u>
58-89-9	Lindane	0.00050	<u>U</u>
72-43-5	Methoxychlor	0.0010	U
8001-35-2	Toxaphene	0.020	<u> </u>

Data File: /var/chem/gc3.i/2170-E.b/c-a2369.d

Report Date: 19-May-2000 12:40

### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2369.d

Lab Smp Id: DD6A5103 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 19-MAY-2000 12:00 Operator : 1891 Smp Info : DD6A5103,2170-E.b,,PEST.sub,,, Inst ID: gc3.i

Misc Info : 130194002

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 19-May-2000 11:01 eppinged Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-a2300 Cal File: c-a2300.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

### Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF	1 000	Dilution Factor
۷t		Volume of final extract (uL)
Vo		Volume of sample extracted (mL)
Vi	1.000	Volume injected

#### CONCENTRATIONS

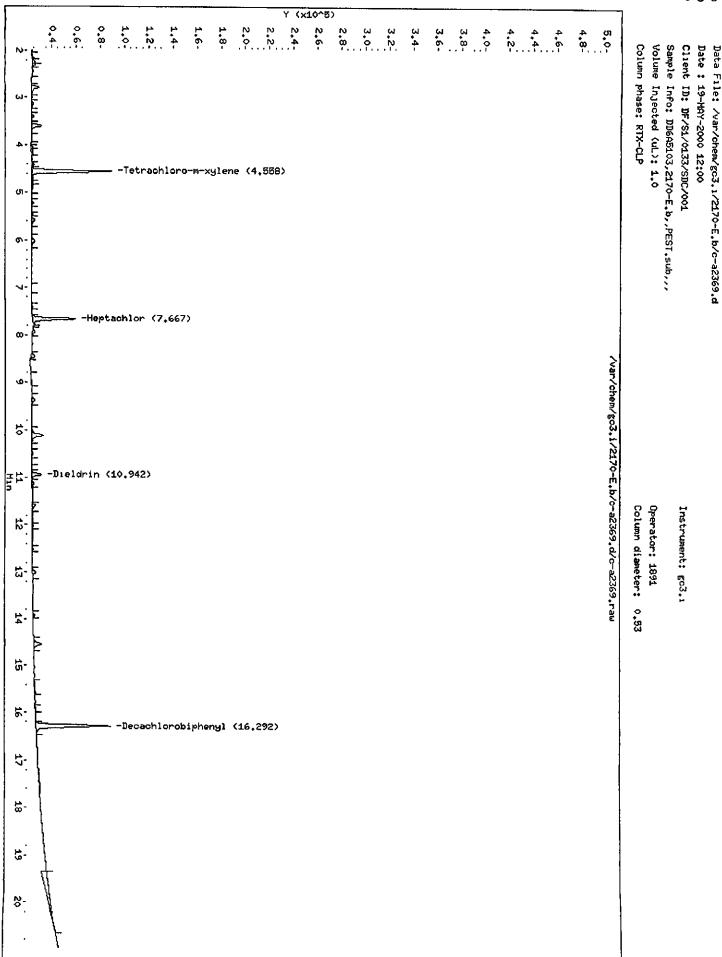
		ON-COLUMN FINAL
Compounds	RT EXP RT DLT RT RESPONSE	( ng) ( mg/L)
	共杂 四元公正司号 杂物质素素 医多种多种多种	
5 alpha-BHC	Compound Not Detected.	
4 HEXACHLOROBENZENE	Compound Not Detected.	
6 gamma-BHC (Lindane)	Compound Not Detected.	
7 beta-BHC	Compound Not Detected.	
9 Chlordane	Compound Not Detected.	
10 Heptachlor	7.667 7 675 -0 008 35494	0.0074 0.000740528(a)
8 delta-BHC	Compound Not Detected	
11 Aldrin	Compound Not Detected	
12 Heptachlor epoxide	Compound Not Detected.	
13 gamma-Chlordane	Compound Not Detected	
14 alpha-Chlordane	Compound Not Detected	
15 Endosulfan I	Compound Not Detected.	
16 4,4'-DDE	Compound Not Detected.	
17 Dieldrin	10 942 10.942 0.000 7967	0 00197 0 000196509(a)
20 Endrin	Compound Not Detected.	
18 Toxaphene	Compound Not Detected.	

Data File: /var/chem/gc3.i/2170-E.b/c-a2369.d Report Date: 19-May-2000 12:40

		CONCENTRATIONS		
		ON-COLUMN FINAL		
Compounds	RT EXP RT DLT RT RESPONSE	( ng) ( mg/L)		
平平产产 电电子电子电子电子电子电子 计计算电话 电线电		, ====================================		
21 4,4'-DDD	Compound Not Detected.			
22 Endosulfan II	Compound Not Detected.			
23 4,4'-DDT	Compound Not Detected.			
24 Endrin aldehyde	Compound Not Detected.			
26 Endosulfan sulfate	Compound Not Detected.			
25 Methoxychlor	Compound Not Detected.			
27 Endrin ketone	Compound Not Detected.			
<pre>\$ 1 Tetrachloro-m-xylene</pre>	4.558 4.558 0.000 66099	0.01632 0 00163247(aR)		
\$ 30 Decachlorobiphenyl	16 292 16.300 -0.008 61079	0.01854 0.00185406 (aR)		

# QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).
   R Spike/Surrogate failed recovery limits.



Data File: /var/chem/gc3.i/2170-E.b/c-b2369.d

Report Date: 19-May-2000 12:46

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2369.d

Client Smp ID: DF/S1/0133/SDC/001 Lab Smp Id: DD6A5103

Inj Date : 19-MAY-2000 12:00

Inst ID: gc3.i Operator : 1891

Smp Info : DD6A5103,2170-E.b,,PEST.sub,,,

Misc Info: 130194002

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTB.m Meth Date : 19-May-2000 11:24 eppinged Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-b2300 Cal File: c-b2300.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF Vt		Dilution Factor Volume of final extract (uL)
Vo	100.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected

#### CONCENTRATIONS

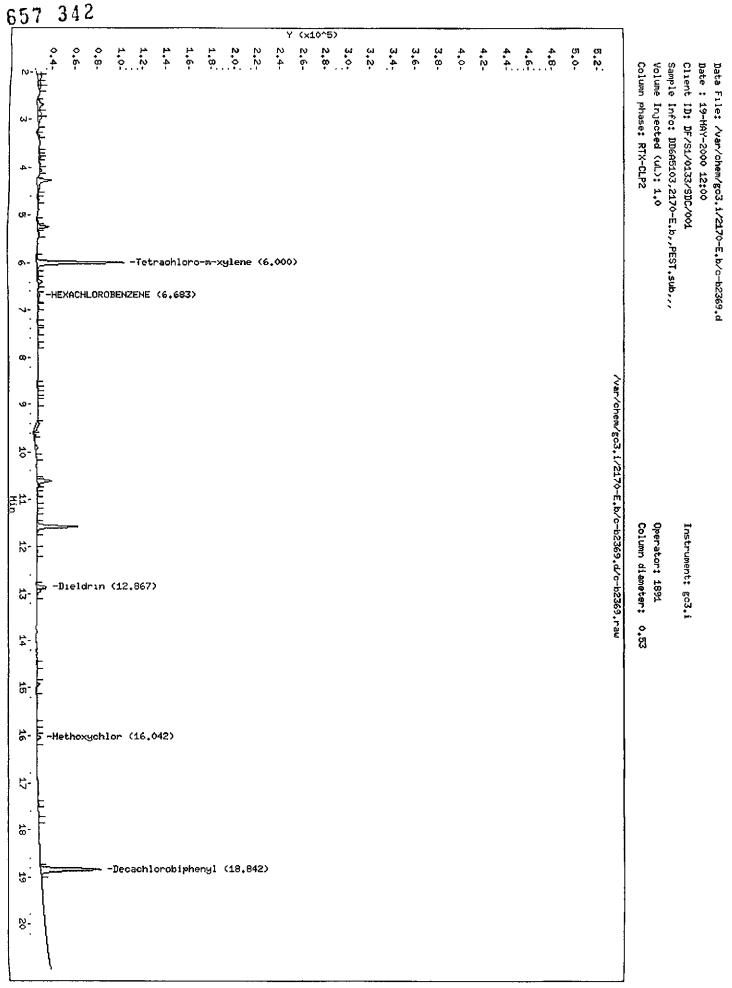
				ON-COLUMN	FINAL
Compounds	RT	EXP RT DLT RT	RESPONSE	( ng)	( mg/L)
*****	==	~~~~==	****	======	======
5 alpha-BHC	Com	pound Not Dete	cted.		
4 HEXACHLOROBENZENE	6.683	6.650 0.03	3 1730		(a)
6 gamma-BHC (Lindane)	Com	pound Not Dete	cted		
7 beta-BHC	Com	pound Not Dete	cted.		
9 Chlordane	Com	pound Not Dete	cted		
10 Heptachlor	Com	pound Not Dete	ected		
8 delta-BHC	Com	pound Not Deta	cted		
11 Aldrın	Com	pound Not Dete	cted		
12 Heptachlor epoxide	Com	pound Not Dete	cted.		
13 gamma-Chlordane	Com	pound Not Dete	cted.		
14 alpha-Chlordane	Com	pound Not Dete	cted		
15 Endosulfan I	Com	pound Not Dete	cted.		
16 4,4'-DDB	Com	pound Not Date	cted		
17 Dieldrin	12 867	12 867 0.00	00 8488	0 00204	0.000204452(a)
20 Endrin	Com	pound Not Dete	ected.		
18 Toxaphene	Com	pound Not Dete	cted.		

Data File: /var/chem/gc3.i/2170-E.b/c-b2369.d Report Date: 19-May-2000 12:46

		CONCENTRATIONS
		on-column final
Compounds	RT EXP RT DLT RT RESPONSE	( ng) ( mg/L)
21 4,4'-DDD	Compound Not Detected.	
22 Endosulfan II	Compound Not Detected.	
23 4,4'-DDT	Compound Not Detected.	
24 Endrin aldehyde	Compound Not Detected.	
26 Endosulfan sulfate	Compound Not Detected	
25 Methoxychlor	16 042 16.017 0.025 3517	0.00281 0.000281357(a)
27 Endrin ketone	Compound Not Detected	
\$ 1 Tetrachloro-m-xylene	6.000 6.000 D DOO 75446	0.0183 0.00183586 (aR)
\$ 30 Decachlorobiphenyl	18.842 18.842 0.000 52438	0.01939 0.00193854 (aR)

# QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).R - Spike/Surrogate failed recovery limits.



### UXB INTERNATIONAL

Lab Name: Severn Trent Laboratories, Inc. SDG Number

Matrix: (soil/water) SOUTH VARL Lab Sample ID COE130194 003

Method: SW846 8081A Um 5/03/00

Pesticides (8081A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A6103 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/19/00

QC Batch: 0139460

Client Sample Id: DF/S1/0133/SDC/001 C

### CONCENTRATION UNITS.

CAS NO.	COMPOUND (ug/L o	r ug/kg) mg/L 💢	<u> </u>
57-74-9	Chlordane (technical)	0.0050	<u>u</u>
72-20-8	Endrin	[0.00050 [	ן ט
76-44-8	Heptachlor	0.00050	וט
1024-57-3	Heptachlor epoxide	0.00050	<u> </u>
58-89-9	Lindane	0.00050	<u> </u>
72-43-5	Methoxychlor	0.0010	ן <u>ט</u>
8001-35-2	Toxaphene	0.020	ן ט

Data File: /var/chem/gc3.i/2170-E.b/c-a2370.d

Report Date: 19-May-2000 12:53

### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2370.d

Lab Smp Id: DD6A6103 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 19-MAY-2000 12:26

Operator : 1891 Inst ID: gc3.i

Smp Info : DD6A6103,2170-E.b,,PEST.sub,,,
Misc Info : 130194003

Comment

: /var/chem/gc3.i/2170-E.b/PESTA.m Method

Meth Date: 19-May-2000 11:01 eppinged Quant Type: ESTD Cal Date: 18-MAY-2000 01:57 Cal File: c-a230 Cal File: c-a2300.d

Als bottle: 1

Dil Factor: 1.00000 Integrator: Falcon

Compound Sublist: PEST.sub

Target Version: 3.40

# Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF	1.000	Dilution Factor
Vt	10.000	Volume of final extract (uL)
Vo	100.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected

COI	NCENTRA	TIONS
ON - C0	DLUMN	FINA

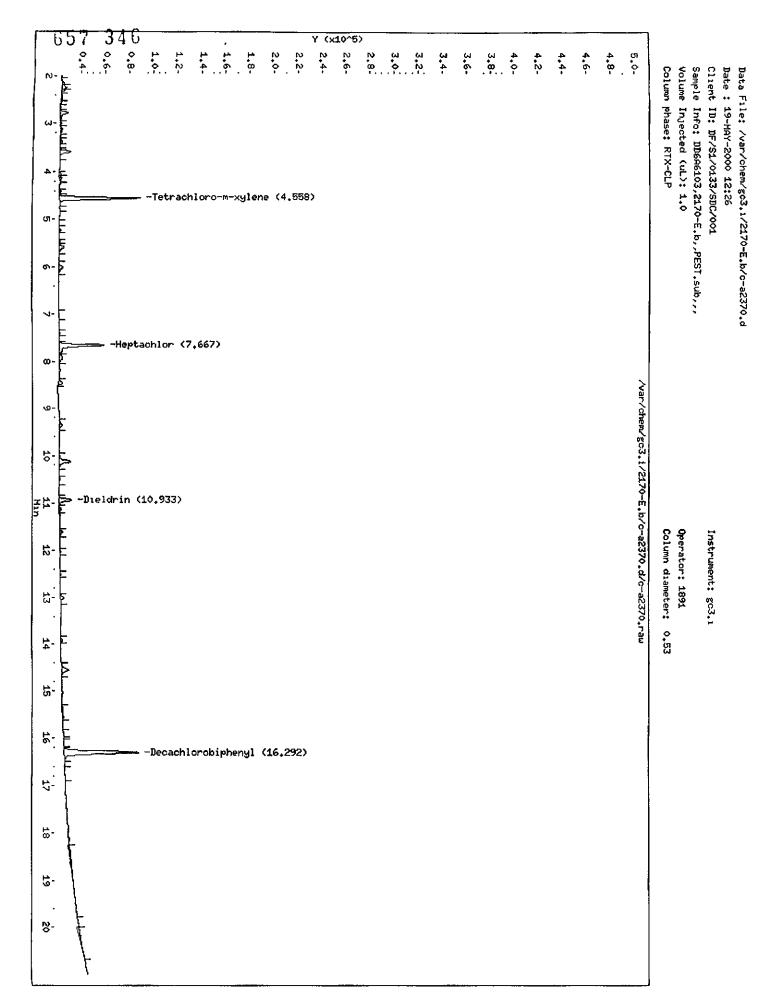
					ON-C	COLUMN	FINAL
Compounds	RT	EXP RT D	LT RT	RESPONSE	(	ng)	( mg/l)
	==	======================================		*******			<b>医医学</b> 耳草二二
5 alpha-BHC	Co	mpound Not	Detecte	1.			
4 HEXACHLOROBENZENE	Co	mpound Not	Detecte	i i			
6 gamma-BHC (Lindane)	Co	mpound Not	Detecte	i.			
7 beta-BHC	Co	mpound Not	Detecte	1			
9 Chlordane	Co	mpound Not	Detecte	i.			
10 Heptachlor	7.667	7 675	-0.008	36811	0 0	0768	0.000768005(a)
8 delta-BHC	Co	mpound Not	Detecte	i.			
11 Aldrin	Co	mpound Not	Detected	1.			
12 Heptachlor epoxide	Co	mpound Not	Detected	i.			
13 gamma-Chlordane	Co	mpound Not	Detected	i.			
14 alpha-Chlordane	Co	mpound Not	Detected	1.			
15 Endosulfan I	Co	mpound Not	Detected	i.			
16 4,4'-DDE	Co	mpound Not	Detecte	i.			
17 Dieldrin	10.933	10.942	-0.009	10116	0.0	0250	0.000249515(a)
20 Endrin	Co	mpound Not	Detected	i.			
18 Toxaphene	Co	mpound Not	Detecte	1.			

Data File: /var/chem/gc3.i/2170-E.b/c-a2370.d Report Date: 19-May-2000 12:53

		CONCENTRATIONS
Compounds	RT EXP RT DLT RT RESPONSE	ON-COLUMN FINAL ( ng) ( mg/L)
医多多性医院医院 化苯基苯基 医二苯基苯基 医二苯基苯基	医医 电自动性电影 医医过滤性 医医电阻性抗坏	
21 4,4'-DDD	Compound Not Detected	
22 Endosulfan II	Compound Not Detected.	
23 4,4'-DDT	Compound Not Detected.	
24 Endrin aldehyde	Compound Not Detected.	
26 Endosulfan sulfate	Compound Not Detected.	
25 Methoxychlor	Compound Not Detected.	
27 Endrin ketone	Compound Not Detected.	
\$ 1 Tetrachloro-m-xylene	4 558 4.558 0 000 66158	0.01634 (0.00163392 (aRL)
\$ 30 Decachlorobiphenyl	16 292 16.300 -0 008 61506	0.01867 0.00186703(aR)

# QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).
   R Spike/Surrogate failed recovery limits.



Data File: /var/chem/gc3.i/2170-E.b/c-b2370.d

Report Date: 19-May-2000 12:54

### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-b2370.d

Lab Smp Id: DD6A6103 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 19-MAY-2000 12:26

Operator: 1891 Inst ID: gc3.i

Smp Info : DD6A6103,2170-E.b,,PEST.sub,,,

Misc Info: 130194003

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date: 19-May-2000 11:24 eppinged Quant Type: ESTD Cal Date: 18-MAY-2000 01:57 Cal File: c-b2300.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF Vt Vo Vi	10.000 100.000	Dilution Factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected

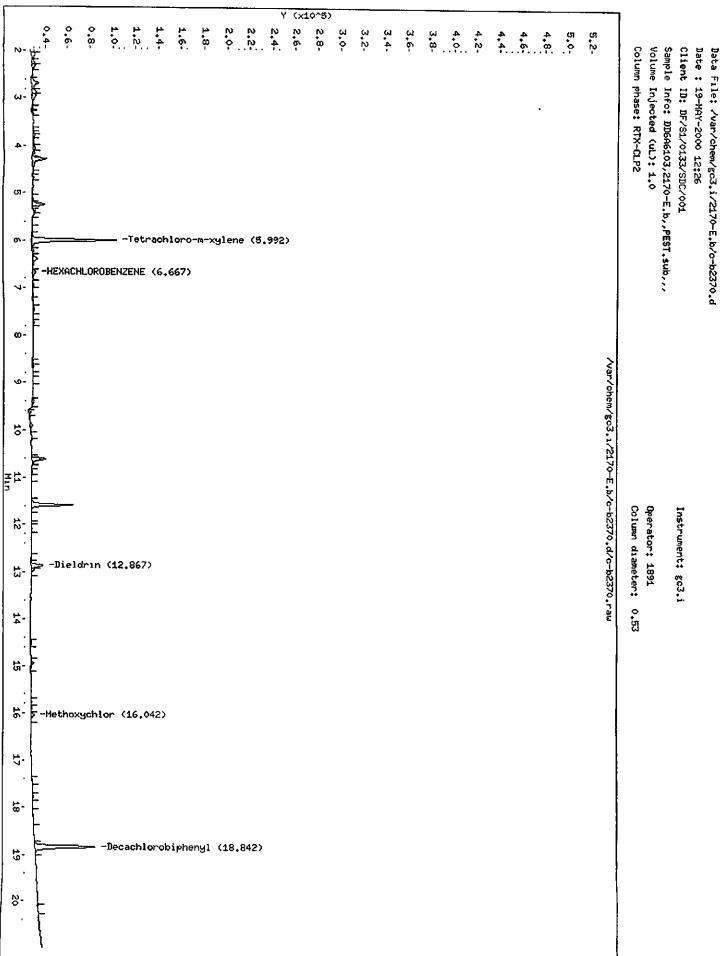
				CONCENTRATIONS						
							ON-	COLUMN	FINAL	
Compo	unds	RT	EXP I	er di	LT RT	RESPONSE	(	ng)	( mg/L)	
99E0E	中央共成四回电阻电影器器器器器器型型型计算符	**						****		•
s	alpha-BHC	Com	pound	Not	Detected	ì.				
4	HEXACHLOROBENZENE	6.667	6.6	50	0.017	2731				(a)
6	gamma-BHC (Lindane)	Com	pound	Not	Detected	<del>1</del>				
7	beta-BHC	Com	pound	Not	Detected	<del>i</del>				
9	Chlordane	Comp	pound	Not	Detected	i.				
10	Heptachlor	Com	pound	Not	Detected	i.				
8	delta-BHC	Com	pound	Not	Detected	3				
11	Aldrin	Com	pound	Not	Detected	<b>1.</b>				
12	Heptachlor epoxide	Com	pound	Not	Detected	à				
13	gamma-Chlordane	Com	pound	Not	Detected	d.				
14	alpha-Chlordane	Com	pound	Not	Detected	d				
15	Endosulfan I	Com	pound	Not	Detected	d.				
16	4,4'-DDB	Com	pound	Not	Detected	đ.				
17	Dieldrin	12.867	12 8	67	0 000	10703	0	00258	0 0002576	305 (a)
20	Bndrin	Com	pound	Not	Detected	d.				
18	Toxaphene	Com	pound	Not	Detected	d.				

657 348
Data File: /var/chem/gc3.i/2170-E.b/c-b2370.d
Report Date: 19-May-2000 12:54

		CONCENTRATIONS		
		on-column final		
Compounds	RT EXP RT DLT RT RESPONSE	( ng) ( mg/L)		
2 <b></b>	그는 토로프변화를 발달했다고로 무늬무실심장계약	BBBBBBB		
21 4,4'-DDD	Compound Not Detected			
22 Endosulfan II	Compound Not Detected			
23 4,4'-DDT	Compound Not Detected.			
24 Bndrin aldehyde	Compound Not Detected.			
26 Endosulfan sulfate	Compound Not Detected			
25 Methoxychlor	16 042 16.017 0.025 3165	0.00253 0.000253197(a)		
27 Endrin ketone	Compound Not Detected			
\$ 1 Tetrachloro-m-xylene	5.992 6 000 -0.008 74388	0 01810 0.00181012(aR)		
\$ 30 Decachlorobiphenyl	18.842 18 842 0.000 53240	0.01968 0.00196819 (aR)		

# QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).R Spike/Surrogate failed recovery limits.



#### UXB INTERNATIONAL

Lab Name · Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID-WATEL

Method: SW846 8081A Un \$3300

Pesticides (8081A)

Sample WT/Vol: 100 / mL

Work Order: DD6A7103

Date Received: 05/13/00

Date Extracted:05/18/00

Date Analyzed: 05/19/00

QC Batch · 0139460

Lab Sample ID COE130194 004

Client Sample Id: DF/S1/0133/SDC/001 D

### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L or	ug/kg) mg/L Q	
57-74-9	Chlordane (technical)	0.0050	<u>יי</u>
72-20-8	Endrin	0.00050	U
76-44-8	Heptachlor	0.00050	<u>U</u>
1024-57-3	Heptachlor epoxide	0.00050	U
58-89-9	Lindane	0.00050	U
72-43-5	Methoxychlor	0.0010	U
8001-35-2	Toxaphene	0.020	U

Data File: /var/chem/gc3.i/2170-E.b/c-a2371.d

Report Date: 19-May-2000 13:55

### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2371.d

Lab Smp Id: DD6A7103 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 19-MAY-2000 12:52

Operator: 1891 Inst ID: qc3.i

Smp Info : DD6A7193,2170-E.b,,PEST.sub,,,

Misc Info : 130192004

Comment: Up School

Method: /var/chem/gc3.i/2170-E.b/PESTA.m

Meth Date: 19-May-2000 11:01 eppinged Quar

Cal Date: 18-MAY-2000 01:57 Cal Quant Type: ESTD Cal File: c-a2300.d

Als bottle: 1

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF Vt Vo Vi	10.000 100.000	Dilution Factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected

CONCENTRATIONS

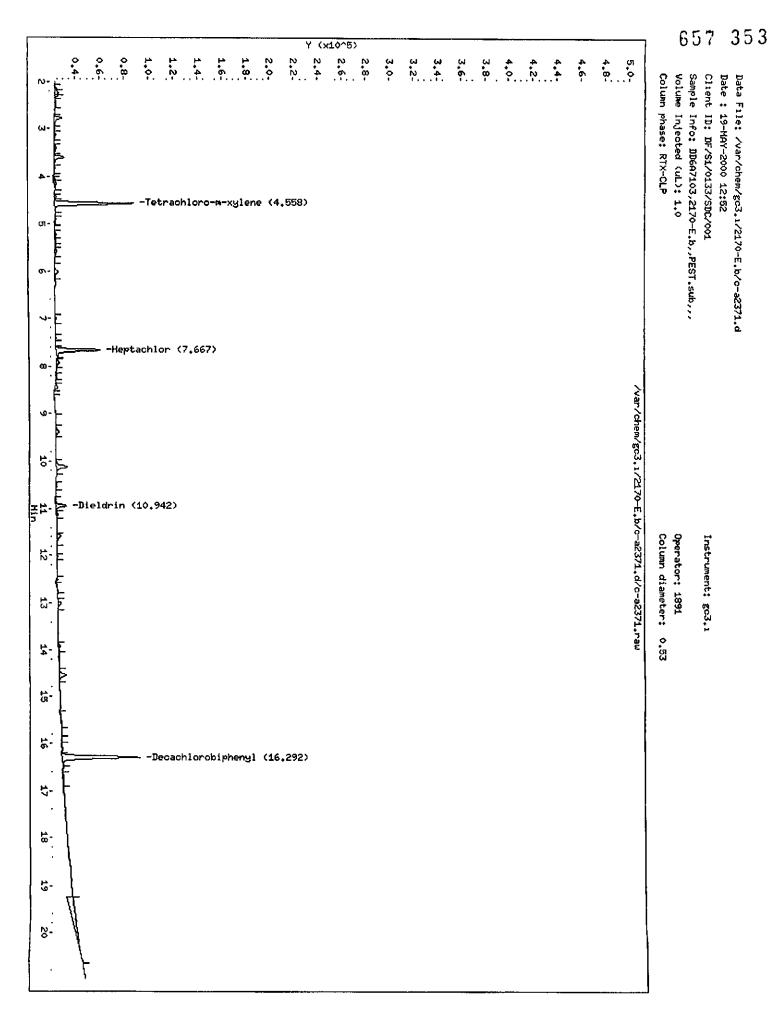
		CONCENTRATIONS						ATIONS
						ON-C	COLUMN	FINAL
Compo	unds	RT	EXP RT D	LT RT	RESPONSE	(	ng)	( mg/L)
					apapp323	<b>= n</b> :		
5	alpha-BHC	Comp	pound Not	Detected	1.			
4	HEXACHLOROBENZENE	Com	pound Not	Detected	i			
6	gamma-BHC (Lindane)	Com	pound Not	Detected	i.			
7	beta-BHC	Comp	pound Not	Detected	i.			
9	Chlordane	Comp	pound Not	Detected	î			
10	Heptachlor	7.667	7.675	-0.008	36639	0 0	00764	0.000764417(a)
8	delta-BHC	Comp	pound Not	Detected	<b>3</b> .			
11	Aldrin	Com	pound Not	Detected	<b>3</b> .		_	
12	Heptachlor epoxide	Com	pound Not	Detected	i.			
13	gamma-Chlordane	Comp	pound Not	Detected	à.			
14	alpha-Chlordane	Comp	pound Not	Detected	a.			
15	Endosulfan I	Comp	pound Not	Detected	<b>1.</b>			
16	4,4'-DDE	Comp	pound Not	Detected	đ.			
17	Dieldrin	10.942	10 942	0.000	8565	0.6	00211	0.000211259(a)
20	Endrin	Comp	pound Not	Detected	d.			
18	Toxaphene	Com	pound Not	Detected	d.			

Data File: /var/chem/gc3.i/2170-E.b/c-a2371.d Report Date: 19-May-2000 13:55

		CONCENTRATIONS	;
		ON-COLUMN FIN	IAL
Con	npounds	RT EXP RT DLT RT RESPONSE ( ng) ( mg	l\r)
===		カン ならかなこれ おかのからま まままれなどと スココニコエ のじょど	
	21 4,4'-DDD	Compound Not Detected	
	22 Endosulfan II	Compound Not Detected	
	23 4,4'-DDT	Compound Not Detected	
	24 Endrin aldehyde	Compound Not Detected.	
	26 Endosulfan sulfate	Compound Not Detected.	
	25 Methoxychlor	Compound Not Detected.	
	27 Endrin ketone	Compound Not Detected.	
\$	1 Tetrachloro-m-xylene	4.558 4 558 0 000 64866 0 01602 0.0016	0202 (aR)
\$	30 Decachlorobiphenyl	16.292 16.300 -0.008 64175 0 01948 0.0019	4804 (AR)

# QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).
   R Spike/Surrogate failed recovery limits.



Data File: /var/chem/gc3.i/2170-E.b/c-b2371.d

Report Date: 19-May-2000 13:58

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2371.d

Client Smp ID: DF/S1/0133/SDC/001 Lab Smp Id: DD6A7103

Inj Date : 19-MAY-2000 12:52

Inst ID: qc3.i Operator : 1891

Smp Info : DD6A7103,2170-E.b,,PEST.sub,,,

Misc Info: 130192004 054940

Method : /var/chem/gc3.i/2170-E.b/PESTB.m Meth Date : 19-May-2000 11:24 eppinged Quar Quant Type: ESTD Cal File: c-b2300.d Cal Date : 18-MAY-2000 01:57

Als bottle: 1

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF Vt Vo Vi	10.000 100.000	Dilution Factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected

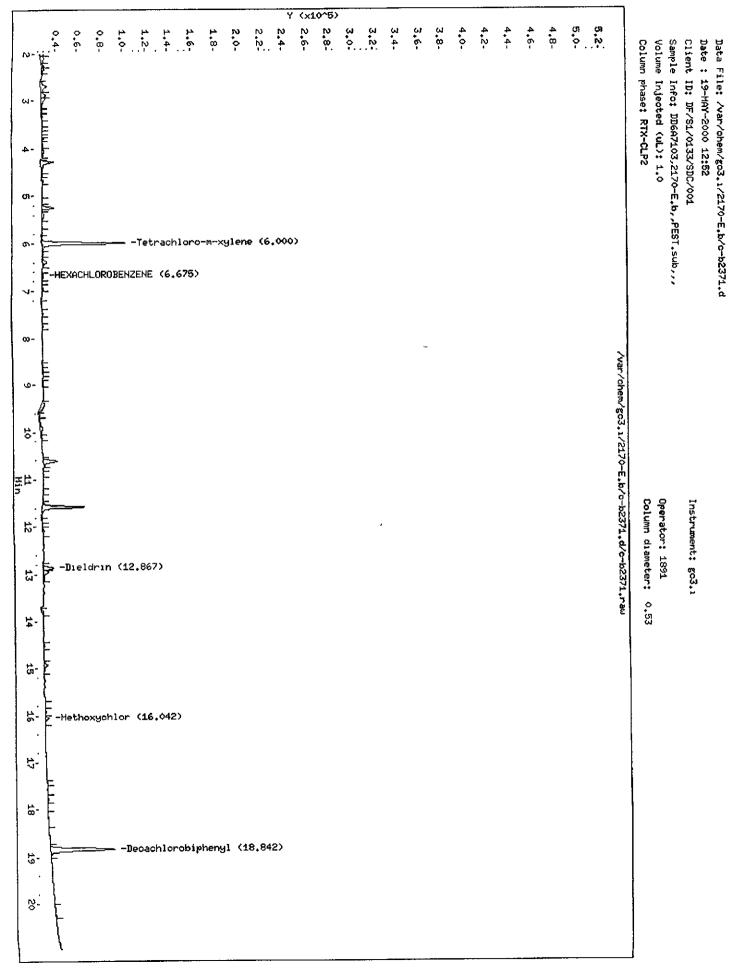
		CONCENTRATIONS
		on-column final
Compounds	RT EXP RT DLT RT RESPONSE	( ng) ( mg/L)
	## ####################################	# <b>====</b>
5 alpha-BHC	Compound Not Detected.	
4 HEXACHLOROBENZENE	6.675 6.650 0.025 782	(a)
6 gamma-BHC (Lindane)	Compound Not Detected.	
7 beta-BHC	Compound Not Detected.	
9 Chlordane	Compound Not Detected	
10 Heptachlor	Compound Not Detected.	
8 delta-BHC	Compound Not Detected	
11 Aldrin	Compound Not Detected.	
12 Heptachlor epoxide	Compound Not Detected	
13 gamma-Chlordane	Compound Not Detected.	
14 alpha-Chlordane	Compound Not Detected	
15 Endosulfan I	Compound Not Detected.	
16 4,4'-DDE	Compound Not Detected.	
17 Dieldrin	12.867 12.867 0 000 8880	0.00214 0 000213894(a)
20 Endrin	Compound Not Detected.	
18 Toxaphene	Compound Not Detected.	

Data File: /var/chem/gc3.i/2170-E.b/c-b2371.d Report Date: 19-May-2000 13:58

		CONCENTRATIONS		
		ON-COLUMN FINAL		
Compounds	RT EXP RT DLT RT RESPONSE	( ng) ( mg/L)		
21 4,4'-DDD	Compound Not Detected			
22 Endosulfan II	Compound Not Detected			
23 4,4'-DDT	Compound Not Detected.			
24 Endrin aldehyde	Compound Not Detected.			
26 Endosulfan sulfate	Compound Not Detected.			
25 Methoxychlor	16.042 16 017 0.025 3201	0.00256 0.000256077(a)		
27 Bndrin ketone	Compound Not Detected.			
\$ 1 Tetrachloro-m-xylene	6.000 6.000 0 000 73353	0.01785 0.00178493 (aR)		
\$ 30 Decachlorobiphenyl	18.842 18.842 0 000 56461	0 02087 0.00208726(aR)		

# QC Flag Legend

- a Target compound detected but, quantitated amount
   Below Limit Of Quantitation(BLOQ).
   R Spike/Surrogate failed recovery limits.



PESTICIDE CALIBRATION DATA

Report Date : 19-May-2000 11:01

### STL-PITTSBURGH

## COMPOUND LISTING

Method file : /var/chem/gc3.i/2170-E.b/PESTA.m
Quant Method : ESTD Tan
Last Update : 19-May-2000 11:00 Num
Data Type : GC MULTI COMP Target Version : 3.40 Number of Cpnds : 30

Global Integrator : Falcon

Chromat Events	Values
Initial:Start Threshold	40.00000
Initial:End Threshold	20.000000
Initial:Area Threshold	1000.000000
Initial:P-P Resolution	1.000000
Initial:Bunch Factor	1.00000
Initial:Negative Peaks	ON
Initial: Tension	0.00000

Compound	RT	RT Window	RF
\$ 1 Tetrachloro-m-xylene 2 Diallate A 3 Diallate B 4 HEXACHLOROBENZENE 5 alpha-BHC 6 gamma-BHC (Lindane) 7 beta-BHC 8 delta-BHC 9 Chlordane	4.558 5.725 6.033 6.650 5.883 6.625 6.908 7.258 7.492 7.675	5.983-6.083 6.600-6.700 5.833-5.933 6.575-6.675 6.858-6.958 7.208-7.308 7.442-7.542	4.049e+06 1.305e+05 3.835e+04 5.563e+06 5.103e+06 2.423e+06 4.700e+06 1.521e+05 2.832e+05 4.148e+05
10 Heptachlor 11 Aldrin 12 Heptachlor epoxide 13 gamma-Chlordane 14 alpha-Chlordane 15 Endosulfan I 16 4,4'-DDE 17 Dieldrin 18 Toxaphene	10.200 7.675 8.300 9.650 9.925 10.200 10.425 10.550 10.942 11.458	10.150-10.250 7.625-7.725 8.250-8.350 9.600-9.700 9.875-9.975 10.150-10.250 10.375-10.475 10.500-10.600 10.892-10.992 11.408-11.508 11.783-11.883 12.525-12.625	5.833e+05 4.793e+06 4.458e+06 4.174e+06 4.337e+06 4.289e+06 3.858e+06 3.642e+06 4.054e+06 5.625e+04 6.398e+04 7.043e+04

Report Date : 19-May-2000 11:01

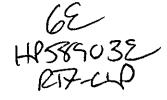
STL-PITTSBURGH

COMPOUND LISTING

Method file : /var/chem/gc3.i/2170-E.b/PESTA.m

Compound	RT	RT Window	RF
19 Isodrin 20 Endrin 21 4,4'-DDD 22 Endosulfan II 23 4,4'-DDT	11.767 11.850 12.317	11.325-11.425 11.717-11.817 11.800-11.900 12.267-12.367	2.918e+06 3.542e+06 2.772e+06
24 Endrin aldehyde 25 Methoxychlor 26 Endosulfan sulfate 27 Endrin ketone 28 Chlorobenzilate 29 Kepone \$ 30 Decachlorobiphenyl	13.450 13.583 14.083 11.425 11.867	12.667-12.767 13.400-13.500 13.533-13.633 14.033-14.133 11.375-11.475 11.817-11.917 16.250-16.350	1.393e+06 3.064e+06 3.451e+06 2.041e+06 3.543e+04

Report Date : 18-May-2000 10:47



STL-PITTSBURGH

# INITIAL CALIBRATION DATA

Start Cal Date : 17-MAY-2000 17:15 End Cal Date : 18-MAY-2000 01:57

Quant Method : ESTD
Origin : Disabled
Target Version : 3.40
Integrator : Falcon

Method file : /var/chem/gc3.i/2170-E.b/PESTA.m

Cal Date : 18-May-2000 10:47 eppinged

Curve Type : Average

Calibration File Names:

Level 1: /var/chem/gc3.i/2170-E.b/c-a2296.d Level 2: /var/chem/gc3.i/2170-E.b/c-a2297.d Level 3: /var/chem/gc3.i/2170-E.b/c-a2298.d Level 4: /var/chem/gc3.i/2170-E.b/c-a2299.d Level 5: /var/chem/gc3.i/2170-E.b/c-a2300.d

	0.00500	0 01000	0.02500	0.05000	0.10000	!	
Compound	Level 1	Level 2		Level 4		RRF	* RSD
		22722EEE					
2 Diallate A	180370		123042		104129	130546	23.837
3 Diallate B	50940	41328	36832	31827	30832	38352	21.373
4 HEXACHLOROBENZENE	+++++	+++++	+++++	+++++	+++++	++++	++++
5 alpha-BHC	6201600	5940500	5497520	5292360	4883520	5563100	9 -382
6 gamma-BHC (Lindane)	5748800	5480200	5068640	4821620	4394410	5102734	10/470
7 beta-BHC	2788800	2670100	2413960	2249220	1991370	2422690	13.23
8 delta-BHC	4790400	5047700	4731720	4637920	4291070	4699762	5.838
9 Chlordane(1)	1 +++++	+++++	152132	+++++	+++++	152132	0.00
(2)	1 +++++	+++++	283184	++++ [	++++	283184	0.00
(3)	+++++	+++++	414792	+++++ 1	+++++	414792	0.00
(4)	1 +++++	1 +++++	583336	+++++	+++++	583336	0.00
10 Heptachlor	5807600	5321900	4677120	4312460	3846250	4793066	16.32
11 Aldrin	4870200	4937500	4453960	4230020	3797330	4457802	10.57
12 Heptachlor epoxide	4770200	4701500	4116600	3865260	3414080	4173528	
13 gamma-Chlordane	4875800	4838500	4287960	4051200	3630790	4336850	12 22
14 alpha-Chlordane	4880800	4841000	4217680	3948900	3558280	4289332	13 34
15 Endosulfan I	4606400	4316300	3744800	3517560	3106900	3858392	
16 4.4'-DDE	3849800	3936500	3610960	3544360	3267940	3641912	7 26
17 Dieldrin	4622600	4444200	4014280	3785980	3404230	4054258	
18 Toxaphene(1)	+++++	+++++	56250	++++	+++++	56250	
(2)	1 +++++	+++++	63978	+++++	+++++	63978	
(3)	<b>++++</b>	+++++	70425	+++++	+++++	70425	
(4)	++++	+++++	73090	+++++	+++++	73090	
19 Isodrin	4271900	3746920	3640820	3194413	3155760	3601963	
20 Endrin	4185800	3758000	3460640	3294520	2947770	-	
21 4,4'-DDD	3119600	3110600	2854360	2853120	2651800	2917896	6.78

Report Date: 18-May-2000 10:47

STL-PITTSBURGH

#### INITIAL CALIBRATION DATA

Start Cal Date : 17-MAY-2000 17:15 End Cal Date : 18-MAY-2000 01:57

: ESTD : Disabled Quant Method Origin Target Version : 3.40

Integrator : Falcon
Method file : /var/chem/gc3.i/2170-E.b/PESTA.m
Cal Date : 18-May-2000 10:47 eppinged
Curve Type : Average

Curve Type

		i	0 00500	0.01000	0.02500	0.05000	0 10000	1	
Compos	und	1	Level 1	Level 2	Level 3	Level 4	Level 5	RRF	* RSD
	医医阴茎 经共享 医自己 化邻苯基甲基苯甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基	.===  #	=====					34444444 <b>(</b>	3366
22 En	dosulfan II	ì	3946200	3968400	3487600	3316420	2991890	3542102	11 824
23 4,	4 ' - DDT	1	2949600	2925300	2746560	2717480	2519290	2771646	6.315
24 En	drın aldehyde	1	3352600	3295100	2877080	2726800	2438460	2938008	13.152
25 Me	thoxychlor	1	1666900	1563400	1329900	1281770	1121910	1392776	15 808
26 En	dosulfan sulfate	- 1	3413800	3386800	2994040	2882520	2641020	3063636	10 866
27 En	drin ketone	}	3784000	3813300	3439400	3278340	2938130	3450634	10.600
28 Ch	lorobenzilate	1	1264800	2168120	1616060	2304987	2852540	2041301	30.273
29 Ke	pone	ļ	42040	38388	35702	31049	29991	35434	
	· 对于人名意西西拉拉 医巴萨氏征 计自由 医中心			_			******	_	
1 Te	etrachloro-m-xylene	1	5067000					•	
30 De	cachlorobiphenyl		4141400	3727100	3140920	2898000	2564230	3294330	19 309

Lab Name: STL-PITTSBURGH

Contract:

SAS No.: 40325 SDG No.: SDGA18632 Lab Code: STLPIT Case No.:

GC Column: RTX-CLP ID: 0.53 (mm) Init. Calib. Date(s): 05/17/00 05/18/00

Date Analyzed :\_\_\_\_ EPA Sample No. (PIBLK): \_\_\_\_\_

Time Analyzed : Lab Sample ID (PIBLK): \_\_\_\_\_

Date Analyzed :05/17/00 EPA Sample No. (PEM):

Time Analyzed: 1649 Lab Sample ID (PEM): EVALB

PEM COMPOUND	RT	RT WI FROM	MOOM TO	CALC AMOUNT (ng)	NOM AMOUNT (ng)	%D
Endrin 4,4'-DDT	11.37 12.31	11.32 12.27		0.024091 0.024457		

114

4,4'-DDT % breakdown (1):

11.80 Endrin % breakdown (1):

Combined % breakdown (1):

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.:

SAS No.: 40325 SDG No.: SDGA18632

GC Column: RTX-CLP ID: 0.53 (mm) Init. Calib. Date(s): 05/17/00 05/18/00

EPA Sample No. (PIBLK):

Date Analyzed :\_\_\_\_\_

Lab Sample ID (PIBLK):

Time Analyzed:

EPA Sample No. (PEM):

Date Analyzed:05/18/00

Lab Sample ID (PEM): EVALB

Time Analyzed:0315

PEM COMPOUND	RT	RT WI FROM	WOOM OT	CALC AMOUNT (ng)	NOM AMOUNT (ng)	%D
Endrin 4,4'-DDT	11.37			0.024288		

4,4'-DDT % breakdown (1):

Endrin % breakdown (1):

Combined % breakdown (1):

OLM03.0

# PESTICIDE CALIBRATION VERIFICATION SUMMARY

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.:

SAS No.: 40325 SDG No.: SDGA18632

GC Column: RTX-CLP ID: 0.53 (mm) Init. Calib. Date(s): 05/17/00 05/18/00

EPA Sample No. (PIBLK):

Date Analyzed :\_\_\_\_\_

Lab Sample ID (PIBLK):

Time Analyzed:

EPA Sample No. (PEM):

Date Analyzed:05/19/00

Lab Sample ID (PEM): EVALB

Time Analyzed: 0832

PEM. COMPOUND	RT	RT W	TO	CALC AMOUNT (ng)	NOM AMOUNT (ng)	%D
Endrin 4,4'-DDT	11.38 12.32	11.32 12.27			0.025000 0.025000	

4,4'-DDT % breakdown (1): 10:01 Endrin % breakdown (1):

Combined % breakdown (1):

FORM VII PEST-1

OLM03.0

Data File: /var/chem/gc3.i/2170-E.b/c-a2362.d Report Date: 19-May-2000 10:12

STL-PITTSBURGH

## CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i Lab File ID: c-a2362.d

Analysis Type:

Lab Sample ID: MEDTOX

Quant Type: ESTD

Injection Date: 19-MAY-2000 08:58

Init. Calibration Date(s): 05/17/0 05/18/0
Init. Calibration Times: 17:15 01:57
Method File: /var/chem/gc3.i/2170-E.b/PESTA.m

		1	'		MIN	1	XAM
	COMPOUND	1	RRF	RFO	RRF	♦D	<b>*</b> D ∫
-					-		
	18 Toxaphene(1)	- 1	56250.000	57622	000 0 010	-2 4	15 0
	(2)	ı	63978 000	65656	000 0 010	-2.6	15.0
	(3)	ŀ	70425.000	71716	000 0 010	-1.8]	15.0
	(4)	t	73090 000	75594	000 0 010	-3.4	15.0
\$	1 Tetrachloro-m-xylene	1	4049022 000	3869640	00010.0001	4.4	15.0
\$	30 Decachlorobiphenyl	1	3294330.000	3097480	000[0.010]	6.0	15.0
i		1	ı		1 1	1	

72 HP585035 657 366
Data File: /var/chem/gc3.i/2170-E.b/c-a2363.d
Report Date: 19-May-2000 10:12

## STL-PITTSBURGH

## CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i

Lab File ID: c-a2363.d

Analysis Type: Lab Sample ID: MEDCHLOR

Quant Type: ESTD

Injection Date: 19-MAY-2000 09:24 Init. Calibration Date(s): 05/17/0 05/18/0 Init. Calibration Times: 17:15 01:57 Method File: /var/chem/gc3.i/2170-E.b/PESTA.m

ı —					MIN	MAX
	COMPOUND	į	RRF	RFO	RRF	%D   %D
		=====]		**********		
1	9 Chlordane(1)	1	152132 000	156056.000	[0.010]	-2.6  15.0
	(2)	- 1	283184.000	288544.000	0.010	-1.9  15.0
ļ	(3)	- 1	414792.000	404980.000	0.010	2.4  15.0
ł	(4)	1	583336.000	573588 000	0.010	1.7[ 15.0]
\$	1 Tetrachloro-m-xylene	1	4049022.000	3956600 000	0.000	2.3  15.0
\$	30 Decachlorobiphenyl	- 1	3294330.000	3209280.000	0 010	2.6  15.0
1		1	ļ	l	.	1

72 HP58

657 367

Data File: /var/chem/gc3.i/2170-E.b/c-a2364.d

Report Date: 19-May-2000 10:12

## STL-PITTSBURGH

## CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i Lab File ID: c-a2364.d

Analysis Type:

Lab Sample ID: MEDA Quant Type: ESTD

Injection Date: 19-MAY-2000 09:50

Init. Calibration Date(s): 05/17/0 05/18/0
Init. Calibration Times: 17:15 01:57
Method File: /var/chem/gc3.i/2170-E.b/PESTA.m

 		1			MIN	I	MAX
•	COMPOUND	1	RRF	RF0	RRF	₹D	*D
				******	-		
\$	1 Tetrachloro-m-xylene	1 -	4049022.000	4129800.0	000.000	-2.0	15.0
I	5 alpha-BHC	1 :	5563100.000	5822840.0	00 0 010	-4 7	15.0
1	6 gamma-BHC (Lindane)	1	5102734.000	5249360.0	00 0.010	-2.9	15.0
1	10 Heptachlor		4793066.000	4849720.0	00 0.010	-1.2	15.0
1	15 Endosulfan I		3858392 000	3853200.0	00 0.010	0.1	15.0
	17 Dieldrin	1	4054258.000	4045560.0	00 0.010	0.2	15.0
ł	20 Endrin	- 1	3529346 000	3574560.0	00 0.010	-1.3	15.0
1	21 4,4°-DDD	- 1	2917896.000	2946240.0	00 0.010	-1.0	15.0
1	23 4,4'-DDT	- 1	2771646.000	2747880 0	000 0 010	0 9	15.0
1	25 Methoxychlor	1	1392776.000	1341360.0	000 0.010	3 7	15.0
\$	30 Decachlorobiphenyl	1	3294330 000	3125680 0	000 0.010	5 1	15.0
l_				l		1	1

STL Pittsburgh

3047

72 HP58903E

657ta3file: /var/chem/gc3.i/2170-E.b/c-a2365.d

Report Date: 19-May-2000 11:00

## STL-PITTSBURGH

#### CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i Lab File ID: c-a2365.d

Analysis Type:

Lab Sample ID: MEDB Quant Type: ESTD Injection Date: 19-MAY-2000 10:16

Init. Calibration Date(s): 05/17/0 05/18/0 Init. Calibration Times: 17:15 01:57 Method File: /var/chem/gc3.i/2170-E.b/PESTA.m

	ı	MIN	MAX
COMPOUND	RRF	RFO RRF	4D   4D
			======   ====
11 Aldrin	4457802 000	4491000.000[0.010	-0 7  15 0
7 beta-BHC	2422690.000	2373680 000 0 010	2 0  15 0
8 delta-BHC	4699762.000	4727200 000 0.010	-0.6  15 0
12 Heptachlor epoxide	4173528.000	4187720.000 0.010	-0.3  15.0
13 gamma-Chlordane	4336850 000	4266360.000 0.010	1.6  15.0
14 alpha-Chlordane	4289332 000	4193800.000 0.010	2 2 15.0
16 4,4'-DDE	3641912 000	3647520.000 0.010	-0 2  15.0
22 Endosulfan II	3542102.000	3533280.000 0 010	0 2   15.0
24 Endrin aldehyde	2938008.000	2911280 000 0 010	0 9  15 0
26 Endosulfan sulfate	3063636.000	3071920.000 0 010	-0 3  15 0
27 Endrin ketone	3450634.000	3486840 000 0.010	-1.0  15.0
	1	1	1 1

7E HP589038

Data File: /var/chem/gc3.i/2170-E.b/c-a2374.d Report Date: 19-May-2000 14:32

## STL-PITTSBURGH

## CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i Lab File ID: c-a2374.d

Analysis Type:

Lab Sample ID: MEDA Quant Type: ESTD

Injection Date: 19-MAY-2000 14:10 Init. Calibration Date(s): 05/17/0 05/18/0
Init. Calibration Times: 17:15 01:57
Method File: /var/chem/gc3.i/2170-E.b/PESTA.m

Ì	·		1	MIN		MAX	
1	COMPOUND	RRF	RF0	RRF	<b>%</b> D	<b>%</b> D	
ł	\$ 1 Tetrachloro-m-xylene	4049022.000	4176840 000	0.000	-3.2	15.0	
1	5 alpha-BHC	5563100.000	5999760 000	0.010	-7.8	15.0	
1	6 gamma-BHC (Lindane)	5102734.000	5376040.000	0 010	-5.4	15.0	
l	10 Heptachlor	4793066 000	4985280.000	0 010	-4.0	15.0	
-	15 Endosulfan I	3858392.000	3947720.000	0.010	-2 3	15.0	
į	17 Dieldrin	4054258.000	4134760.000	0.010	-2 0	15.0	
-	20 Endrin	3529346.000	3659160 000	0 010	-3.7	15.0	
-	21 4,4'-DDD	2917896 000	2880880.000	0 010	1 3	15.0	
١	23 4,4'-DDT	2771646.000	2867800.000]	0 010	-3.5	15 0	
I	25 Methoxychlor	1392776.000	1400940.000	0.010	-0 6	15.0	
-	\$ 30 Decachlorobiphenyl	3294330.000	3223680.000	0 010	2.1	15.0	
1	1		ı i	1		i	

Data File: /var/chem/gc3.i/2170-E.b/c-a2375.d

Report Date: 19-May-2000 15:20

STL-PITTSBURGH

# CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i Lab File ID: c-a2375.d

Analysis Type:

Lab Sample ID: MEDB Quant Type: ESTD

Injection Date: 19-MAY-2000 14:36 Init. Calibration Date(s): 05/17/0 05/18/0 Init. Calibration Times: 17:15 01:57 Method File: /var/chem/gc3.i/2170-E.b/PESTA.m

MAX MIN | RRF | %D | %D | COMPOUND RRF | 4457802.000| 4539520.000|0 010| -1 8| 15.0| 11 Aldrin 2422690.000 | 2397400.000 | 0.010 | 1.0 | 15.0 7 beta-BHC | 4699762 000 | 4854040.000 | 0.010 | -3.3 | 15.0 | 8 delta-BHC 4173528.000 4259040.000 0.010 -2.0 15.0 12 Heptachlor epoxide 4336850.000 4360040.000 0.010 -0.5 15.0 13 gamma-Chlordane

4289332.000 4314960.000 0.010 -0.6 15.0 14 alpha-Chlordane | 3641912 000 | 3672480.000 | 0.010 | -0.8 | 15.0 | 16 4,4'-DDE 3542102.000 3537920 000 0.010 0.1 15.0 22 Endosulfan II 2938008 000 2979720.000 0.010 -1 4 15.0 24 Endrin aldehyde | 3063636.000| 3094680 000|0.010| -1 0| 15.0| 26 Endosulfan sulfate | 3450634 000| 3532360 000|0 010| -2.4| 15.0| 27 Endrin ketone

Lab Name: STL-PITTSBURGH Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: C0E130194

GC Column: RTX-CLP ID: 0.53 (mm) Init. Calib. Date(s): 05/17/00 05/18/00

Instrument ID: GC3

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS, SAMPLES, AND STANDARDS IS GIVEN BELOW:

	MEAN SURROC TCX: 4.54	ATE RT FROM 1 DCB: 1	NITIAL CALI 16.27	BRATION		
	EPA	LAB	DATE	TIME	TCX	DCB
	SAMPLE NO.	SAMPLE ID	ANALYZED	ANALYZED	RT #	RT #
		=========		=========	=======	======
01		EVALB	05/17/00	1649	4.55	16.28
02		MEDTOX	05/17/00	1715	4.55	16.28
03		MEDCHLOR	05/17/00	1741	4.55	16.28
04	***	LOWAPPX9	05/17/00	1807		
05	<b>&gt;</b>	MLOWAPPX9	05/17/00	1833		
06	W38/W	MEDAPPX9	05/17/00	1859		
07		MHIGHAPPX9	05/17/00	1925		
08	_	HIGHAPPX9	05/17/00	1951		····
09		LOWA	05/17/00	2017	4.55	16.28
10		MILOWA	05/17/00	2230	4.56	16.29
11		MEDA	05/17/00	2256	4.56	16.28
12		MHIGHA	05/17/00	2322	4.56	16.28
13		HIGHA	05/17/00	2348	4.55	16.28
14		LOWB	05/18/00	0014		
15		MLOWB	05/18/00	0040	·	
16		MEDB	05/18/00	0106		
17		MHIGHB	05/18/00	0132	<u> </u>	
18		HIGHB	05/18/00	0157		
19		2NID A	05/18/00	0223	4.56	16.28
20		2ND B	05/18/00	0249		
21		EVALB	05/18/00	0315	4.56	16.28
22		EVALB	05/19/00	0832	4.56	16.29
23		MEDTOX	05/19/00	0858	4.56	16.29
24		MEDCHLOR	05/19/00	0924	4.56	16.29
25		MEDA	05/19/00	0950	4.56	16.30
26		MEDB	05/19/00	1016		
	DF/S1/0133/S		05/19/00	1042	4.56	16.29
	DF/S1/0133/S		05/19/00	1108	4.57	16.29
29	DF/S1/0133/S		05/19/00	1134	4.56	16.29
	DF/S1/0133/S		05/19/00	1200	4.56	16.29
	DF/S1/0133/S		05/19/00	1226	4.56	16.29
32	DF/S1/0133/S	טע6A7103	05/19/00	1252	4.56	16.29
]			<u> </u>		l <u></u>	·

QC LIMITS

TCX = Tetrachloro-m-xylene (+/- 0.05 MINUTES)
DCB = Decachlorobiphenyl (+/- 0.05 MINUTES)

\* Values outside of QC limits.

page 1 of 2

FORM VIII PEST

OLM03.0

<sup>#</sup> Column used to flag retention time values with an asterisk.

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT

Case No.:

SAS No.: 40325 SDG No.: C0E130194

GC Column: RTX-CLP ID: 0.53 (mm) Init. Calib. Date(s): 05/17/00 05/18/00

Instrument ID: GC3

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS, SAMPLES, AND STANDARDS IS GIVEN BELOW:

	MEAN SURROUTCX: 4.54	GATE RT FROM I DCB: 1	NITIAL CALI 16.27	BRATION		
	EPA SAMPLE NO.	LAB SAMPLE ID	DATE ANALYZED	TIME ANALYZED	TCX RT #	DCB RT #
01 02 03 04 05 06 07 08 09 10 11 12 13 14 15 16	PBLK	SAMPLE ID  ========  DDE2E101  DDE2E102  MEDA  MEDB	ANALYZED  ===================================	ANALYZED  ===================================		- 1
20 21 22						
23 24 25 26						
27 28 29 30 31						

QC LIMITS

TCX = Tetrachloro-m-xylene  $(+/-\tilde{0}.05 \text{ MINUTES})$ 

DCB = Decachlorobiphenyl

(+/- 0.05 MINUTES)

# Column used to flag retention time values with an asterisk.

\* Values outside of QC limits.

page 2 of 2

FORM VIII PEST

OLM03.0

Data File: /var/chem/gc3.i/2170-E.b/c-a2281.d

Report Date: 18-May-2000 10:59

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2281.d

Lab Smp Id: EVALB

Inj Date : 17-MAY-2000 16:49

Inst ID: gc3.i

Operator: 1891 In Smp Info: EVALB, 2170-E.b, , EVALBR. sub, , 3, 1

Misc Info: 190-88-8

Comment

: /var/chem/gc3.i/2170-E.b/PESTA.m Method

Meth Date : 18-May-2000 10:58 g Cal Date : 18-MAY-2000 01:57 Quant Type: ESTD Cal File: c-a2300.d

QC Sample: PEM Als bottle: 1

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: EVALBR.sub

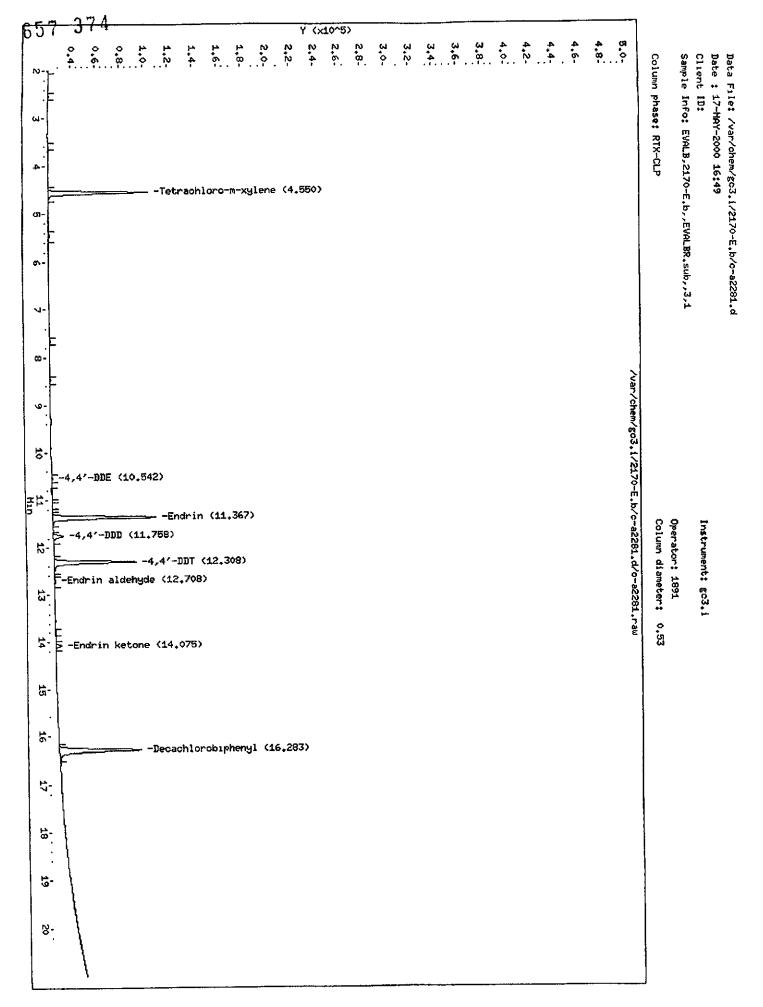
Target Version: 3.40

							С	ONCENT	RATIO	NS
							ON-	COLUMN	. F	INAL
Co	mpo	unds	RT	EXP RT	DLT RT	RESPONSE	(	ng)	(	ng)
	***	20. 20. 20. 20. 20. 20. 20. 20. 20. 20.	40.00						=0	
\$	1	Tetrachloro-m-xylene	4.550	4.558	-0.008	82259	0,	02032	0.02	03158(R)
	16	4,4'-DDE	10.542	10.542	0.000	717	0 0	00197	0.00	0196875
	20	Endrin	11 367	11.367	0.000	85027	٥.	02409	0.02	40914
	21	4,4'-DDD	11 758	11.758	0.000	8031	0	00275	0.00	275232
	23	4,4'-DDT	12.308	12.308	0.000	67787	Ο.	02446	0.02	44573
	24	Endrin aldehyde	12.708	12 708	0.000	1187	0 0	00404	0.00	0404015
	27	Endrin ketone	14 075	14 075	0.000	4767	٥.	00138	0.00	138148
\$	30	Decachlorobiphenyl	16.283	16.283	0.000	66957	٥.	02032	0.02	03249(R)

QC Flag Legend

R - Spike/Surrogate failed recovery limits.

OD 11:470 ENDRIN 6:570



Data File: /var/chem/gc3.i/2170-E.b/c-a2282.d

Report Date: 18-May-2000 10:48

## STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2282.d

Lab Smp Id: MEDTOX

Inj Date : 17-MAY-2000 17:15

Operator : 1891 Inst ID: gc3.i

Smp Info : MEDTOX,2170-E.b,,1-TOX.sub,,1,3
Misc Info : 190-84-13

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 18-May-2000 10:48 eppinged Quar Cal Date : 17-MAY-2000 17:15 Cal Quant Type: ESTD Cal File: c-a2282.d

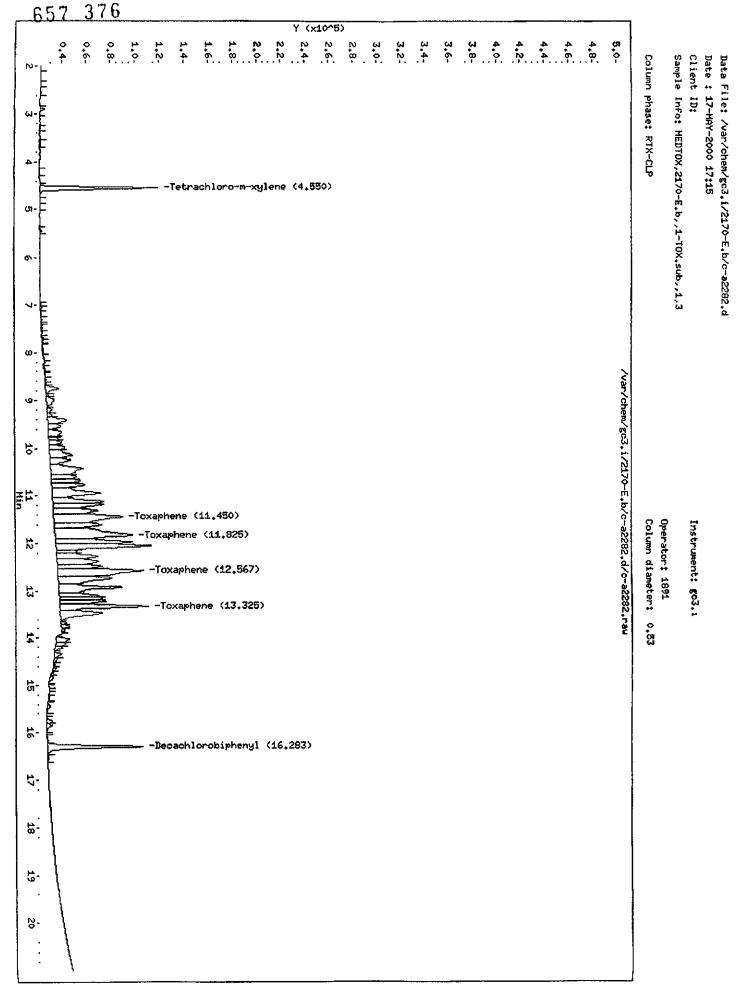
Als bottle: 1 Calibration Sample, Level: 3

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 1-TOX.sub

Target Version: 3.40

				AMOU	NTS
				CAL-AMT	ON-COL
Compounds	RT	EXP RT DLT RT	RESPONSE	( ng)	( ng)
	***	****** ****		*****	
18 Toxaphene	11.450	11.450 0 000	56250	1.00000	1.00000
5 1 Tetrachloro-m-xylene	4.550	4 558 -0.008	97585	0.02500	0 0250000
\$ 30 Decachlorobiphenyl	16.283	16.283 0.000	78886	0.02500	0 0250000
* ** **********************************				******	



Data File: /var/chem/gc3.i/2170-E.b/c-a2283.d

Report Date: 18-May-2000 10:48

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2283.d

Lab Smp Id: MEDCHLOR

Inj Date : 17-MAY-2000 17:41

Inst ID: gc3.i Operator : 1891

Smp Info : MEDCHLOR, 2170-E.b, , 2-CHLO.sub, , 1, 3

Misc Info : 190-85-10

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 18-May-2000 10:48 eppinged Quar Cal Date : 17-MAY-2000 17:41 Cal Quant Type: ESTD Cal File: c-a2283.d

Calibration Sample, Level: 3 Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: 2-CHLO.sub Integrator: Falcon

Target Version: 3.40

					AMOU	nts
					CAL-AMT	ON-COL
Compounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
李章里是是这种不同时代与李章的心理的	**		38888	**=====	*=====	======
9 Chlordane	7.483	7 483	0.000	38033	0.25000	0.250000
\$ 1 Tetrachloro-m-xylene	4.550	4.558	-0.008	97514	0.02500	0.0250000
s 30 Decachlorobiohenvi	16.283	16.283	0.000	79795	0.02500	0 0250000

Data File: /var/chem/gc3.i/2170-E.b/c-a2289.d

Report Date: 18-May-2000 10:49

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-a2289.d

Lab Smp Id: LOWA

Inj Date : 17-MAY-2000 20:17 Operator : 1891 In Smp Info : LOWA,2170-E.b,,3-INDA.sub,,1,1 Inst ID: gc3.i

Misc Info: 190-84-1

Comment

Method: /var/chem/gc3.i/2170-E.b/PESTA.m

Meth Date: 18-May-2000 10:48 eppinged Quant Type: ESTD

Cal Date: 17-MAY-2000 20:17 Cal File: c-a2289

Als bottle: 1 Calibration Samp Cal File: c-a2289.d

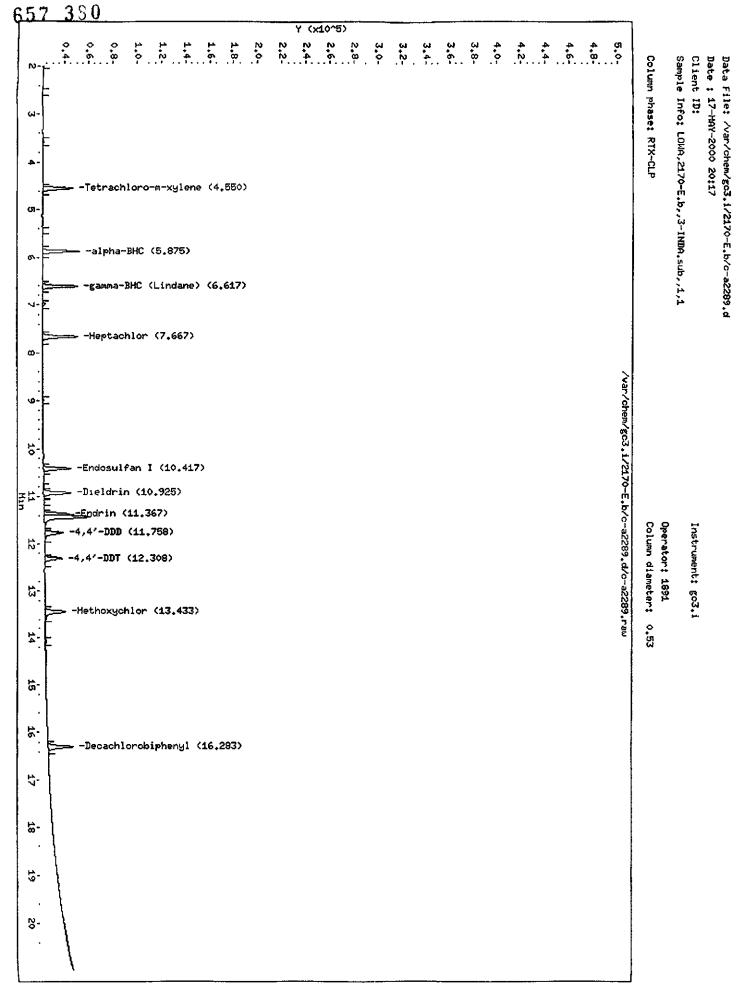
Calibration Sample, Level: 1

Dil Factor: 1.00000

Compound Sublist: 3-INDA.sub Integrator: Falcon

Target Version: 3.40

					AMOU	NTS
					CAL-AMT	ON-COL
Compounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
医乳球性动脉 医克朗氏 医克朗氏 医克朗氏 医克朗氏 计	==		**====	********	2424	
\$ 1 Tetrachloro-m-xylene	4.550	4.558	-0.008	25335	0.00500	0 00\$65037
5 alpha-BHC	5.87 <b>5</b>	5.883	-0.008	31008	0.00500	0 00500000
6 gamma-BHC (Lindane)	6.617	6.617	0 000	28744	0.00500	0.00500000
10 Heptachlor	7.667	7.675	-0.008	29038	0.00500	0 00500000
15 Endosulfan I	10.417	10 425	-0.008	23032	0.00500	0.00500000
17 Dieldrin	10.925	10.933	-0.008	23113	0.00500	0.00500000
20 Endrin	11 367	11 367	0.000	20929	0 00500	0 00500000
21 4,4'-DDD	11 758	11.758	0 000	15598	0.00500	0.00500000
23 4,4'-DDT	12 308	12.308	0 000	14748	0.00500	0.00500000
25 Methoxychlor	13.433	13.442	-0.009	16669	0.01000	0 0100000
\$ 30 Decachlorobiphenyl	16 283	16.283	0.000	20707	0.00500	0.00564747



Data File: /var/chem/gc3.i/2170-E.b/c-a2292.d

Report Date: 18-May-2000 10:49

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2292.d

Lab Smp Id: MLOWA

Inj Date : 17-MAY-2000 22:30 Operator : 1891 Ins Smp Info : MLOWA, 2170-E.b,,3-INDA.sub,,1,2 Inst ID: gc3.i

Misc Info: 190-84-2

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 18-May-2000 10:48 eppinged Quar Cal Date : 17-MAY-2000 22:30 Cal Quant Type: ESTD Cal File: c-a2292.d

Als bottle: 1 Calibration Sample, Level: 2

Dil Factor: 1.00000

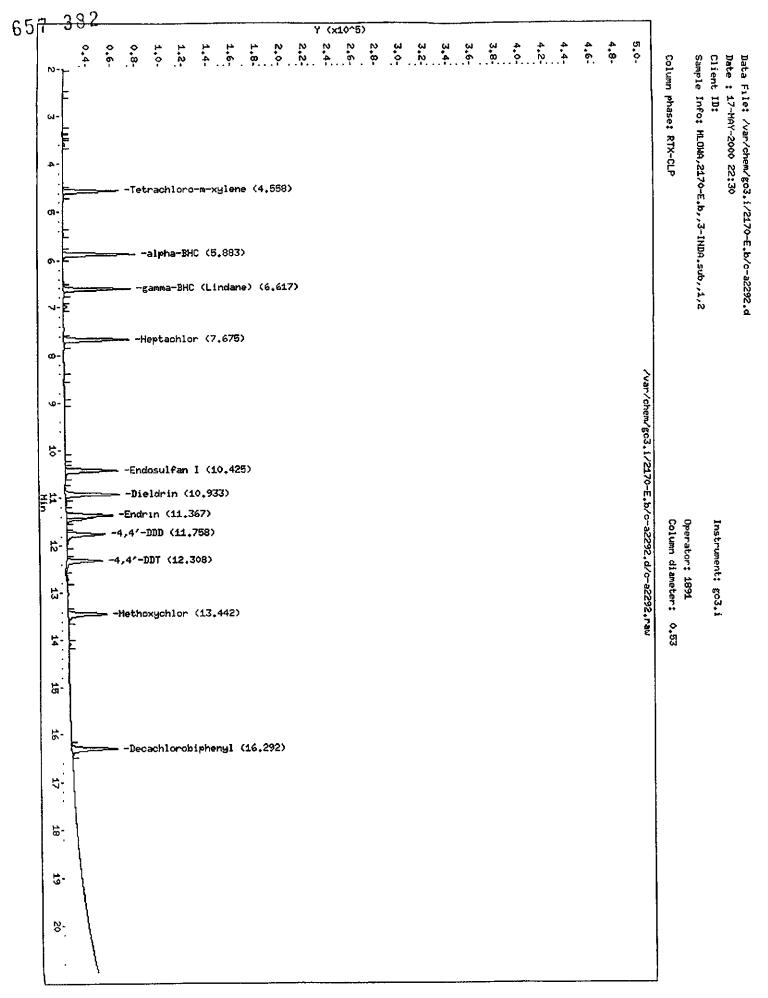
Integrator: Falcon Compound Sublist: 3-INDA.sub

Target Version: 3.40

						AMOU	nts
						CAL-AMT	ON-COL
Co	ompounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	(ng)
	· 完全要要 双耳角 医多角性 医 经过 经可有要集 首 医 医 基 经	••					
\$	1 Tetrachloro-m-xylene	4.558	4.558	0.000	45629	0.01000	0 0101170
	5 alpha-BHC	5.883	5.883	0.000	59405	0.01000	0.00978496
	6 gamma-BHC (Lindane)	6.617	6.617	0 000	54802	0.01000	0.00976080
	10 Heptachlor	7.675	7.675	0 000	53219	0.01000	0.00956359
	15 Endosulfan I	10.425	10.425	0 000	43163	0.01000	0.00967487
	17 Dieldrin	10.933	10.933	0 000	44442	0 01000	0 00980324
	20 Endrin	11.367	11.367	0 000	37580	0.01000	0.00946147(M)
	21 4,4'-DDD	11.758	11.758	0.000	31106	0.01000	0.00998555
	23 4,4'-DDT	12.308	12 308	0 000	29253	0.01000	0.00995864
	25 Methoxychlor	13 442	13.442	0.000	31268	0.02000	0 0193592
Ş	30 Decachlorobiphenyl	16.292	16.283	0.009	37271	0.01000	0.0101094

## QC Flag Legend

M - Compound response manually integrated.



Data File: /var/chem/gc3.i/2170-E.b/c-a2293.d

Report Date: 18-May-2000 10:49

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2293.d

Lab Smp Id: MEDA

Inj Date : 17-MAY-2000 22:56

Operator: 1891 In Smp Info: MEDA,2170-E.b,,3-INDA.sub,,1,3 Inst ID: qc3.i

Misc Info : 190-84-3

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m

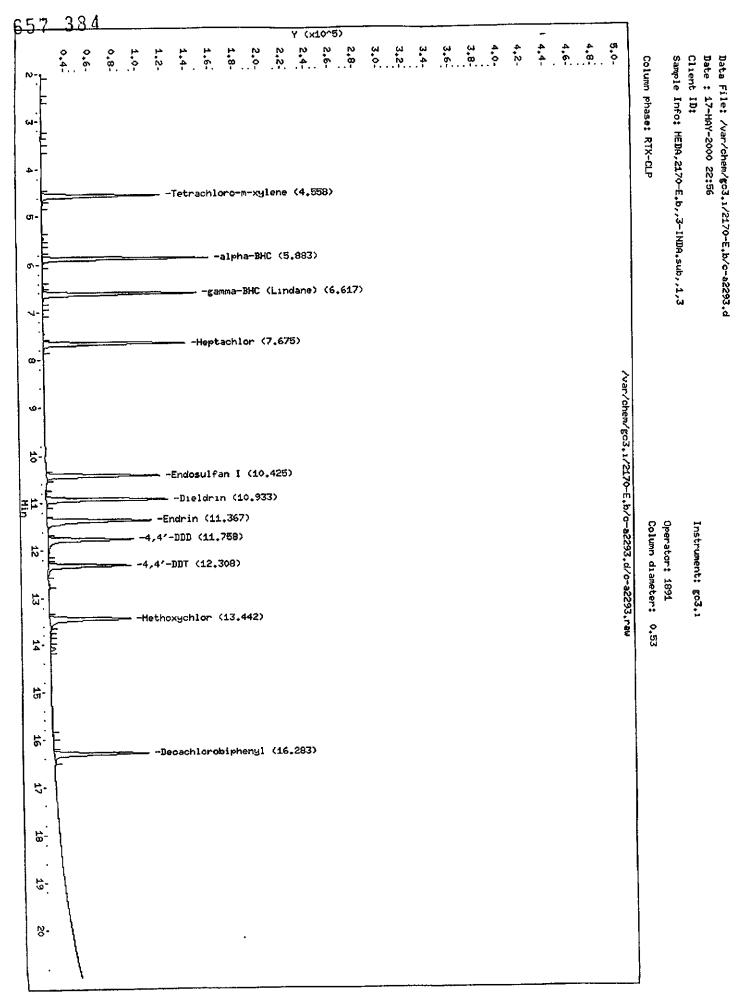
Meth Date: 18-May-2000 10:48 eppinged Quant Type: ESTD Cal File: c-a2293.d

Cal Date : 17-MAY-2000 22:56 Als bottle: 1 Dil Factor: 1.00000 Calibration Sample, Level: 3

Integrator: Falcon Compound Sublist: 3-INDA.sub

Target Version: 3.40

						AMOU	nts
						CAL-AMT	ON-COL
Co	mpounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
==	自由 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2				4588848	******	
\$	1 Tetrachloro-m-xylene	4.558	4.558	0 000	98378	0.02500	0 0217570
	5 alpha-BHC	5.883	5.883	0.000	137438	0.02500	0 0233743
	6 gamma-BHC (Lindane)	6.617	6.617	0.000	126716	0.02500	0 0233253
	10 Heptachlor	7.675	7.675	0.000	116928	0.02500	0 0221922
	15 Endosulfan I	10.425	10.425	0.000	93620	0.02500	0.0221717
	17 Dieldrin	10.933	10.933	0 000	100357	0 02500	0.0230158
	20 Endrin	11.367	11.367	0.000	86516	0 02500	0 0227585
	21 4,4'-DDD	11.758	11.758	0.000	71359	0.02500	0.0235649
	23 4.4'-DDT	12.308	12 308	0 000	68664	0 02500	0.0238929
	25 Methoxychlor	13 442	13.442	0 000	66495	0.05000	0.0437448
\$	30 Decachlorobiphenyl	16.283	16,283	0.000	78523	0.02500	0.0213970



Data File: /var/chem/gc3.i/2170-E.b/c-a2294.d

Report Date: 18-May-2000 10:49

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-a2294.d

Lab Smp Id: MHIGHA

Inst ID: gc3.i

Inj Date : 17-MAY-2000 23:22 Operator : 1891 Inst Smp Info : MHIGHA,2170-E.b,,3-INDA.sub,,1,4

Misc Info : 190-84-4

Comment

: /var/chem/gc3.i/2170-E.b/PESTA.m Method

Meth Date: 18-May-2000 10:48 eppinged Quant Type: ESTD Cal Date: 17-MAY-2000 23:22 Cal File: c-a2294.d

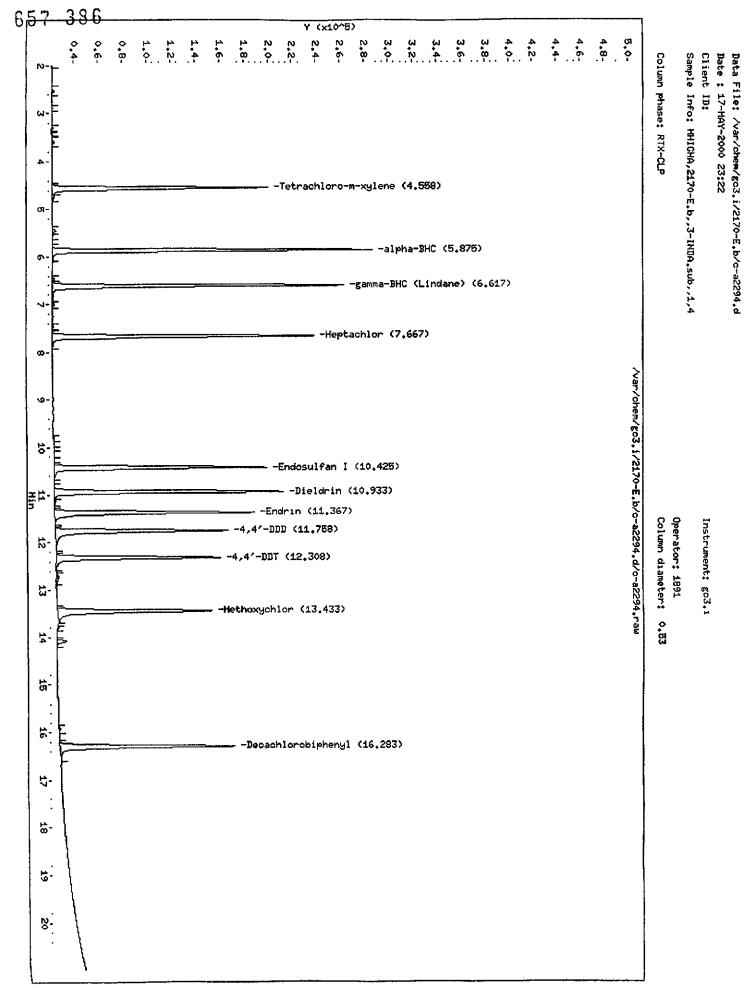
Als bottle: 1 Calibration Sample, Level: 4

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 3-INDA.sub

Target Version: 3.40

							AMOU	NTS
							CAL-AMT	ON-COL
Co	mpo	unds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
	2 <b>2 2</b>	3 2 2 2 2 2 2 2 2 2 2 2 3 3 3 3 3 2 2 2 2 2 2 3 3 3 3 3 2 2 2 2 2 2 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3		2 2 2 2 2 2 3 A	***	****		
\$	1	Tetrachloro-m-xylene	4.558	4,558	0.000	178492	0.05000	0.0416676
	5	alpha-BHC	5.875	5 883	-0 008	264618	0.05000	0.0461570
	6	gamma-BHC (Lindane)	6 617	6.617	0.000	241081	0 05000	0.0456609
	10	Heptachlor	7.667	7.675	-0.008	215623	0 05000	0.0428694
	15	Endosulfan I	10.425	10.425	0.000	175878	0 05000	0.0434668
	17	Dieldrin	10 933	10.933	0.000	189299	0.05000	0.0448920
	20	Endrin	11.367	11.367	0 000	164726	0.05000	0 0448266
	21	4,4'-DDD	11.758	11 758	0.000	142656	0.05000	0.0478002
	23	4,4'-DDT	12.308	12.308	0.000	135874	0.05000	0.0479318
	25	Methoxychlor	13.433	13.442	-0.009	128177	0.10000	0.0877629
\$	3.0	Decachlorobiphenyl	16.283	16.283	0.000	144900	0.05000	0.0416756



Data File: /var/chem/gc3.i/2170-E.b/c-a2295.d

Report Date: 18-May-2000 10:49

## STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2295.d

Lab Smp Id: HIGHA

Inj Date : 17-MAY-2000 23:48

Operator: 1891 Ins Smp Info: HIGHA, 2170-E.b,,3-INDA.sub,,1,5 Inst ID: gc3.i

Misc Info: 190-84-5

Comment

: /var/chem/gc3.i/2170-E.b/PESTA.m Method

Meth Date: 18-May-2000 10:48 eppinged Quant Type: ESTD Cal File: c-a2295.d Cal Date : 17-MAY-2000 23:48

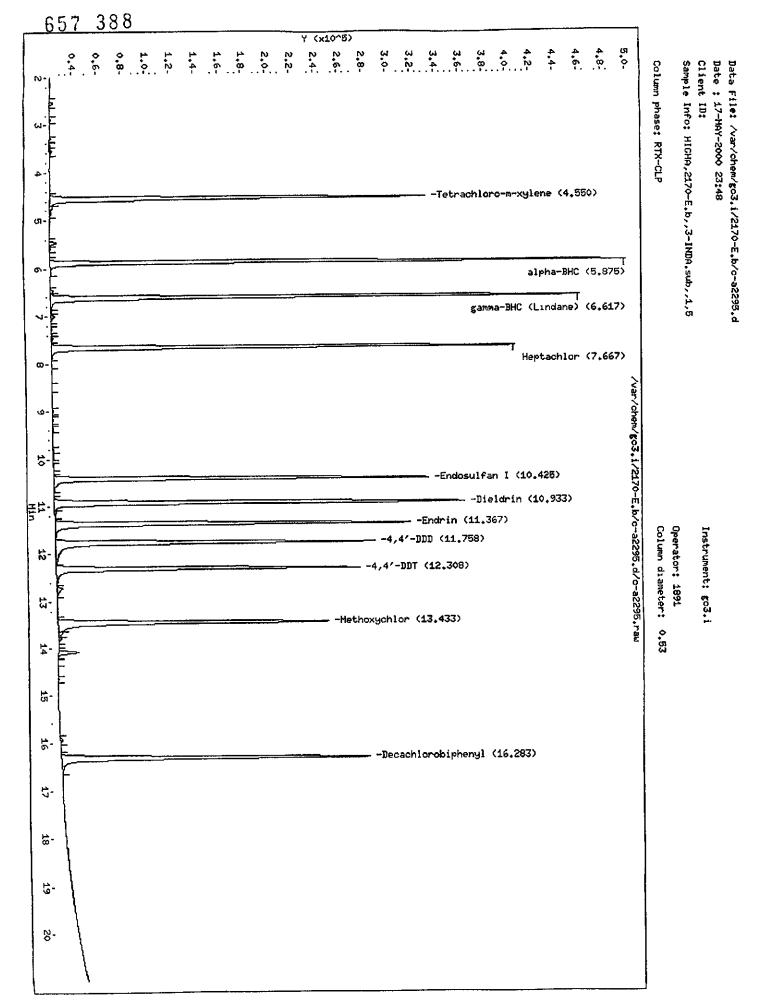
Calibration Sample, Level: 5

Als bottle: 1
Dil Factor: 1.00000
Integrator: Falcon

Compound Sublist: 3-INDA.sub

Target Version: 3.40

						AMOU	NTS
						CAL-AMT	ON-COL
Co	mpounds	ŔŦ	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
		==			*****	£ 0 0 0 0 0 0 0	EMMODER
\$	1 Tetrachloro-m-xylene	4.550	4.558	-0.008	311025	0.10000	0.0768148
	5 alpha-BHC	5 875	5 883	-0.008	488352	0.10000	0.0877841
	6 gamma-BHC (Lindane)	6.617	6.617	0.000	439441	0.10000	0.0861187
	10 Heptachlor	7.667	7 675	-0 008	384625	0.10000	0.0802461
	15 Endosulfan I	10.425	10.425	0.000	310690	0.10000	0 0805232
	17 Dieldrin	10.933	10 933	0.000	340423	0.10000	0.0839668
	20 Endrin	11.367	11 367	0 000	294777	0.10000	0 0835217
	21 4,4'-DDD	11 758	11.758	0.000	265180	0.10000	0.0908806
	23 4,4'-DDT	12.308	12.308	0.000	251929	0 10000	0.0908951
	25 Methoxychlor	13.433	13.442	-0 009	224382	0.20000	0.161104
s	30 Decachlorobiphenvl	16.283	16 283	0 000	256423	0 10000	0.0778377



Data File: /var/chem/gc3.i/2170-E.b/c-a2296.d

Report Date: 18-May-2000 10:49

## STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-a2296.d

Lab Smp Id: LOWB

Inj Date : 18-MAY-2000 00:14

Operator: 1891 In Smp Info: LOWB,2170-E.b,,4-INDB.sub,,1,1 Misc Info: 190-84-7 Inst ID: gc3.i

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 18-May-2000 10:48 eppinged Quant Type: ESTD Cal File: c-a2296.d Cal Date : 18-MAY-2000 00:14

Calibration Sample, Level: 1 Als bottle: 1

Dil Factor: 1.00000 Integrator: Falcon

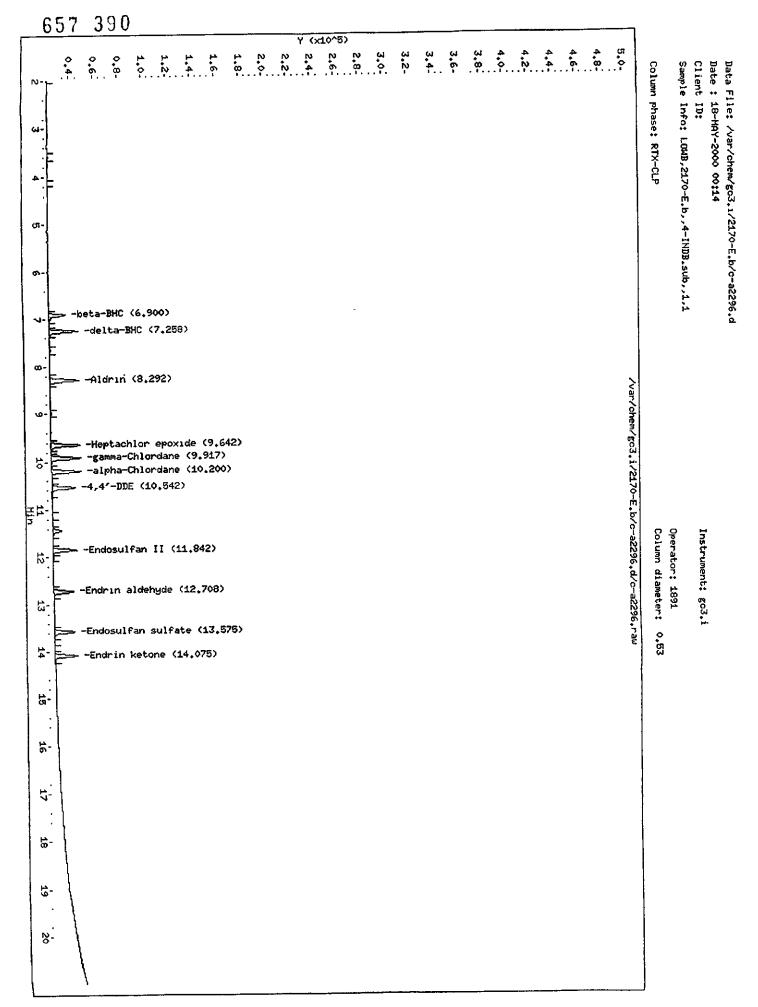
Compound Sublist: 4-INDB.sub

Target Version: 3.40

					AMOU	nts
					CAL-AMT	ON-COL
Compounds	RT	EXP RT	DLT RT	RESPONSE	(ng)	( ng)
· · · · · · · · · · · · · · · · · · ·	**		*****	4440-046	50×7000	
11 Aldrin	8.292	8.292	0.000	24351	0.00500	0.00500000
7 beta-BHC	6 900	6.900	0.000	13944	0 00500	0.00500000
8 delta-BHC	7.258	7.258	0 000	23952	0.00500	0.00500000
12 Heptachlor epoxide	9.642	9 642	0.000	23851	0 00500	0.00500000
13 gamma-Chlordane	9.917	9.917	0.000	24379	0 00500	0.00500000
14 alpha-Chlordane	10.200	10.200	0.000	24404	0.00500	0.00500000
16 4,4'-DDE	10 542	10.542	0.000	19249	0.00500	0.00500000
22 Endosulfan II	11.842	11.842	0.000	19731	0.00500	0.00500000
24 Endrin aldehyde	12,708	12.708	0.000	16763	0.00500	0.00500000
26 Endosulfan sulfate	13 575	13.575	0 000	17069	0.00500	0.00500000
27 Endrin ketone	14.075	14.075	0.000	18920	0 00500	0.00500000

STL Pittsburgh 3069

.--.



Data File: /var/chem/gc3.i/2170-E.b/c-a2297.d

Report Date: 18-May-2000 10:49

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-a2297.d

Lab Smp Id: MLOWB

Inj Date : 18-MAY-2000 00:40

Inst ID: gc3.i Operator : 1891

Smp Info : MLOWB,2170-E.b,,4-INDB.sub,,1,2
Misc Info : 190-84-8

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 18-May-2000 10:48 eppinged Quar Quant Type: ESTD Cal Date : 18-MAY-2000 00:40 Cal File: c-a2297.d

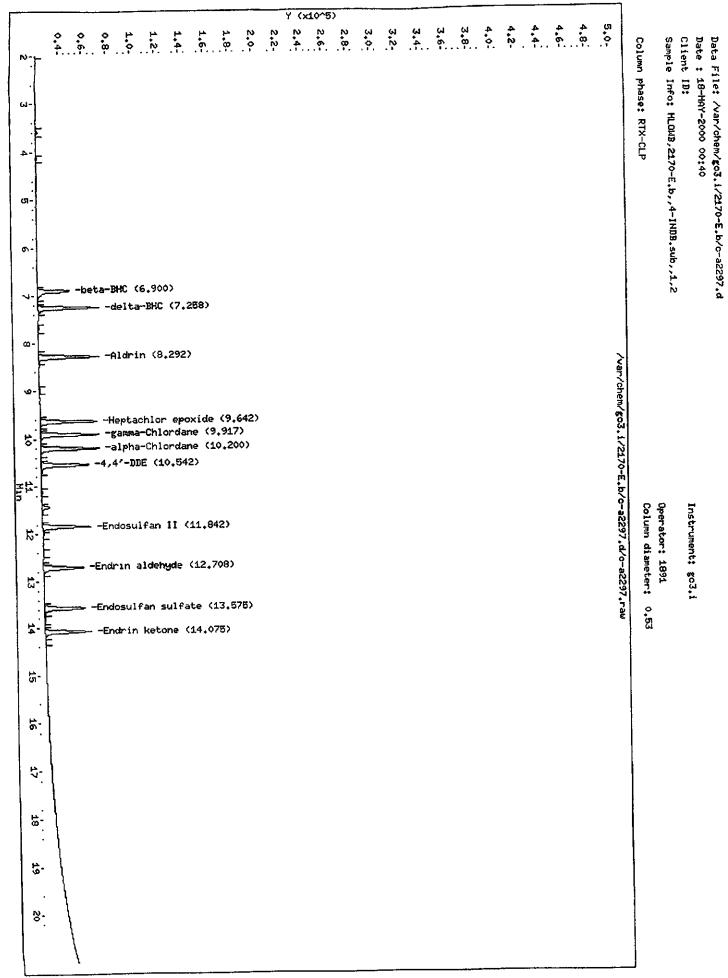
Calibration Sample, Level: 2 Als bottle: 1

Dil Factor: 1.00000 Integrator: Falcon

Compound Sublist: 4-INDB.sub

Target Version: 3.40

						AMOU	NTS
						CAL-AMT	ON-COL
Compo	unds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
		**		*****			***
11	Aldrin	8.292	8.292	0.000	49375	0.01000	0.0100686
7	beta-BHC	6.900	6.900	0.000	26701	0.01000	0.00978256
8	delta-BHC	7.258	7.258	0 000	50477	0.01000	0 0102615
12	Heptachlor epoxide	9 642	9.642	0.000	47015-	0.01000	0.00992747
13	gamma-Chlordane	9 917	9 917	0.000	48385	0.01000	0.00996160
14	alpha-Chlordane	10.200	10.200	0.000	48410	0 01000	0 00995906
16	4,4'-DDE	10 542	10.542	0.000	39365	0.01000	0.0101113
22	Endosulfan II	11.842	11.842	0 000	39684	0.01000	0 0100280
24	Endrin aldehyde	12.708	12.708	0.000	32951	0.01000	0 00991350
26	Endosulfan sulfate	13.575	13.575	0.000	33868	0.01000	0.00996030
27	Endrin ketone	14.075	14 075	0.000	38133	0 01000	0.0100386



Data File: /var/chem/gc3.i/2170-E.b/c-a2298.d

Report Date: 18-May-2000 10:49

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-a2298.d

Lab Smp Id: MEDB

Inj Date : 18-MAY-2000 01:06

Operator: 1891 Ir Smp Info: MEDB,2170-E.b,,4-INDB.sub,,1,3 Inst ID: gc3.i

Misc Info : 190-84-9

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m

Meth Date: 18-May-2000 10:48 eppinged Cal Date: 18-MAY-2000 01:06 Quant Type: ESTD Cal File: c-a2298.d

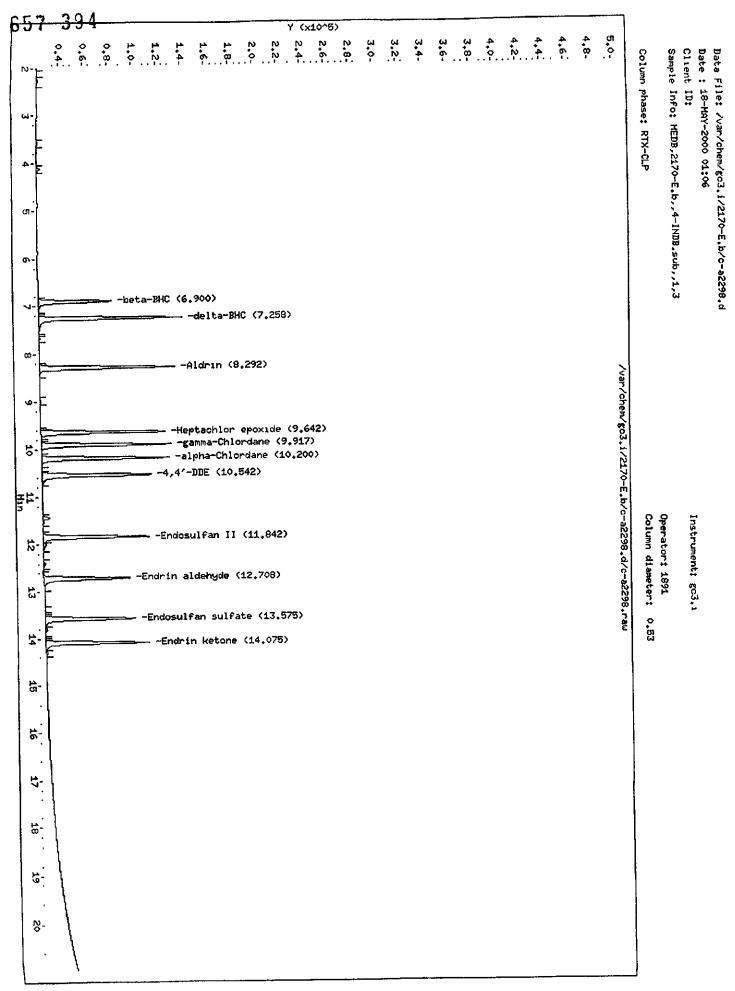
Als bottle: 1 Calibration Sample, Level: 3

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 4-INDB.sub

Target Version: 3.40

						AMOU	DUNTS	
						CAL-AMT	ON-COL	
Compounds		RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)	
****		<b>20</b>		****			*****	
11	Aldrin	8.292	8.292	0 000	111349	0.02500	0.0234227	
7	beta-BHC	6 900	6.900	0.000	60349	0.02500	0.0229963	
8	delta-BHC	7.258	7.258	0.000	118293	0.02500	0.0243571	
12	Heptachlor epoxide	9.642	9.642	0.000	102915	0.02500	0.0227214	
13	gamma-Chlordane	9.917	9.917	0.000	107199	0.02500	0.0229675	
14	alpha-Chlordane	10.200	10.200	0.000	105442	0.02500	0+0226928	
16	4,4'-DDE	10 542	10 542	0.000	90274	0.02500	0.0237620	
22	Endosulfan II	11.842	11.842	0.000	87190	0.02500	0.0229403	
24	Endrin aldehyde	12.708	12 708	0.000	71927	0.02500	0.0226547	
26	Endosulfan sulfate	13.575	13.575	0.000	74851	0.02500	0.0229261	
27	Endrin ketone	14.075	14.075	0.000	85985	0.02500	0.0233725	



Data File: /var/chem/gc3.i/2170-E.b/c-a2299.d

Report Date: 18-May-2000 10:49

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2299.d

Lab Smp Id: MHIGHB
Inj Date : 18-MAY-2000 01:32

Operator : 1891 Inst ID: gc3.i

Smp Info : MHIGHB, 2170-E.b, ,4-INDB.sub, ,1,4

Misc Info: 190-84-10

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 18-May-2000 10:48 eppinged Quar Cal Date : 18-MAY-2000 01:32 Cal Quant Type: ESTD Cal File: c-a2299.d

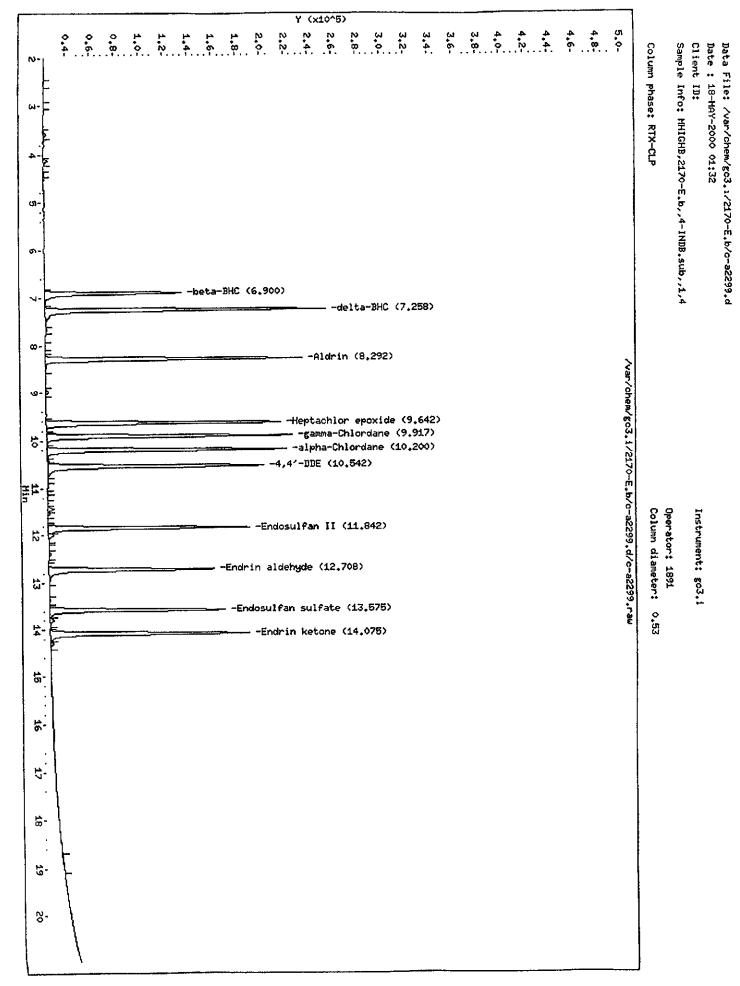
Als bottle: 1 Calibration Sample, Level: 4

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 4-INDB.sub

Target Version: 3.40

						AMOU	NTS
						CAL-AMT	ON-COL
Compo	un <b>ds</b>	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
=====	, # = = = = <del> </del>		**====		20078888		******
11	Aldrin	8.292	8.292	0.000	211501	0.05000	0.0457505
7	beta-BHC	6.900	6 900	0.000	112461	0.05000	0 0444418
8	delta-BHC	7.258	7 258	0.000	231896	0.05000	0.0482922
12	Heptachlor epoxide	9.642	9.642	0.000	193263	0.05000	0.0442919
13	gamma-Chlordane	9.917	9.917	0.000	202560	0.05000	0.0448800
14	alpha-Chlordane	10.200	10.200	0.000	197445	0.05000	0 0441504
16	4,4'-DDE	10.542	10.542	0.000	177218	0.05000	0.0474428
22	Endosulfan II	11.842	11.842	0.000	165821	0.05000	0.0450643
24	Endrin aldehyde	12.708	12.708	0.000	136340	0.05000	0.0445134
26	Endosulfan sulfate	13.575	13.575	0.000	144126	0 05000	0.0454758
27	Endrin ketone	14.075	14.075	0.000	163917	0.05000	0.0458027



Data File: /var/chem/gc3.i/2170-E.b/c-a2300.d

Report Date: 18-May-2000 10:49

## STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-a2300.d Lab Smp Id: HIGHB

Inj Date : 18-MAY-2000 01:57

Inst ID: gc3.i Operator : 1891

Smp Info : HIGHB, 2170-E.b,, 4-INDB.sub,, 1,5 Misc Info : 190-84-11

Comment :

: /var/chem/gc3.i/2170-E.b/PESTA.m Method

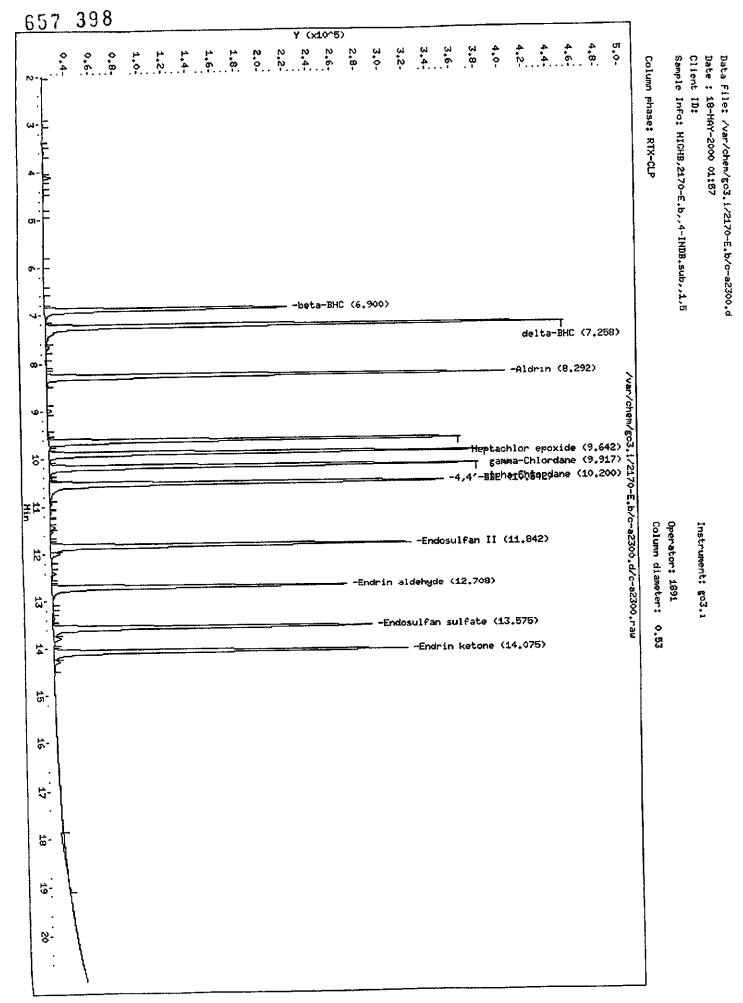
Meth Date: 18-May-2000 10:48 eppinged Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-a2300.d

Calibration Sample, Level: 5 Als bottle: 1

Dil Factor: 1.00000 Integrator: Falcon Compound Sublist: 4-INDB.sub

Target Version: 3.40

					AMOU	nts
					CAL-AMT	ON-COL
Compounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
<b>建国祖国民主众教育党与自由进行公司来等国建设公司等</b>		*****		********	医医医氏氏结节	
11 Aldrin	8.292	8.292	0 000	379733	0 10000	0 0851839
7 beta-BHC	6.900	6.900	0.000	199137	0.10000	0.0821966
8 delta-BHC	7.258	7.258	0.000	429107	0.10000	0 0913040
12 Heptachlor epoxide	9.642	9.642	0 000	341408	0 10000	0 0818032
13 gamma-Chlordane	9.917	9.917	0 000	363079	0 10000	0 0837195
14 alpha-Chlordane	10.200	10.200	0.000	355828	0.10000	0.0829565
16 4,4'-DDE	10.542	10.542	0.000	326794	0.10000	0 0897314
22 Endosulfan II	11.842	11 842	0.000	299189	0.10000	0.0844665
24 Endrin aldehyde	12 708	12.708	0 000	243846	0.10000	0.0829970
26 Endosulfan sulfate	13.575	13.575	0.000	264102	0.10000	0.0862054
27 Endrin ketone	14.075	14 075	0.000	293813	0.10000	0.0851475



Data File: /var/chem/gc3.i/2170-E.b/c-a2301.d

Report Date: 18-May-2000 10:59

### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-a2301.d Lab Smp Id: 2ND A Inj Date : 18-MAY-2000 02:23

Inst ID: gc3.i Operator : 1891

Smp Info : 2ND A,2170-E.b,,INDA.sub,,2,3

Misc Info : 190-82-2

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m

Quant Type: ESTD Meth Date: 18-May-2000 10:58 g Cal Date : 18-MAY-2000 01:57 Cal File: c-a2300.d

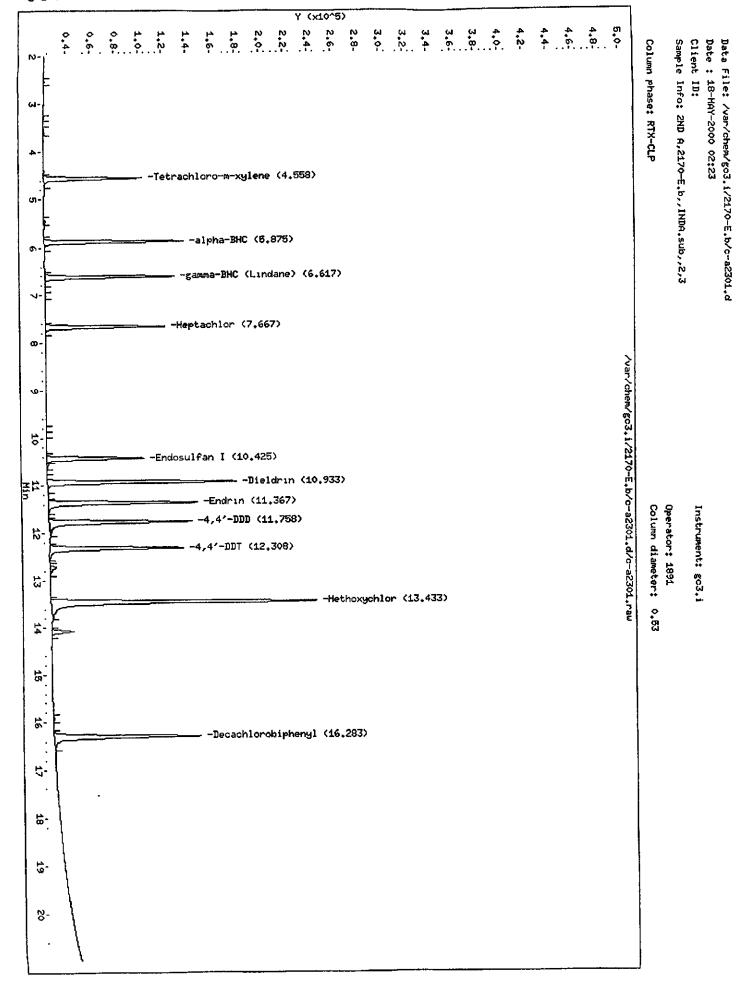
Continuing Calibration Sample Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: INDA.sub Integrator: Falcon

Target Version: 3.40

					AMOU	NTS
					CAL-AMT	ON-COL
Compounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
		*****	******	******	*********	****
s 1 Tetrachloro-m-xylene	4.558	4.558	0.000	81097	0.02500	0 0200288
5 alpha-BHC	5.875	5.883	-0.008	114953	0.02500	0 0206635
6 gamma-BHC (Lindane)	6.617	6 617	0.000	107249	0.02500	0.0210179
10 Heptachlor	7.667	7 675	-0.008	98867	0.02500	0.0206271
15 Endosulfan I	10.425	10 425	0.000	80119	0 02500	0.0207649
17 Dieldrin	10.933	10.933	0 000	155908	0.02500	0.0384554
20 Endrin	11.367	11 367	0.000	123356	0 02500	0.0349515
21 4,4'-DDD	11.758	11.758	0.000	117992	0.02500	0.0404374
23 4,4'-DDT	12.308	12.308	0.000	110738	0.02500	0.0399539
25 Methoxychlor	13.433	13.442	-0.009	219957	0 05000	0 157927
\$ 30 Decachlorobiphenyl	16.283	16.283	0.000	120967	0 02500	0.0367198



Data File: /var/chem/gc3.i/2170-E.b/c-a2302.d

Report Date: 18-May-2000 10:59

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2302.d

Lab Smp Id: 2ND B

Inj Date : 18-MAY-2000 02:49

Operator: 1891 Smp Info: 2ND B,2170-E.b,,INDB.sub,,2,3 Inst ID: gc3.i

Misc Info : 190-82-5

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 18-May-2000 10:58 g Quar Cal Date : 18-MAY-2000 01:57 Cal Quant Type: ESTD Cal File: c-a2300.d

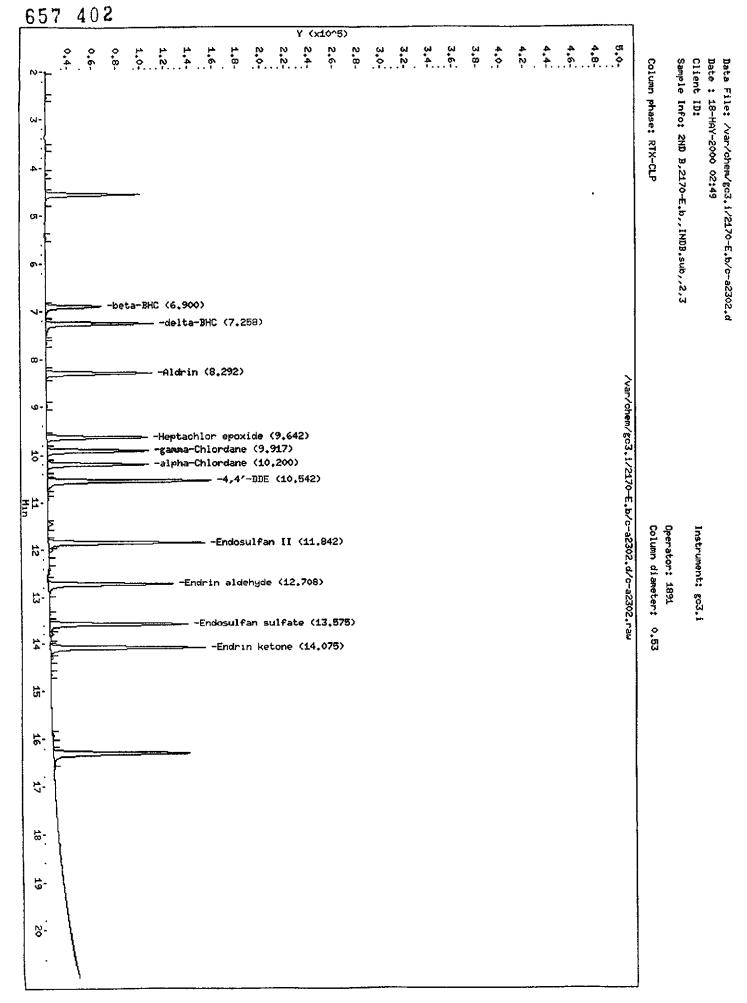
Continuing Calibration Sample Als bottle: 1

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: INDB.sub

Target Version: 3.40

					AMOU	NTS
					CAL-AMT	ON-COL
Compounds	RT	EXP RT	DLT RT	response	( ng)	( ng)
医全球性蛋白医胃抗性血管医疗性皮肤性结合性	<b>4 17</b>		590000	*****	222844	
11 Aldrin	8.292	8.292	0.000	87079	0.02500	0.0195341
7 beta-BHC	6.900	6.900	0.000	45737	0 02500	0.0188786
8 delta-BHC	7.258	7 258	0.000	88859	0 02500	0 0189071
12 Heptachlor epoxide	9.642	9.642	0.000	83488	0.02500	0.0200042
13 gamma-Chlordane	9 917	9.917	0.000	83894	0.02500	0.0193444
14 alpha-Chlordane	10 200	10 200	0.000	84129	0 02500	0.0196135
16 4,4'-DDE	10.542	10.542	0.000	135583	0.02500	0 0372285
22 Endosulfan II	11 842	11 842	0.000	129381	0.02500	0.0365266
24 Endrin aldehyde	12,708	12.708	0 000	104257	0 02500	0 0354856
26 Endosulfan sulfate	13.575	13.575	0.000	114472	0 02500	0.0373648
27 Endrin ketone	14.075	14.075	0.000	128548	0.02500	0 0372534



Data File: /var/chem/gc3.i/2170-E.b/c-a2303.d

Report Date: 18-May-2000 10:59

### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-a2303.d

Lab Smp Id: EVALB

Inj Date : 18-MAY-2000 03:15

Operator: 1891 Inst ID: gc3.i

Smp Info : EVALB, 2170-E.b, , EVALBR. sub, , 3, 1

Misc Info : 190-88-8

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTA.m

Meth Date: 18-May-2000 10:58 g Quant Type: ESTD Cal Date: 18-MAY-2000 01:57 Cal File: c-a2300.d

Als bottle: 1 QC Sample: PEM

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: EVALBR.sub

Target Version: 3.40

						CONCENT	RATIONS
						ON-COLUMN	FINAL
Co	mpounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	(ng)
30	医咽性病毒性分类 医肾炎 计过程 医自己性 法国 医 电电	21 di		*****	========		6345222
\$	1 Tetrachloro-m-xylene	4.558	4.558	0.000	82955	0.02049	0.0204877(R)
	16 4,4'-DDE	10.542	10 542	0.000	729	0.00020	0.000200170
	20 Endrin	11.367	11.367	0.000	85719	0.02429	0.0242875
	21 4,4'-DDD	11.758	11.758	0.000	7838	0.00269	0.00268618
	23 4,4'-DDT	12.308	12.308	0.000	67934	0.02451	0.0245103
	24 Endrin aldehyde	12.717	12.708	0.009	2199	0.000748	0.000748466
	27 Endrin ketone	14.075	14.075	0.000	5329	0.00154	0 00154435
\$	30 Decachlorobiphenyl	16.283	16.283	0 000	67410	0.02046	0.0204624(R)

QC Flag Legend

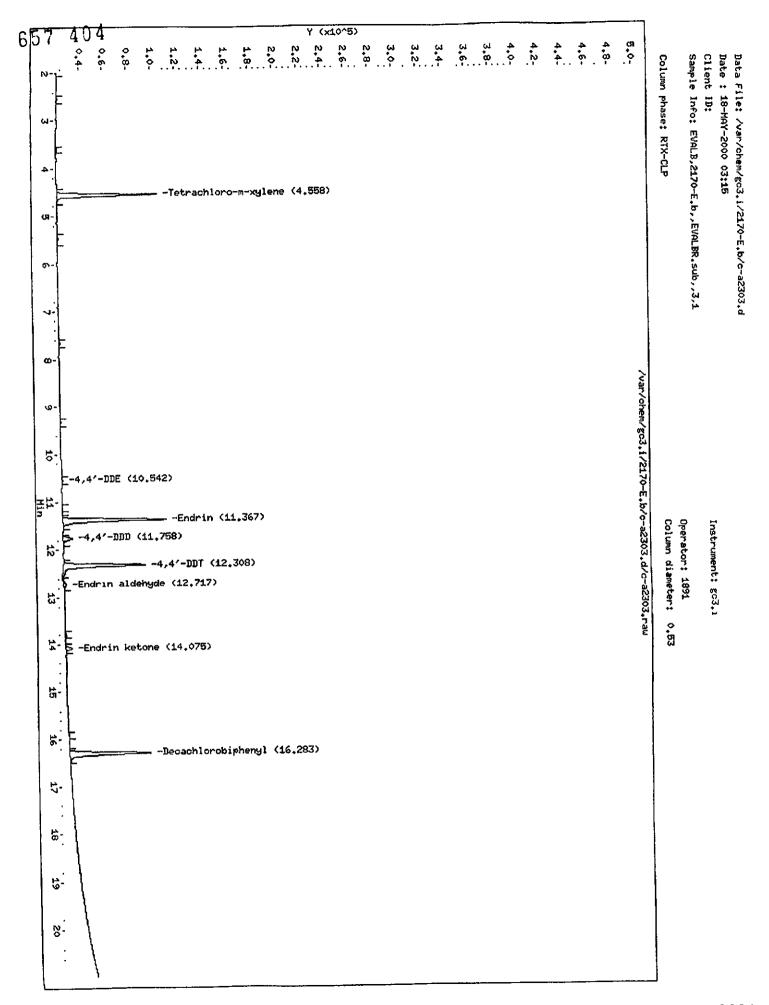
R - Spike/Surrogate failed recovery limits.

97021 A.170

Endrin Beentedon: (2199+5339)x10ch 9.106 (2199+5339+053-19)

DDT Brankda - (738 + 7435) (105-11.81)

3083



Data File: /var/chem/gc3.i/2170-E.b/c-a2361.d

Report Date: 19-May-2000 14:21

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2361.d

Lab Smp Id: EVALB

Inj Date : 19-MAY-2000 08:32

Operator : 1891 Inst ID: qc3.i

Smp Info : EVALB, 2170-E.b, , EVALBR. sub, , 3, 1 Misc Info : 190-88-8

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m

Meth Date: 19-May-2000 11:01 eppinged Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-a2300.d

Als bottle: 1 QC Sample: PEM

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: EVALBR.sub

Target Version: 3.40

							CONCENT	RATIONS
							ON-COLUMN	FINAL
Co	οqπ	unds	RT	EXP RT	DLT RT	RESPONSE	( ng)	(ng)
	<b>22</b>				F20960	========		
\$	1	Tetrachloro-m-xylene	4.558	4.558	0.000	80031	0.01977	0.0197655(R)
	16	4,4'-DDE	10.550	10.550	0.000	653	0.000179	0.000179301(M)
	20	Endrin	11.375	11 375	0.000	84290	0 02388	0.0238826
	21	4,4'-DDD	11.767	11.767	0 000	6776	0 00232	0.00232222
	23	4,4'-DDT	12.317	12.317	0.000	67440	0.02433	0.0243321
	24	Endrın aldehyde	12 717	12.717	0 000	1257	0 000428	0.000427841
	27	Endrin ketone	14.083	14 083	0.000	3829	0 00111	0.00110965
\$	30	Decachlorobiphenyl	16.292	16.300	-0.008	64670	0.01963	0.0196307(R)

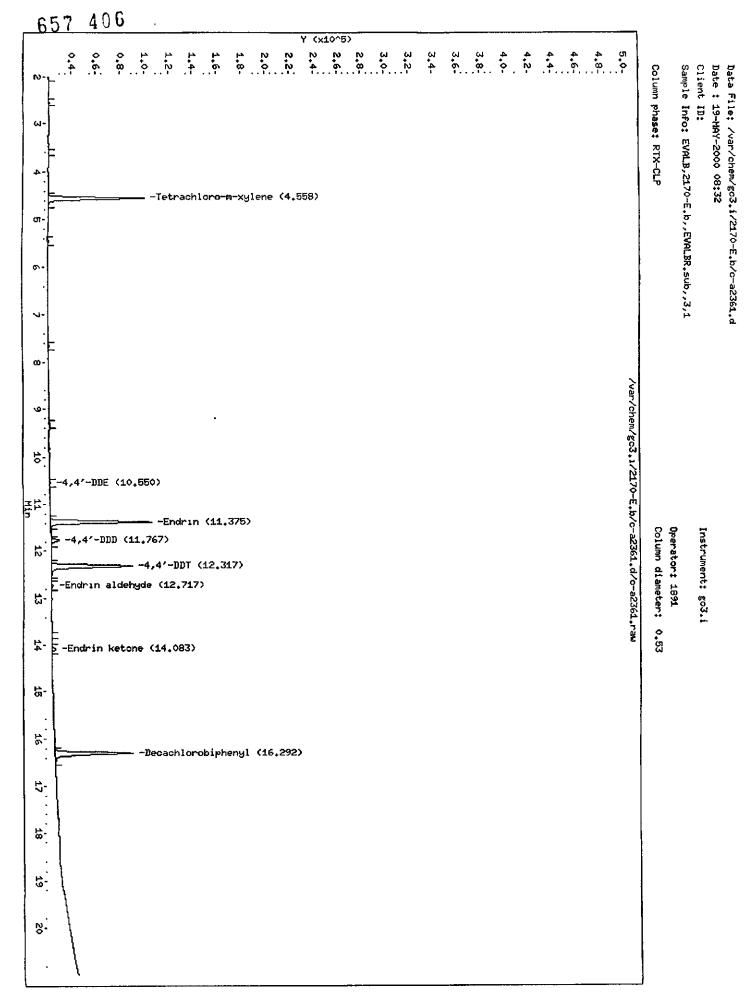
## QC Flag Legend

R - Spike/Surrogate failed recovery limits.

M - Compound response manually integrated.

DOT 9.907. EXDRIN 5.77.

Endrum Biceatabr\_ (1267+309) Xcoc 5,7%



Data File: /var/chem/gc3.i/2170-E.b/c-a2362.d

Report Date: 19-May-2000 10:12

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-a2362.d Lab Smp Id: MEDTOX

Inj Date : 19-MAY-2000 08:58

Operator: 1891 Inst ID: gc3.i

Smp Info : MEDTOX,2170-E.b,,1-TOX.sub,,2,3

Misc Info: 190-84-13

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 19-May-2000 10:12 g Quar Cal Date : 18-MAY-2000 01:57 Cal Quant Type: ESTD Cal File: c-a2300.d

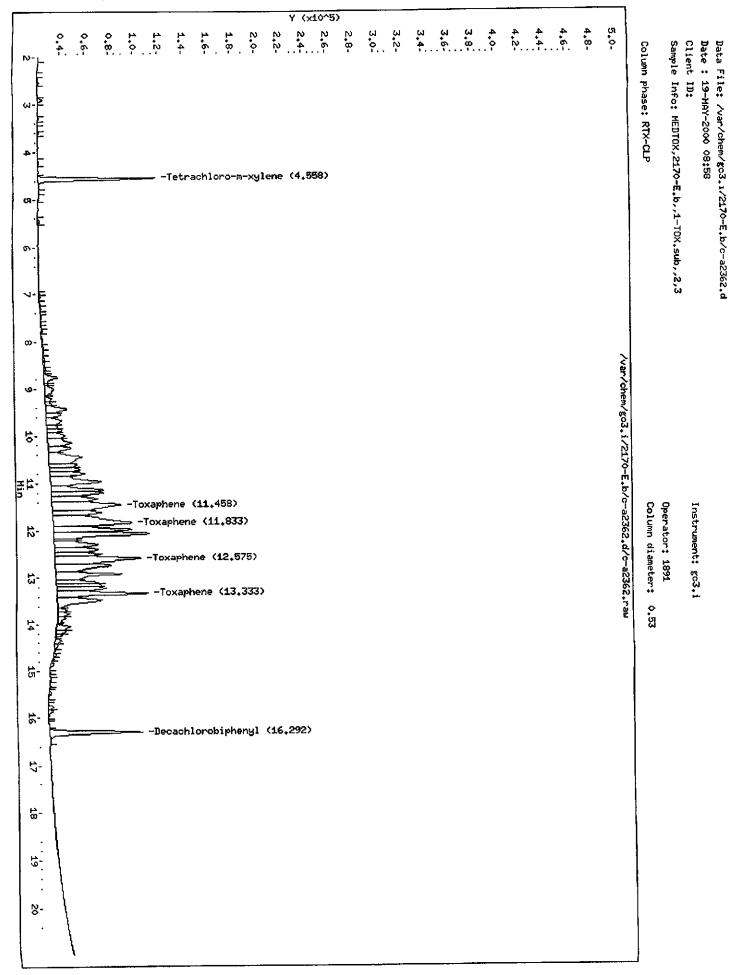
Als bottle: 1 Continuing Calibration Sample

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 1-TOX.sub

Target Version: 3.40

					AMOU	nts
					CAL-AMT	ON-COL
Compounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
	===		E G O O O O		*****	******
18 Toxaphene	11.458	11.458	0.000	57622	1.00000	1.02439
<pre>\$ 1 Tetrachloro-m-xylene</pre>	4.558	4.558	0.000	96741	0.02500	0.0238924
\$ 30 Decachlorobiphenyl	16.292	16.300	-0.008	77437	0.02500	0.0235061



Data File: /var/chem/gc3.i/2170-E.b/c-a2363.d

Report Date: 19-May-2000 10:13

### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2363.d

Lab Smp Id: MEDCHLOR

Inj Date : 19-MAY-2000 09:24

Operator : 1891 Inst ID: gc3.i

Smp Info : MEDCHLOR, 2170-E.b,, 2-CHLO.sub,, 2,3 Misc Info : 190-85-10

Comment

: /var/chem/gc3.i/2170-E.b/PESTA.m Method

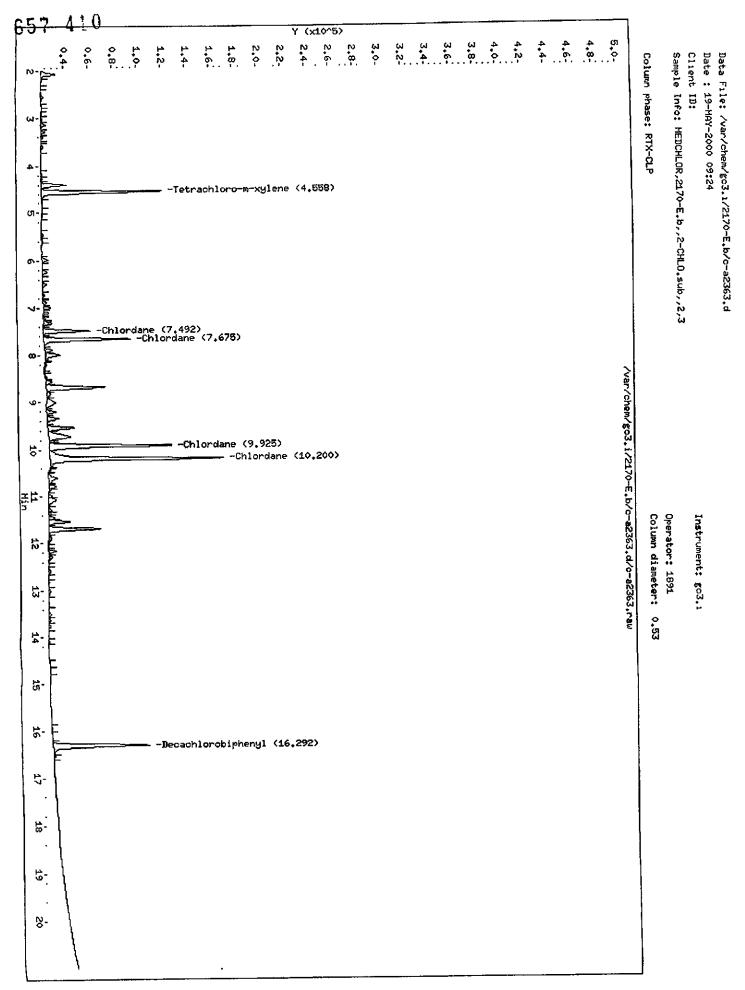
Meth Date: 19-May-2000 10:12 g Quant Type: ESTD Cal Date :  $18-MA\bar{Y}-2000 01:57$ Cal File: c-a2300.d

Als bottle: 1 Continuing Calibration Sample

Dil Factor: 1.00000 Integrator: Falcon Compound Sublist: 2-CHLO.sub

Target Version: 3.40

						MOUA	nts
						CAL-AMT	ON-COL
Co	mpounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
==	*	# =		a	****		*****
	9 Chlordane	7.492	7.492	0 000	39014	0.25000	0.256448
\$	1 Tetrachloro-m-xylene	4.558	4.558	0.000	98915	0 02500	0.0244294
\$	30 Decachlorobiphenyl	16.292	16.300	-0.008	80232	0.02500	0.0243546



Data File: /var/chem/gc3.i/2170-E.b/c-a2364.d

Report Date: 19-May-2000 10:13

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2364.d

Lab Smp Id: MEDA

Inj Date : 19-MAY-2000 09:50

Operator : 1891 Inst ID: qc3.i

Smp Info : MEDA,2170-E.b,,INDA.sub,,2,3

Misc Info: 190-84-3

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 19-May-2000 10:12 g Quar Cal Date : 18-MAY-2000 01:57 Cal Quant Type: ESTD Cal File: c-a2300.d

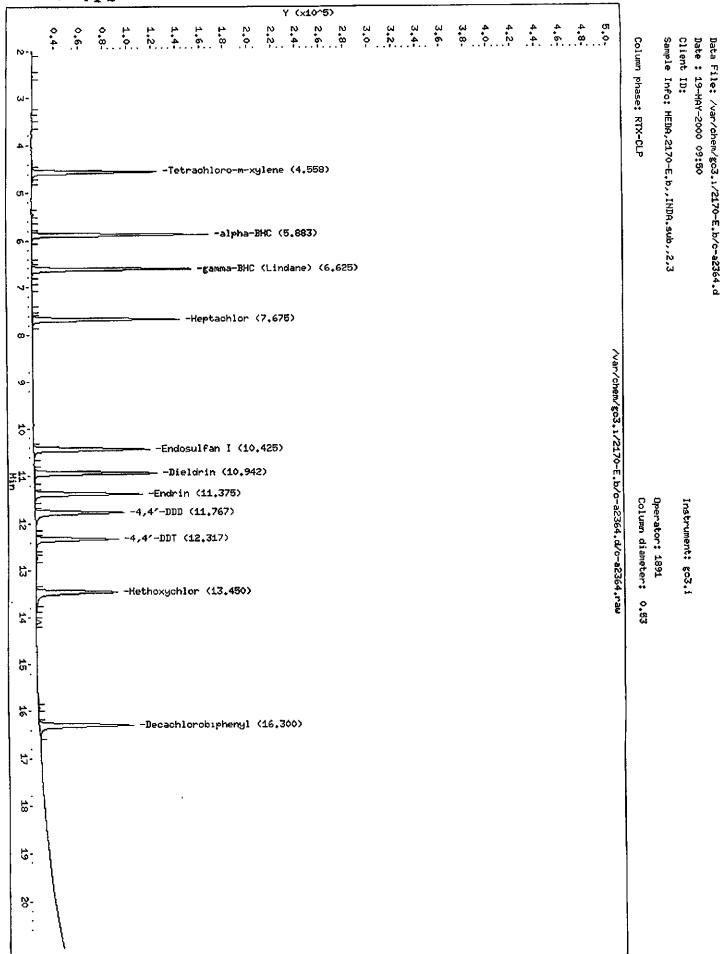
Als bottle: 1 Continuing Calibration Sample

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: INDA.sub

Target Version: 3.40

						AMOU	NTS
						CAL-AMT	ON-COL
Com	pounds	RT	EXP RT	DLT RT	response	( ng)	( ng)
		90	***	======	******		
\$	1 Tetrachloro-m-xylene	4.558	4 558	0 000	103245	0.02500	0.0254988
	5 alpha-BHC	5 883	5.883	0.000	145571	0.02500	0.0261672
	6 gamma-BHC (Lindane)	6.625	6.625	0 000	131234	0.02500	0.0257184
	10 Heptachlor	7.675	7.675	0.000	121243	0.02500	0.0252955
	15 Endosulfan I	10.425	10.425	0.000	96330	0 02500	0.0249664
	17 Dieldrin	10.942	10.942	0.000	101139	0 02500	0.0249464
	20 Endrin	11.375	11.375	0.000	89364	0 02500	0.0253203
	21 4,4'-DDD	11.767	11.767	0.000	73656	0.02500	0.0252428
	23 4,4'-DDT	12.317	12.317	0.000	68697	0 02500	0.0247856
	25 Methoxychlor	13.450	13.450	0 000	67068	0.05000	0.0481542
\$	30 Decachlorobiphenyl	16.300	16.300	0.000	78142	0.02500	0.0237201



Data File: /var/chem/gc3.i/2170-E.b/c-a2365.d

Report Date: 19-May-2000 11:00

### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2365.d

Lab Smp Id: MEDB

Inj Date : 19-MAY-2000 10:16

Operator: 1891 Inst ID: gc3.i

Smp Info : MEDB,2170-E.b,,INDB.sub,,2,3
Misc Info : 190-84-9

Comment

: /var/chem/gc3.i/2170-E.b/PESTA.m Method

Meth Date: 19-May-2000 11:00 g Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-a2300.d.

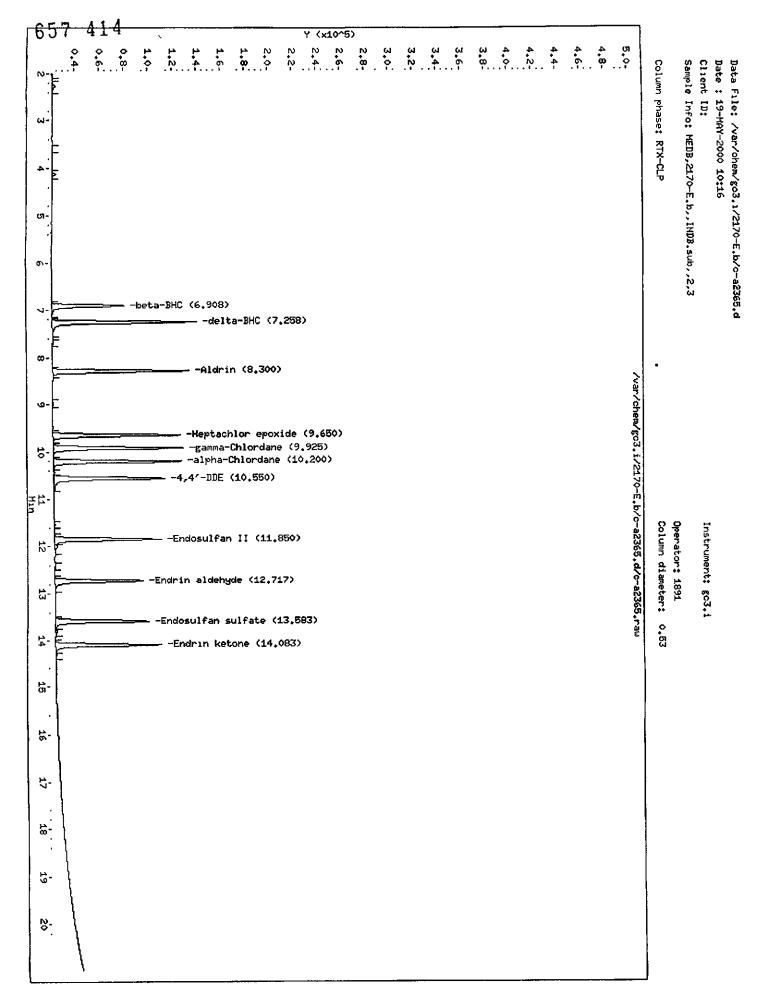
Als bottle: 1 Continuing Calibration Sample

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: INDB.sub

Target Version: 3.40

						AMOU.	nts
						CAL-AMT	ON-COL
Compour	nds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
				======	*******		
11 /	Aldrin	8.300	8.300	0 000	112275	0.02500	0.0251862
7 1	beta-BHC	6.908	6.908	0.000	59342	0.02500	0 0244943
8 (	delta-BHC	7.258	7.258	0.000	118180	0.02500	0.0251460
12	Heptachlor epoxide	9.650	9.650	0.000	104693	0.02500	0.0250850
13 9	gamma-Chlordane	9.925	9.925	0.000	106659	0.02500	0.0245936
14	alpha-Chlordane	10.200	10.200	0 000	104845	0.02500	0.0244432
16	4,4'-DDE	10.550	10.550	0.000	91188	0.02500	0.0250385
22	Endosulfan II	11.850	11.850	0.000	88332	0.02500	0.0249377
24	Endrin aldehyde	12.717	12 717	0.000	72782	0.02500	0.0247726
26	Endosulfan sulfate	13.583	13 583	0.000	76798	0.02500	0.0250676
27	Endrin ketone	14.083	14 083	0.000	87171	0.02500	0.0252623



Data File: /var/chem/gc3.i/2170-E.b/c-a2374.d

Report Date: 19-May-2000 14:32

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2374.d Lab Smp Id: MEDA

Inj Date : 19-MAY-2000 14:10

Operator : 1891 Inst ID: gc3.i

Smp Info : MEDA, 2170-E.b, , INDA. sub, , 2, 3

Misc Info: 190-84-3

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 19-May-2000 14:32 g Quar Cal Date : 18-MAY-2000 01:57 Cal Quant Type: ESTD Cal File: c-a2300.d

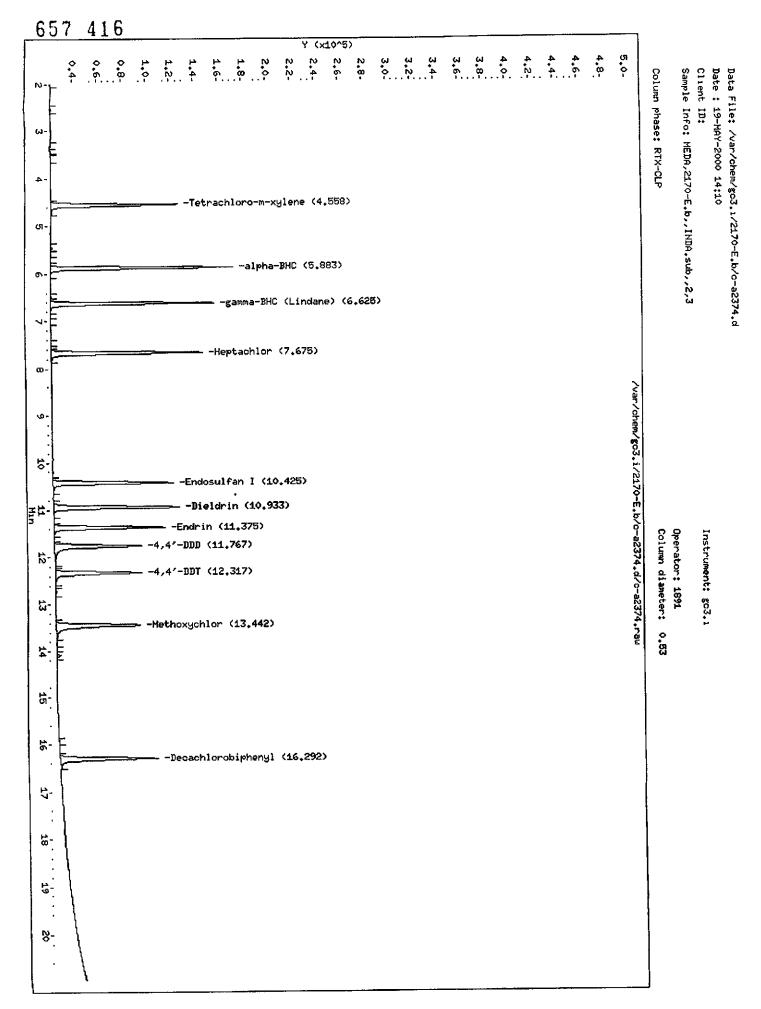
Als bottle: 1 Continuing Calibration Sample

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: INDA.sub

Target Version: 3.40

							AMO	UNTS
							CAL-AMT	ON-COL
Co	ogmo	unds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
		医克里氏征 医甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基甲基	==	222277		2927000		*****
\$	1	Tetrachloro-m-xylene	4.558	4.558	0 000	104421	0.02500	0.0257892
	5	alpha-BHC	5.883	5.883	0.000	149994	0.02500	0.0269623
	6	gamma-BHC (Lindane)	6.625	6 625	0.000	134401	0.02500	0.0263390
	10	Heptachlor	7.675	7.675	0.000	124632	0.02500	0.0260026
	15	Endosulfan I	10 425	10.425	0 000	98693	0.02500	0.0255788
	17	Dieldrin	10.933	10.942	-0.009	103369	0.02500	0.0254964
	20	Endrin	11.375	11.375	0.000	91479	0 02500	0.0259195
	21	4,4'-DDD	11.767	11 767	0.000	72022	0.02500	0.0246828
	23	4,4'-DDT	12.317	12 317	0.000	71695	0.02500	0 0258673
	25	Methoxychlor	13.442	13 450	-0 008	70047	0.05000	0 0502931
\$	30	Decachlorobiphenyl	16.292	16.300	-0.008	80592	0.02500	0.0244638



Data File: /var/chem/gc3.i/2170-E.b/c-a2375.d

Report Date: 19-May-2000 15:20

### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2375.d

Lab Smp Id: MEDB

Inj Date : 19-MAY-2000 14:36 Operator : 1891 Smp Info : MEDB,2170-E.b,,INDB.sub,,2,3 Inst ID: gc3.i

Misc Info: 190-84-9

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 19-May-2000 15:20 g Quar Cal Date : 18-MAY-2000 01:57 Cal Quant Type: ESTD Cal File: c-a2300.d

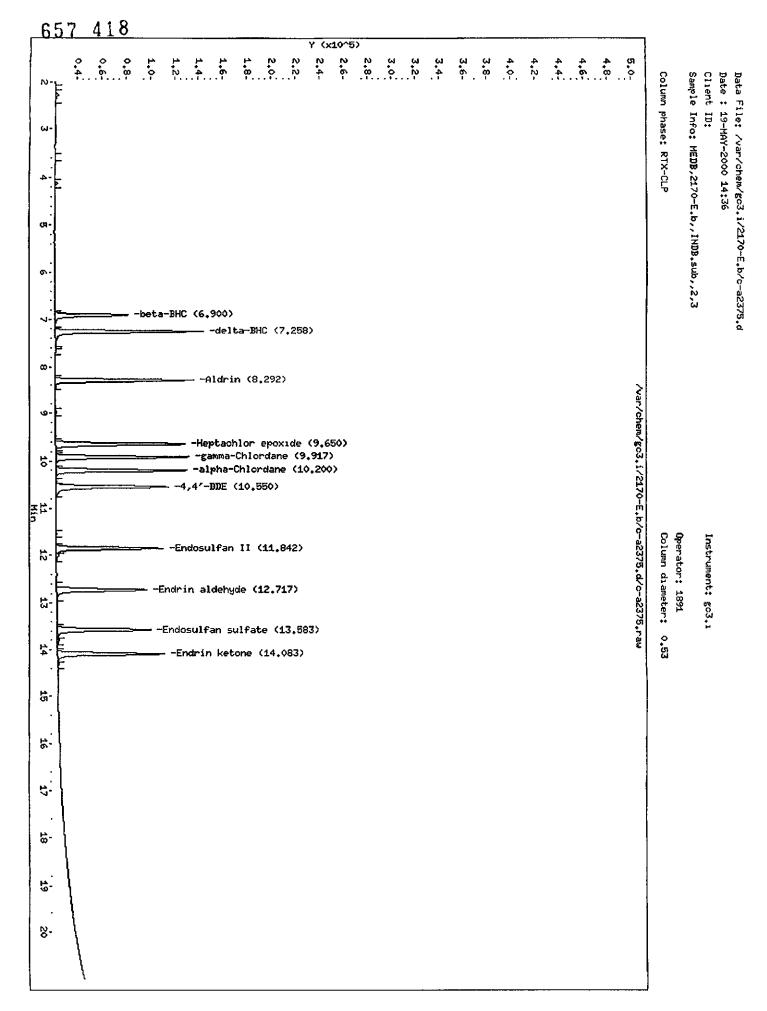
Als bottle: 1 Continuing Calibration Sample

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: INDB.sub

Target Version: 3.40

						AMOU	NTS
						CAL-AMT	ON-COL
Compo	unds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
		<b>10</b> 10			*******	*****	0=4=558
11	Aldrin	8.292	8.300	-0.008	113488	0.02500	0.0254583
7	beta-BHC	6.900	6.908	-0 008	59935	0.02500	0.0247390
8	delta-BHC	7 258	7.258	0.000	121351	0.02500	0 0258207
12	Heptachlor epoxide	9 650	9.650	0.000	106476	0.02500	0.0255122
13	gamma-Chlordane	9.917	9.925	-0.008	109001	0.02500	0.0251337
14	alpha-Chlordane	10.200	10.200	0 000	107874	0.02500	0.0251494
16	4,4'-DDE	10 550	10.550	0.000	91812	0.02500	0 0252098
22	Endosulfan II	11 842	11.850	-0.008	88448	0.02500	0 0249705
24	Endrin aldehyde	12.717	12.717	0.000	74493	0.02500	0.0253549
26	Endosulfan sulfate	13.583	13 583	0 000	77367	0.02500	0.0252533
27	Endrin ketone	14.083	14.083	0 000	88309	0.02500	0.0255921



Report Date : 19-May-2000 11:24

e7-4-92

## STL-PITTSBURGH

## COMPOUND LISTING

Method file : /var/chem/gc3.i/2170-E.b/PESTB.m Quant Method : ESTD Tar Last Update : 19-May-2000 11:22 Num Target Version : 3.40 Number of Cpnds: 30

Data Type : GC MULTI COMP

Global Integrator : Falcon

Chromat Events	Values
Initial:Start Threshold	40.000000
Initial:End Threshold	20.00000
Initial:Area Threshold	1000.000000
Initial:P-P Resolution	1.000000
Initial:Bunch Factor	1.00000
Initial:Negative Peaks	ON
Initial:Tension	0.00000

Compound	RT	RT Window	RF
\$ 1 Tetrachloro-m-xylene 2 Diallate A 3 Diallate B 4 HEXACHLOROBENZENE 5 alpha-BHC 6 gamma-BHC (Lindane) 7 beta-BHC 8 delta-BHC 9 Chlordane	9.475	7.175-7.275 7.425-7.525 6.600-6.700 7.583-7.683 8.458-8.558 8.750-8.850 9.417-9.517	4.110e+06 1.093e+05 2.882e+04 6.370e+06 5.386e+06 2.737e+06 5.546e+06 1.572e+05 2.993e+05 4.311e+05
10 Heptachlor 11 Aldrin 12 Heptachlor epoxide 13 gamma-Chlordane 14 alpha-Chlordane 15 Endosulfan I 16 4,4'-DDE 17 Dieldrin 18 Toxaphene	12.208 9.475 10.167 11.483 11.892 12.208 12.283 12.692 12.867 13.892 14.117 14.667	12.158-12.258 9.425-9.525 10.117-10.217 11.433-11.533 11.842-11.942 12.158-12.258 12.233-12.333 12.642-12.742 12.817-12.917 13.842-13.942 14.067-14.167 14.617-14.717 15.617-15.717	3.540e+05 5.290e+06 5.126e+06 4.383e+06 4.594e+06 4.478e+06 4.037e+06 3.926e+06 4.152e+06 5.292e+04 8.324e+04 5.234e+04

657 420

Report Date : 19-May-2000 11:24

STL-PITTSBURGH

COMPOUND LISTING

Method file : /var/chem/gc3.i/2170-E.b/PESTB.m

	Compound	RT	RT Window	RF
	19 Isodrin 20 Endrin 21 4,4'-DDD 22 Endosulfan II 23 4,4'-DDT 24 Endrin aldehyde 25 Methoxychlor 26 Endosulfan sulfate 27 Endrin ketone 28 Chlorobenzilate	13.475 13.900 13.950 14.517 14.667 16.017 15.225 16.267	10.967-11.067 13.425-13.525 13.850-13.950 13.900-14.000 14.467-14.567 14.617-14.717 15.967-16.067 15.175-15.275 16.217-16.317 13.608-13.708	3.540e+06 3.012e+06 3.414e+06 2.620e+06 2.781e+06 1.250e+06 2.813e+06 3.087e+06
	29 Kepone 30 Decachlorobiphenyl	16.258	16.208-16.308 18.792-18.892	5.494e+03
<u> </u>				

Report Date : 18-May-2000 11:41

HP58903F

## STL-PITTSBURGH

# INITIAL CALIBRATION DATA

Start Cal Date : 17-MAY-2000 17:15 End Cal Date : 18-MAY-2000 01:57

Quant Method : ESTD : Disabled Origin Target Version : 3.40 Integrator : Falcon

: /var/chem/gc3.i/2170-E.b/PESTB.m Method file

: 18-May-2000 11:40 eppinged Cal Date

: Average Curve Type

Calibration File Names:

Level 1: /var/chem/gc3.i/2170-E.b/c-b2296.d Level 2: /var/chem/gc3.i/2170-E.b/c-b2297.d Level 3: /var/chem/gc3.i/2170-E.b/c-b2298.d Level 4: /var/chem/gc3.i/2170-E.b/c-b2299.d Level 5: /var/chem/gc3.i/2170-E.b/c-b2300.d

	0.00500	0.01000	0.02500	0.05000	0.10000		
Compound	Level 1			•	Level 5	RRF	1 RSD
2 Diallate A	151590   151590	115224	102868	•	87639	109261	23.984
3 Diallate B	1 38910	,	27444	23735	23034	28820	22.466
4 HEXACHLOROBENZENE	1 +++++	++++	+++++	+++++	+++++	+++++	++++
5 alpha-BHC	1 6723400	69788001	6426560	6131440	5591790	6370398	8.458
6 gamma-BHC (Lindane)	5987400		5422400	5115020	4805370	5385878	8.402
7 beta-BHC	3179200		2670720	2510600	2254170	2737018	14.097
8 delta-BHC	5741200		5513440	5488780	5016420	S545788 j	6.387
9 Chlordane(1)	+++++	+++++	157184	+++++	+++++	157184	0.000
(2)	+++++	! +++++	299340	+++++	+++++ 1	299340	0.000
(3)	1 +++++	! +++++	431052	+++++	++++	431052	0.000
(4)		+++++	354008	+++++	+++++	354008	0.000
10 Heptachlor	1 5942200	5881700	5329920	4951000	4346410	5290246	12.618
11 Aldrin	1 5578800	5669700	5115080	4905880	4361710	5126234	10 384
12 Heptachlor epoxide	1 4822000	4791600	4383000	4209560	3708950	4383022	10.481
13 gamma-Chlordane	1 4952400		4555000	4446800	4005850	4593890	8.910
14 alpha-Chlordane	1 4880000	•	4466440	4268740	3865800	4477596	9.763
15 Endosulfan I	1 4496400	•	4025160	3834920	3348880	4036632	11.880
	1 4011800	•	•	3897780	3558410	3926194	6.125
16 4,4'-DDE 17 Dieldrin	4389200	•	4181680	4047900	3597860	4151588	8 740
18 Toxaphene(1)	****	+++++	52924	+++++	+++++	52924	0.000
(2)	. +++++		83240	++++	+++++	83240	0.000
(3)	+++++	+++++	52341	+++++	1 +++++	52341	0.000
(4)	++++	. +++++	52025	+++++	+++++	52025	0.000
19 Isodrin	4341000	3875800	3819980	3402227	3370460	3761893	10.58
20 Endrin	375480	•	3567880	3457840	3104170	3539938	7.98
20 ENGIIN 21 4,4'-DDD	1 310520	•	2981520	2990380	2783470	3011574	5 16

Report Date : 18-May-2000 11:41

657 422

STL-PITTSBURGH

# INITIAL CALIBRATION DATA

Start Cal Date : 17-MAY-2000 17:15 End Cal Date : 18-MAY-2000 01:57 Quant Method : ESTD Origin : Disabled

Target Version : 3.40
Integrator : Falcon

: /var/chem/gc3.i/2170-E.b/PESTB.m Method file

: 18-May-2000 11:40 eppinged Cal Date

Curve Type : Average

	0 00500     Level 1	•		0.05000   Level 4		RRF	* RSD
Compound		•	·-				
22 Endosulfan II	3621800	3674400	3396240			3413566	7.451
23 4,4'-DDT	2584600	2703300	2613640	2696140	2503420	2620220	3.171
24 Endrin aldehyde	3040400	3072200	2709440	2660020	2422400	2780892	9 855
25 Methoxychlor	[ 1340700]	1346800	1220500	1216690	1125375]	1250013	7.495
26 Endosulfan sulfate	2918400	3006300	2766960	2787380	2585310	2812870	5.707
27 Endrin ketone	3226000	3351700	3056760	3015000	2783910	3086674	7.011
28 Chlorobenzilate	1251100	1949760	1471620	2048053	2513620	1846831	26.956
29 Kepone	5650	5288	5726	5277	5530	5494	3.739
B 등 등 후 후 후 등 등 등 후 후 후 후 후 후 후 후 후 후 후			222222222			47,005,001	17.414
1 Tetrachloro-m-xylene	4960800	4686400				4109568	
30 Decachlorobiphenyl	3138600	2973000	2634680	2493380	2285470	2705026	12.880
				ll		···-	¦

657 423

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: SDGA18632

GC Column: RTX-CLP2 ID: 0.53 (mm) Init. Calib. Date(s): 05/17/00 05/18/00

EPA Sample No.(PIBLK): \_\_\_\_\_ Date Analyzed:\_\_\_\_\_

Lab Sample ID (PIBLK): \_\_\_\_\_ Time Analyzed :\_\_\_\_

EPA Sample No. (PEM): Date Analyzed :05/17/00

Lab Sample ID (PEM): EVALB Time Analyzed :1649

PEM COMPOUND	RT	RT WI FROM	NDOW TO	CALC AMOUNT (ng)	NOM AMOUNT (ng)	%D
Endrin 4,4'-DDT	13.47 14.51	1	13.53 14.56	0.023922 0.023365	0.025000 0.025000	-4.3 -6.5

12.2

4,4'-DDT % breakdown (1):

Endrin % breakdown (1):

7\_22\_

Combined % breakdown (1):

15/7

DE 5-19-06

OLMO3.0

PESTICIDE CALIBRATION VERIFICATION SUMMARY 657 424

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT

Case No.:

SAS No.: 40325 SDG No.: SDGA18632

GC Column: RTX-CLP2 ID: 0.53 (mm) Init. Calib. Date(s): 05/17/00 05/18/00

EPA Sample No. (PIBLK):

Date Analyzed :\_\_\_\_\_

Lab Sample ID (PIBLK):

Time Analyzed :\_\_\_\_\_

EPA Sample No. (PEM):

Date Analyzed :05/18/00

Lab Sample ID (PEM): EVALB

Time Analyzed:0315

PEM COMPOUND	RT	RT WI FROM	NDOW TO	CALC AMOUNT (ng)	NOM AMOUNT (ng)	%D
Endrin_ 4,4'-DDT_	13.47 14.51	13.43 14.46	13.53 14.56	0.024992 0.024711	0.025000	-0.0 -1.2

4,4'-DDT % breakdown (1):

121

Endrin % breakdown (1):

Combined % breakdown (1):

02,500

OLM03.0

## PESTICIDE CALIBRATION VERIFICATION SUMMARY

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT

Case No.:

SAS No.: 40325 SDG No.: SDGA18632

GC Column: RTX-CLP2 ID: 0.53 (mm) Init. Calib. Date(s): 05/17/00 05/18/00

EPA Sample No. (PIBLK):

Date Analyzed :\_\_\_\_\_

Lab Sample ID (PIBLK): \_\_\_\_

Time Analyzed:

EPA Sample No. (PEM):

Date Analyzed:05/19/00

Lab Sample ID (PEM): EVALB

Time Analyzed: 0832

PEM		RT W		CALC	NOM	
COMPOUND	RT	FROM	TO	AMOUNT (ng)	AMOUNT (ng)	%D
		=====	=====	======	=======	=====
Endrin	13.47	13.43	13.53	0.027357	0.025000	9.4
4,4'-DDT_	14.52	14.47	14.57	0.027652	0.025000	10.6
			<del></del>			
	ļ ———					<del></del>
						<del></del>
	<del></del>		<del></del>			

10.3

4,4'-DDT % breakdown (1):

10.85 Endrin % breakdown (1):

Combined % breakdown (1):

FORM VII PEST-1

OLM03.0

65744Pile: /var/chem/gc3.i/2170-E.b/c-b2362.d PT-UPZ
Report Date: 19-May-2000 11:21

STL-PITTSBURGH

# CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i

Lab File ID: c-b2362.d

Analysis Type:

Lab Sample ID: MEDTOX

Quant Type: ESTD

Injection Date: 19-MAY-2000 08:58
Init. Calibration Date(s): 05/17/0 05/18/0
Init. Calibration Times: 17:15 01:57

Method File: /var/chem/gc3.i/2170-E.b/PESTB.m

MIN MAX | RRF | %D | %D | RRF COMPOUND 52924.000 | 57579.000 | 0.010 | -8.8 | 15.0 | 18 Toxaphene(1) 83240.000 90902.000 0.010 -9.2 15.0 (2) 52341.000 57794.000 0.010 -10.4 15.0 (3) 52025.000 56303.000 0.010 -8.2 15.0 (4) | 4109568 000| 4440600.000|0.000| -8.1| 15.0| 1 Tetrachloro-m-xylene | 2705026 000| 2739320 000|0.010| -1.3| 15.0| \$ 30 Decachlorobiphenyl \_\_|\_\_\_|

HP58903F 657 421

Data File: /var/chem/gc3.i/2170-E.b/c-b2363.d

Report Date: 19-May-2000 11:22

STL-PITTSBURGH

## CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i Lab File ID: c-b2363.d

Analysis Type:

Lab Sample ID: MEDCHLOR

Quant Type: ESTD

Injection Date: 19-MAY-2000 09:24

Init. Calibration Date(s): 05/17/0 05/18/0
Init. Calibration Times: 17:15 01:57

Method File: /var/chem/gc3.i/2170-E.b/PESTB.m

ı_		1		MIN		MAX	1
ļ	COMPOUND	RRF	RF0	RRF	<b>₹</b> D	%D	1
1			222020222	==   ====			•
t	9 Chlordane(1)	157184.000	188652.0	00 0.010	-20.0	15.0	) <-
1	(2)	299340 000	354316.0	00 0.010	-18.4	15.0	) <-
1	(3)	431052.000	497972.0	00[0.010]	-15.5	15.0	)   <-
	(4)	354008.000	410640.0	00 0.010	-16 0	15 (	) <-
\$	1 Tetrachloro-m-xylene	4109568 000	4565520.0	000 0001	-11 1	15.0	1
į\$	30 Decachlorobiphenyl	2705026.000	2848960.0	00 0.010	-5.3	15.0	)
1_			l			!	-1

STL Pittsburgh

3107

72 HP58903F 657 428

Data File: /var/chem/gc3.i/2170-E.b/c-b2364.d

Report Date: 19-May-2000 11:22

STL-PITTSBURGH

# CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i Lab File ID: c-b2364.d Analysis Type:

Lab Sample ID: MEDA Quant Type: ESTD

Injection Date: 19-MAY-2000 09:50 Init. Calibration Date(s): 05/17/0 05/18/0 Init. Calibration Times: 17:15 01:57 Method File: /var/chem/gc3.i/2170-E.b/PESTB.m

		1 1	MIN	MAX
(	COMPOUND	RRF RF0	RRF   %	D   tD
	· "我们是我们们的,我们们们们们们们们们们们们们们们们们们们们们们们们们们们们们们们			bere
\$	l Tetrachloro-m-xylene	4109568.000 4682440.000	0.000  -1	3.9  15.0
	5 alpha-BHC	6370398 000  7432920.000	0.010  -1	6.7  15.0
	6 gamma-BHC (Lindane)	5385878 000  6451520.000	0.010  -1	9.8  15.0
	10 Heptachlor	5290246 000  6020560.000	0.010  -1	3.8 15.0
	15 Endosulfan I	4036632.000[ 4459520.000]	0.010  -1	0.5  15.0
	17 Dieldrin	4151588.000  4690120.000	0.010  -1	3.0  15.0
	20 Endrin	3539938.000  3964920.000	0.010  -1	2.0  15.0
	21 4,4'-DDD	3011574.000 3314080 000	0.010  -1	0.0  15.0
	23 4,4'-DDT	2620220.000  2841320.000	0.010  -	8.4  15.0
	25 Methoxychlor	1250013.000  1333880.000	0 010  -	6 7   15 0
\$	30 Decachlorobiphenyl	2705026.000  2868680.000	0.010  -	6.0  15 0
		l		

7E HP58903F

Data File: /var/chem/gc3.i/2170-E.b/c-b2365.d @7-ccl2 657 429

Report Date: 19-May-2000 11:22

## STL-PITTSBURGH

# CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i Lab File ID: c-b2365.d

Analysis Type:

Lab Sample ID: MEDB Quant Type: ESTD Injection Date: 19-MAY-2000 10:16

Init. Calibration Date(s): 05/17/0 05/18/0
Init. Calibration Times: 17:15 01:57
Method File: /var/chem/gc3.i/2170-E.b/PESTB.m

I ppp			
RRF	RF0	RRF	%D   %D
5126234.000	5699320.00	00 0 010	-11.2  15.0
5545788.000	6222240.0	00(0.010	-12.2  15.
4383022.000	4851240.0	00 0 010	-10 7] 15
4593890.000	5067720 0	00 0 010	-10.3  15
4477596.000	4931960 0	00 0.010	-10.1 15.
3926194.000	4340120.0	00 0 010	-10.5  15
3413566.000	3693480.0	00 0.010	-8.2  15.
2780892 000	2978880 0	00 0.010	-7.1 15.
2812870.000	3058640.0	00 0.010	-8.7  15.
3086674 000	3281600.0	00 0.010	-6.3  15.
	5126234.000   2737018.000   5545788.000   4383022.000   4593890.000   4477596.000   3926194.000   3413566.000   2780892.000	5126234.000   5699320.00   2737018.000   3000920 00   5545788.000   6222240.0   4383022.000   4851240.0   4593890.000   5067720 0   4477596.000   4931960 0   3926194.000   4340120.0   3413566.000   3693480.0   2780892 000   2978880 0   2812870.000   3058640.0	5126234.000   5699320.000   0 010   2737018.000   3000920 000   0.010   5545788.000   6222240.000   0 010   4383022.000   4851240.000   0 010   4593890.000   5067720 000   0 010   4477596.000   4931960 000   0.010   3926194.000   4340120.000   0 010   3413566.000   3693480.000   0.010   2780892 000   2978880 000   0.010   2812870.000   3058640.000   0.010   3086674 000   3281600.000   0.010

Aus 10.77,0

3109

657 430

Data File: /var/chem/gc3.i/2170-E.b/c-b2374.d

Report Date: 19-May-2000 14:32

7E HP58903F

## STL-PITTSBURGH

## CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i Lab File ID: c-b2374.d

Analysis Type:

Lab Sample ID: MEDA Quant Type: ESTD

Injection Date: 19-MAY-2000 14:10

Init. Calibration Date(s): 05/17/0 05/18/0 Init. Calibration Times: 17:15 01:57 Method File: /var/chem/gc3.i/2170-E.b/PESTB.m

		l		MIN	[ :	MAX
	COMPOUND	RRF	RF0	RRF	%D	<b></b> ₽D
<b>.</b> 12	*********************				-	
3	1 Tetrachloro-m-xylene	4109568.00	0  4778800.0	000.000	-16.3	15.0
	5 alpha-BHC	6370398 00	0  7537800.00	00 0.010	-18.3	15.0
	6 gamma-BHC (Lindane)	5385878.00	0 6495720 0	00 0.010	-20 6	15.0
	10 Heptachlor	5290246.00	0  6159640.00	00 0.010	-16.4	15.0
	15 Endosulfan I	4036632.00	0  4520320.00	00[0.010]	-12.0	15.0
	17 Dieldrin	4151588.00	0 4736360.00	00 0.610	-14.1[	15.0
	20 Endrin	3539938.00	0 4072840.00	00 0 010	-15 1	15 0
	21 4,4'-DDD	3011574 00	0  3298080.00	000 010	-9 5	15.0
	23 4,4'-DDT	2620220.00	0 2967680.00	00[0.010]	-13.3	15.0
	25 Methoxychlor	1250013.00	0   1401320.00	00 0.010	-12 1	15.0
:	30 Decachlorobiphenyl	2705026.00	0  2948880.00	00 0.010	-9.0	15 0

75.d RT-CUPZ

Data File: /var/chem/gc3.i/2170-E.b/c-b2375.d Report Date: 19-May-2000 15:21

657 431

## STL-PITTSBURGH

## CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gc3.i Lab File ID: c-b2375.d

Analysis Type:

Lab Sample ID: MEDB Quant Type: ESTD

Injection Date: 19-MAY-2000 14:36 Init. Calibration Date(s): 05/17/0 05/18/0

Init. Calibration Times: 17:15 01:57
Method File: /var/chem/gc3.i/2170-E.b/PESTB.m

,	COMPOUND	RRF	   RFO	MIN     MAX   RRF   %D   %D
	BB BB BB BB BB BB AB AB AB BB BB BB BB B			.   =====   ====
	11 Aldrin	5126234.000	5772680 000	0 0.010  -12.6  15.
	7 beta-BHC	2737018 000	3048720.000	0 0.010  -11.4  15.
	8 delta-BHC	5545788.000	6307880.000	0 0.010  -13.7  15
	12 Heptachlor epoxide			0 0.010  -10.7  15.
	13 gamma-Chlordane	4593890.000	5110760 000	0 0.010  -11.3  15
	14 alpha-Chlordane	4477596.000	5000800.000	0 0.010  -11 7  15
	16 4,4'-DDE	3926194.000	4404560.000	0 0.010  -12.2  15.
	22 Endosulfan II			0 0.010  -10.3  15.
	24 Endrin aldehyde			0 0.010  -9 0  15.
	26 Endosulfan sulfate			0 0.010  -11.4  15.
	27 Endrin ketone		-	0 0.010  -10 3  15

Ay 12:87.D

### 8D PESTICIDE ANALYTICAL SEQUENCE

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.:

SAS No.: 40325 SDG No.: C0E130194

GC Column: RTX-CLP2 ID: 0.53 (mm) Init. Calib. Date(s): 05/17/00 05/18/00

Instrument ID: GC3

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS, SAMPLES, AND STANDARDS IS GIVEN BELOW:

	MEAN SURROC TCX: 5.97		NITIAL CALI .8.82	BRATION		
-	EPA	LAB	DATE	TIME	TCX	DCB
1	SAMPLE NO.	SAMPLE ID	ANALYZED	ANALYZED	RT #	RT #
			=======================================	========	========	=======
01		EVALB	05/17/00	1649	5.98	18.83
02		MEDTOX	05/17/00	1715	5.98	18.83
03	•	MEDCHLOR	05/17/00	1741	5.99	18.83
04		LOWAPPX9	05/17/00	1807	]	
05	() <1) Rt	MLOWAPPX9	05/17/00	1833		
06	Carron	MEDAPPX9	05/17/00	1859		
07		MH1GHAPPX9	05/17/00	1925		
08		HTCHAPDX9	05/17/00	1951		
09	_	LOWA	05/17/00	2017	5.99	18.83
10		MLOWA	05/17/00	2230	5.99	18.84
11		MEDA	05/17/00	2256	5.99	18.84
12		MHIGHA	05/17/00	2322	5.99	18.84
13		HIGHA	05/17/00	2348	5.99	18.83
14		LOWB	05/18/00	0014		
15		MLOWB	05/18/00	0040		
16		MEDB	05/18/00	0106		
17		MHIGHB	05/18/00	0132		
18		HIGHB	05/18/00	0157		
19		2ND A	05/18/00	0223	5.99	18.83
20		2ND B	05/18/00	0249		
21		EVALB	05/18/00	0315	5.99	18.84
22		EVALB	05/19/00	0832	6.00	18.84
23		MEDTOX	05/19/00	0858	6.00	18.84
24		MEDCHLOR	05/19/00	0924	6.00	18.84
25		MEDA	05/19/00	0950	6.00	18.84
26		MEDB	05/19/00	1016		
	DF/S1/0133/S		05/19/00	1042	6.00	18.84
	DF/S1/0133/S		05/19/00	1108	6.00	18.84
	DF/S1/0133/S		05/19/00	1134	6.00	18.84
	DF/S1/0133/S		05/19/00	1200	6.00	18.84
	DF/S1/0133/S	DD6A6103	05/19/00	1226	5.99	18.84
	DF/S1/0133/S	DD6A7103	05/19/00	1252	6.00	18.84
	,,, -					

QC LIMITS

page 1 of 2

FORM VIII PEST

OLM03.0

TCX = Tetrachloro-m-xylene (+/- 0.05 MINUTES)
DCB = Decachlorobiphenyl (+/- 0.05 MINUTES)

<sup>#</sup> Column used to flag retention time values with an asterisk.

<sup>\*</sup> Values outside of QC limits.

## 8D PESTICIDE ANALYTICAL SEQUENCE

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT

Case No.:

SAS No.: 40325

SDG No.: C0E130194

GC Column: RTX-CLP2 ID: 0.53 (mm) Init. Calib. Date(s): 05/17/00 05/18/00

Instrument ID: GC3

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS, SAMPLES, AND STANDARDS IS GIVEN BELOW:

	MEAN SURRO					
	EPA	DCB: 1	DATE	TIME	TCX	DCB
	SAMPLE NO.	SAMPLE ID	ANALYZED	ANALYZED	RT #	RT #
01 02 03 04 05	PBLK ZZZZZ	DDE2E101 ZZZZZ MEDA MEDB	05/19/00 05/19/00 05/19/00 05/19/00	1318 1344 1410 1436	6.00 6.00 6.00	18.84 18.84 18.84
06 07 08						
09 10 11						
12 13						
14 15 16						
17 18 19						
20 21 22						
23 24						
25 26 27						
28 29						
30 31 32						

QC LIMITS

TCX = Tetrachloro-m-xylene (+/- 0.05 MINUTES)

DCB = Decachlorobiphenyl (+/- 0.05 MINUTES)

# Column used to flag retention time values with an asterisk.

\* Values outside of QC limits.

page 2 of 2

FORM VIII PEST

OLM03.0

Data File: /var/chem/gc3.i/2170-E.b/c-b2281.d

Report Date: 18-May-2000 11:44

## STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2281.d

Lab Smp Id: EVALB

Inj Date : 17-MAY-2000 16:49

Operator : 1891 Inst ID: gc3.i

Smp Info : EVALB, 2170-E.b, , EVALBR. sub, , 3, 1

Misc Info: 190-88-8

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date: 18-May-2000 11:42 g Quant Type: ESTD Cal File: c-b2300.d Cal Date : 18-MAY-2000 01:57

QC Sample: PEM Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: EVALBR.sub Integrator: Falcon

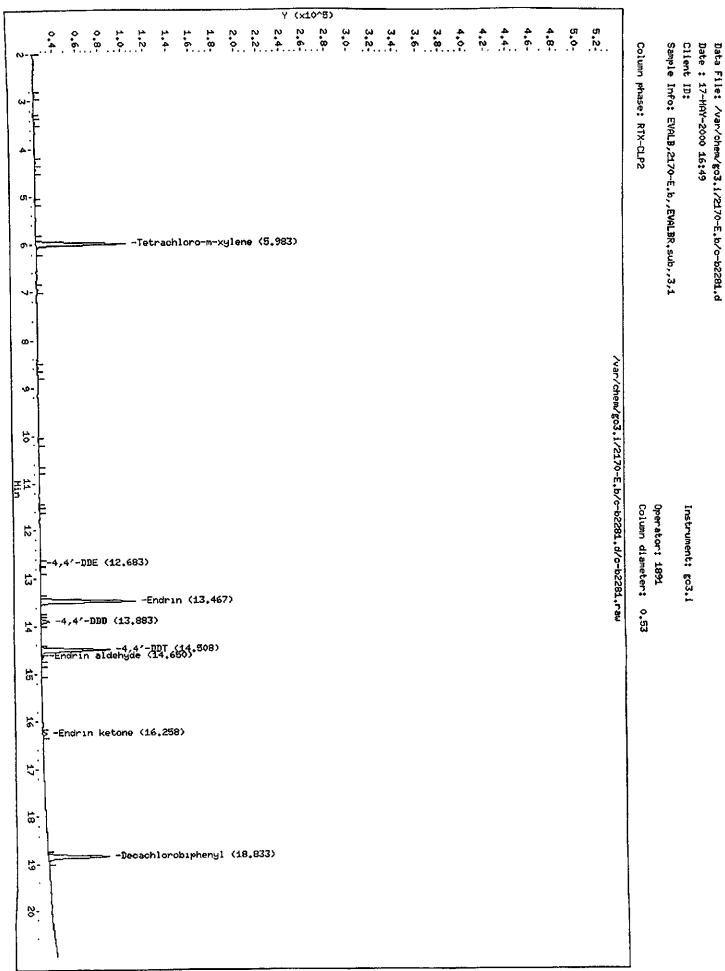
Target Version: 3.40

					CONCENT	RATIONS
					ON-COLUMN	FINAL
Compounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
	==	22222	200205	z=======	H25534E	그그고구르는
§ 1 Tetrachloro-m-xylene	5.983	5.992	-0.009	79968	0 01946	0.0194590(R)
16 4,4'-DDB	12.683	12.692	-0.009	709	0.000181	0.000180582
20 Bndrin	13.467	13.475	-0.008	84683	0 02392	0.0239222
21 4,4'-DDD	13.883	13.892	-0.009	7787	0 00259	0.00258569
23 4,4'-DDT	14.509	14.508	0,000	61222	0.02337	0.0233652
24 Endrin aldehyde	14.650	14 658	-0.008	1412	0.000508	0.000507751
27 Endrin ketone	16.258	16.267	-0.009	3996	0.00129	0.00129460
\$ 30 Decachlorobiphenyl	18.833	18.842	-0.009	53341	0 01972	0.0197192(R)

# QC Flag Legend

R - Spike/Surrogate failed recovery limits.

00 12.200 GNORIN 6.010



Data File: /var/chem/gc3.i/2170-E.b/c-b2282.d

Report Date: 18-May-2000 11:44

### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2282.d

Lab Smp Id: MEDTOX Inj Date : 17-MAY-2000 17:15

Inst ID: gc3.i Operator: 1891

Smp Info : MEDTOX,2170-E.b,,1-TOX.sub,,1,3
Misc Info : 190-84-13

Comment

: /var/chem/gc3.i/2170-E.b/PESTB.m Method

Quant Type: ESTD Meth Date : 18-May-2000 11:42 g Cal Date : 17-MAY-2000 17:15 Cal File: c-b2282.d

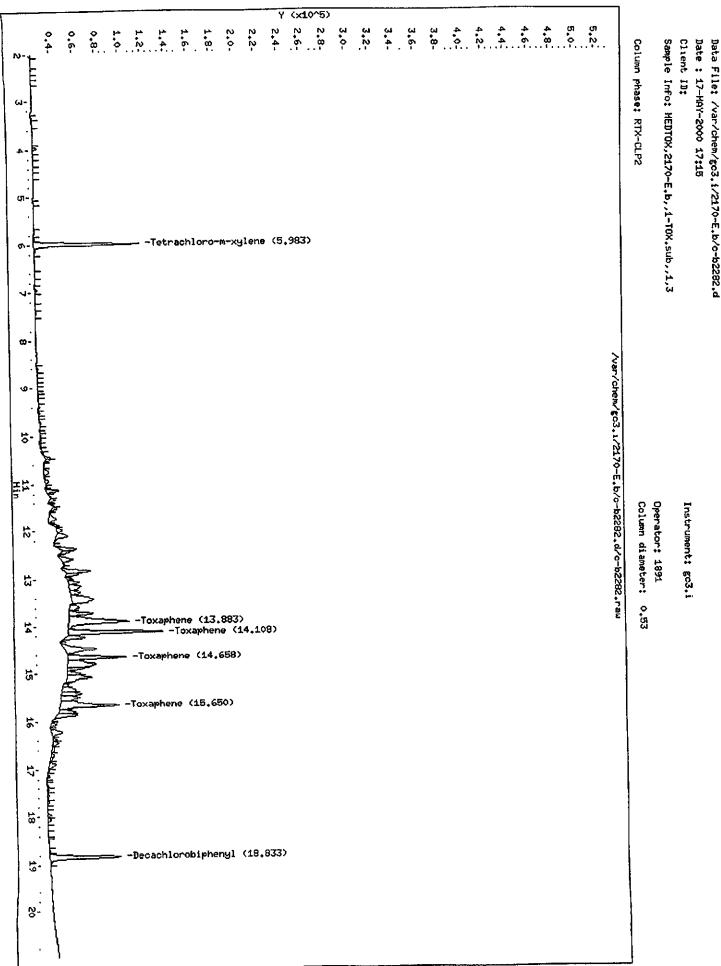
Calibration Sample, Level: 3 Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: 1-TOX.sub Integrator: Falcon

Target Version: 3.40

				amounts		
				CAL-AMT	ON-COL	
Compounds	RT	BXP RT DLT RT	RESPONSE	( ng)	( ng)	
	==	<b>2</b> 0	****	======		
18 Toxaphene	13.883	13.883 0.000	52924	1.00000	1.00000	
s 1 Tetrachloro-m-xylene	5.983	5.992 -0.009	93378	0.02500	0.0250000	
\$ 30 Decachlorobiphenyl	18.633	18.842 -0.009	62869	0.02500	0.0250000	



Data File: /var/chem/gc3.i/2170-E.b/c-b2283.d

Report Date: 18-May-2000 11:44

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-b2283.d

Lab Smp Id: MEDCHLOR

Inj Date : 17-MAY-2000 17:41 Operator : 1891 Inst ID: gc3.i

Smp Info : MEDCHLOR, 2170-E.b., 2-CHLO.sub, ,1,3

Misc Info: 190-85-10

Comment

: /var/chem/gc3.i/2170-E.b/PESTB.m Method

Quant Type: ESTD Meth Date : 18-May-2000 11:42 g Cal File: c-b2283.d Cal Date : 17-MAY-2000 17:41

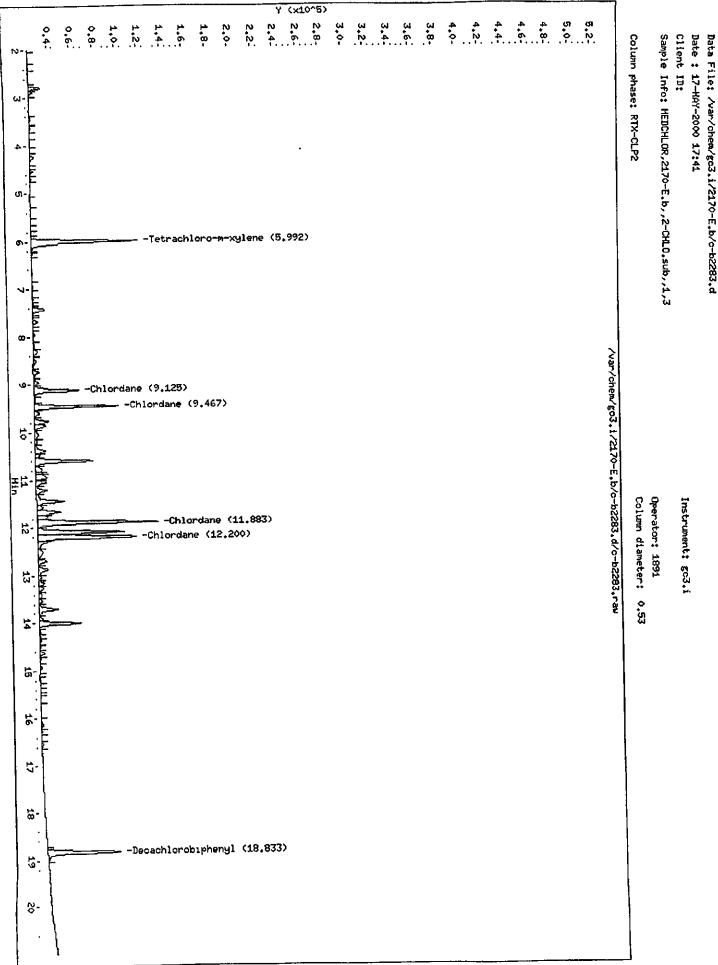
Calibration Sample, Level: 3 Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: 2-CHLO.sub Integrator: Falcon

Target Version: 3.40

					amounts			
					CAL-	AMT	ON	-col
Compounds	RT	EXP RT DL	T RT	RESPONSE	(	ng)	(	ng)
2.2.2.2.2.2.2.2.2.2.2.2.2.2.2.2.2.2.2.		222222 23	====		===:	2222	==	****
9 Chlordane	9 125	9 125	0.000	39296	0.2	5000	0.2	50000
\$ 1 Tetrachloro-m-xylene	5.992	5.992	0.000	94430	0.0	2500	0.02	50000
\$ 30 Decachlorobiphenyl	18.833	18.842 -	0.009	64237	0 0:	2500	0 02	250000



Data File: /var/chem/gc3.i/2170-E.b/c-b2289.d

Report Date: 18-May-2000 11:44

## STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2289.d

Lab Smp Id: LOWA

Inj Date : 17-MAY-2000 20:17

Operator: 1891 Inst ID: gc3.i

Smp Info : LOWA, 2170-E.b, , 3-INDA.sub, , 1, 1

Misc Info : 190-84-1

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date: 18-May-2000 11:42 g Quant Type: ESTD Cal Date: 17-MAY-2000 20:17 Cal File: c-b2289.d

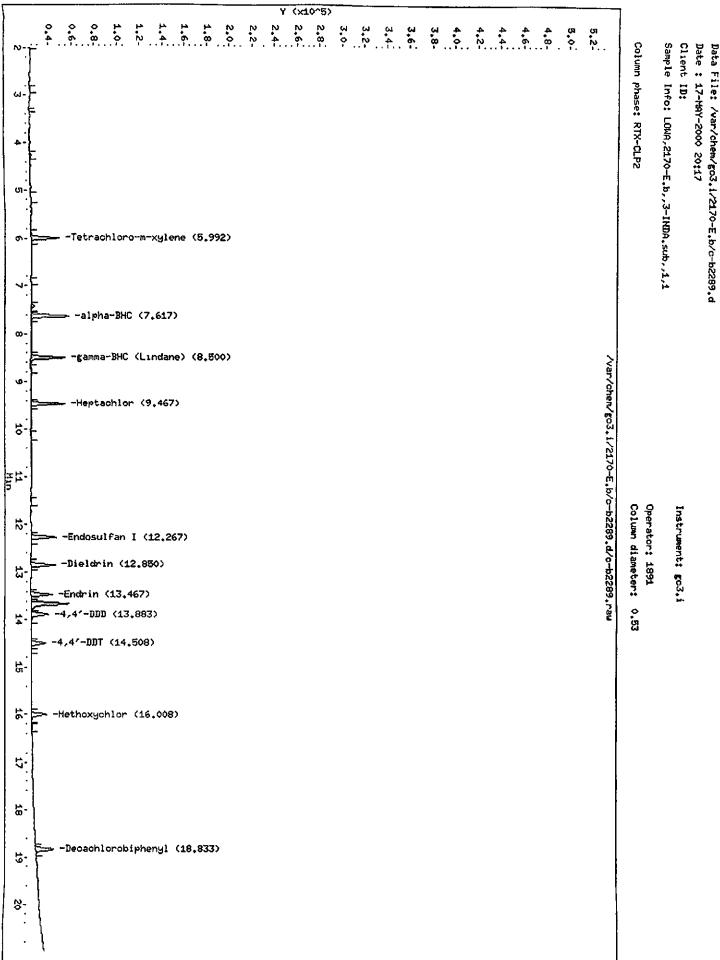
Als bottle: 1 Calibration Sample, Level: 1

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 3-INDA.sub

Target Version: 3.40

						AMOUNTS	
						CAL-AMT	ON-COL
Com	pounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
===	*	==	202=5 <b>2</b>	======	*****	2227555	*****
Ś	1 Tetrachloro-m-xylene	5.992	5 992	0 000	24804	0.00500	0.00567727
•	5 alpha-BHC	7 617	7 625	-0.008	33617	0.00500	0.00500000
	6 gamma-BHC (Lindane)	8.500	8.508	-0.008	29937	0.00500	0.00500000
	10 Heptachlor	9.467	9.475	-0 008	29711	0.00500	0.00500000
	15 Endosulfan I	12 267	12.275	-0 008	22482	0 00500	0.00500000
	17 Dieldrin	12.850	12 858	-0.008	21946	0.00500	0.00500000
	20 Bndrin	13 467	13 475	-0.00B	18774	0.00500	0.00500000
	21 4,4'-DDD	13.883	13.892	-0 009	15526	0.00500	0.00500000
	23 4,4'-DDT	14.508	14.508	0 000	12923	0.00500	0.00500000
	25 Methoxychlor	16.008	16.008	0.000	13407	0.01000	0 0100000
\$	30 Decachlorobiphenyl	18.833	18.842	-0 009	15693	0.00500	0.00549852



Data File: /var/chem/gc3.i/2170-E.b/c-b2292.d

Report Date: 18-May-2000 11:44

## STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2292.d Lab Smp Id: MLOWA

Inj Date : 17-MAY-2000 22:30

Inst ID: gc3.i

Operator: 1891 In: Smp Info: MLOWA, 2170-E.b,,3-INDA.sub,,1,2

Misc Info : 190-84-2

Comment

: /var/chem/gc3.i/2170-E.b/PESTB.m Method

Quant Type: ESTD Meth Date: 18-May-2000 11:42 g Cal File: c-b2292.d Cal Date : 17-MAY-2000 22:30

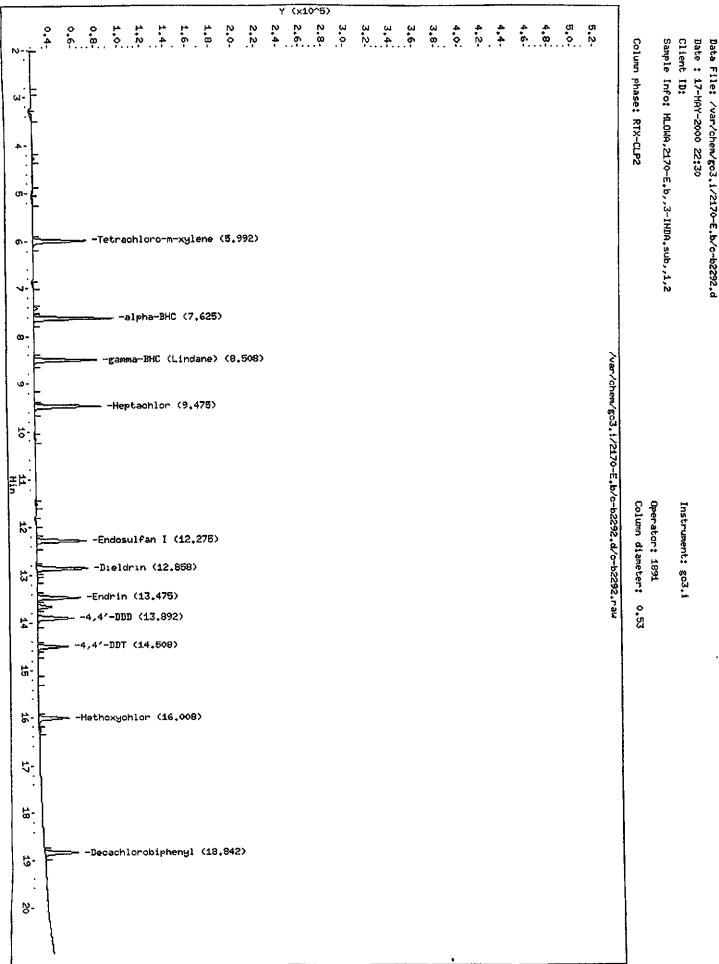
Calibration Sample, Level: 2 Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: 3-INDA.sub Integrator: Falcon

Target Version: 3.40

					amounts		
					CAL-AMT	ON-COL	
Compounds	RT	BXP RT	DLT RT	response	( ng)	(ng)	
	==	<b>55052</b>	828232	===#####	2442777	ge=3553	
\$ 1 Tetrachloro-m-xylene	5.992	5.992	0.000	46864	0 01000	0 0104729	
S alpha-BHC	7.625	7 625	0.000	69788	0.01000	0.0101864	
6 gamma-BHC (Lindane)	8.508	B 508	0.000	55992	0 01000	0.00966496	
10 Heptachlor	9.475	9 475	0.000	58817	0.01000	0.00994883	
15 Endosulfan I	12.275	12 275	0.000	44778	0.01000	0.00997927	
17 Dieldrin	12.858	12 858	0 000	45413	0.01000	0.0101703	
20 Endrin	13.475	13 475	0.000	38150	0.01000	0.0100795	
21 4,4'-DDD	13.892	13.892	0 000	31973	0.01000	0.0101461	
23 4,4'-DDT	14.508	14.508	0 000	27033	0.01000	0.0102245	
	16.008	16.008	0.000	26936	0.02000	0 0200454	
25 Methoxychlor \$ 30 Decachlorobiphenyl	18.842	18 842	0.000	29730	0.01000	0.0102741	



Data File: /var/chem/gc3.i/2170-E.b/c-b2293.d

Report Date: 18-May-2000 11:44

### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2293.d

Lab Smp Id: MEDA

Inj Date : 17-MAY-2000 22:56

Operator: 1891 Inst ID: gc3.i

Smp Info : MEDA, 2170-E.b, , 3-INDA.sub, , 1, 3

Misc Info: 190-84-3

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date : 18-May-2000 11:42 g Quant Type: ESTD Cal Date : 17-MAY-2000 22:56 Cal File: c-b2293.d

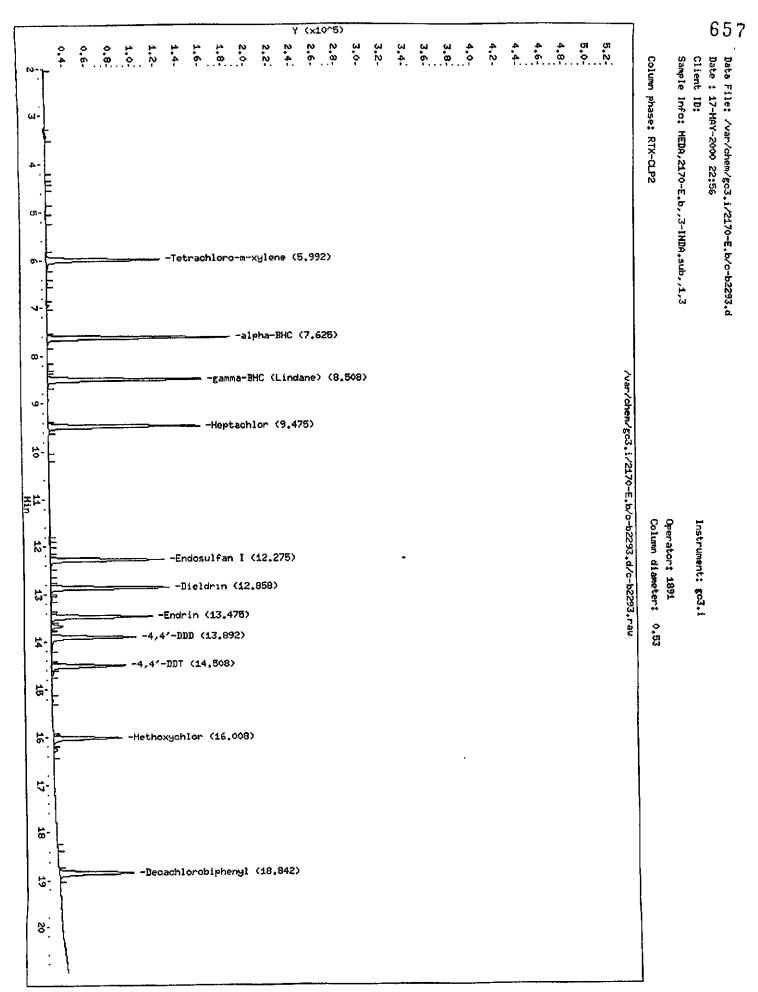
Als bottle: 1 Calibration Sample, Level: 3

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 3-INDA.sub

Target Version: 3.40

						AMOUNTS		
						CAL-AMT	ON-COL	
Col	mpounda	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)	
	- 医乳口状心炎 - 医克森氏 医甲甲基甲基磺胺 经现在日日日	33	202222		=======	5022007	******	
\$	1 Tetrachloro-m-xylene	5 992	5.992	0.000	100176	0.02500	0.0220099	
	5 alpha-BHC	7 625	7.625	0.000	160664	0.02500	0.0239454	
	6 gamma-BHC (Lindane)	8.508	8.508	0.000	135560	0.02500	0.0239097	
	10 Heptachlor	9 475	9.475	0.000	133248	0.02500	0.0233035	
	15 Endosulfan I	12.275	12.275	0.000	100629	0.02500	0.0232232	
	17 Dieldrin	12.858	12.858	0.000	104542	0.02500	0.0239187	
	20 Endrin	13 475	13.475	0.000	89197	0.02500	0.0240257	
	21 4,4'~DDD	13,892	13.892	0.000	74538	0.02500	0.0240859	
	23 4,4'-DDT	14.508	14.508	0.000	65341	0.02500	0.0248082	
	25 Methoxychlor	16.008	16,008	0 000	61025	0.05000	0.0468462	
\$	30 Decachlorobiphenyl	18.842	18.842	0 000	65867	0.02500	0 0225926	



Data Filet /var/chem/gc3.i/2170-E.b/c-b2294.d

Report Date: 18-May-2000 11:44

### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2294.d

Lab Smp Id: MHIGHA

Inj Date : 17-MAY-2000 23:22

Operator: 1891 Inst ID: gc3.i

Smp Info : MHIGHA, 2170-E.b,, 3-INDA.sub,, 1,4

Misc Info : 190-84-4

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date : 18-May-2000 11:42 g Quant Type: ESTD Cal Date : 17-MAY-2000 23:22 Cal File: c-b2294.d

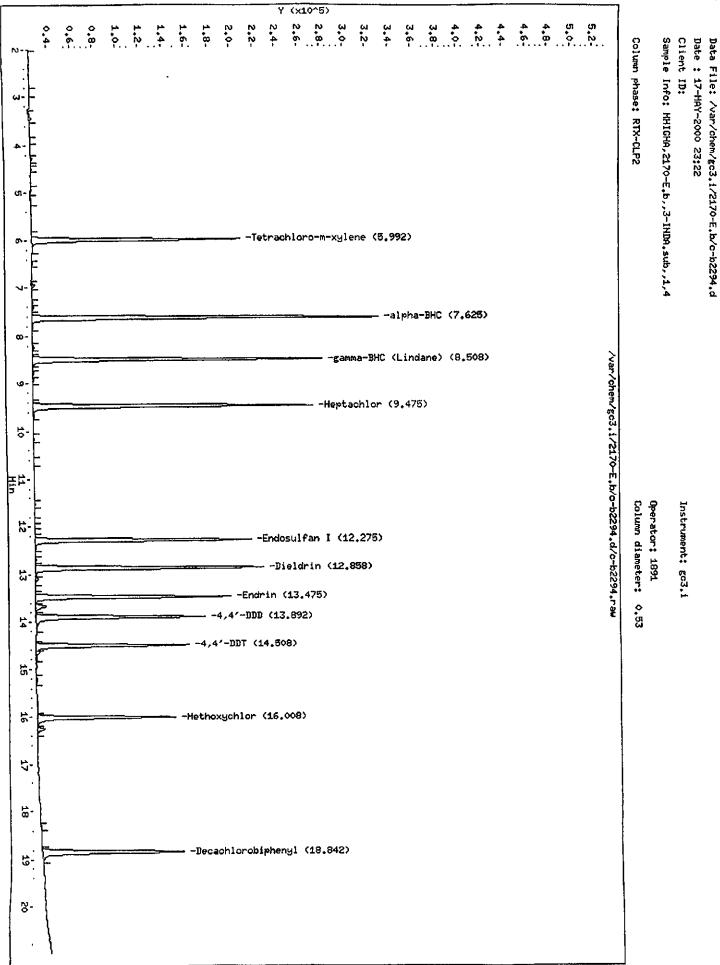
Als bottle: 1 Calibration Sample, Level: 4

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 3-INDA.sub

Target Version: 3.40

							AMOUNTS		
							CAL-AMT	ON-COL	
Cor	npot	ınds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)	
==:	_	##2==#################################	3E	*****	*****	********		******	
\$	1	Tetrachloro-m-xylene	5.992	5 992	0.000	183640	0.05000	0.0423938	
	5	alpha-BHC	7,625	7.625	0.000	306572	0.05000	0.0466976	
	6	gamma-BHC (Lindane)	8.508	8.508	0.000	255751	0.05000	0.0462395	
	10	Heptachlor	9.475	9.475	0.000	247550	0.05000	0.0447956	
	15	Endosulfan I	12.275	12.275	0 000	191746	0.05000	0.0455608	
	17	Dieldrin	12.858	12.858	0.000	202395	0.05000	0.0471781	
	20	Endrin	13.475	13.475	0.000	172892	0 05000	0 0473822	
	21	4,4'-DDD	13.892	13.892	0.000	149519	0.05000	0.0487255	
	23	4,4'-DDT	14.508	14.508	0.000	134807	0.05000	0.0508817	
	25	Methoxychlor	16.008	16.008	0.000	121669	0.10000	0.0949669	
\$	30	Decachlorobiphenyl	18 842	18.842	0.000	124669	0.05000	0.0443675	



657 448
Data File: /var/chem/gc3.i/2170-E.b/c-b2295.d

Report Date: 18-May-2000 11:44

### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2295.d

Lab Smp Id: HIGHA

Inj Date : 17-MAY-2000 23:48

Operator : 1891 Inst ID: gc3.i

Smp Info : HIGHA, 2170-E.b,, 3-INDA.sub,, 1,5

Misc Info : 190-84-5

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date : 18-May-2000 11:42 g Quant Type: ESTD Cal File: c-b2295.d

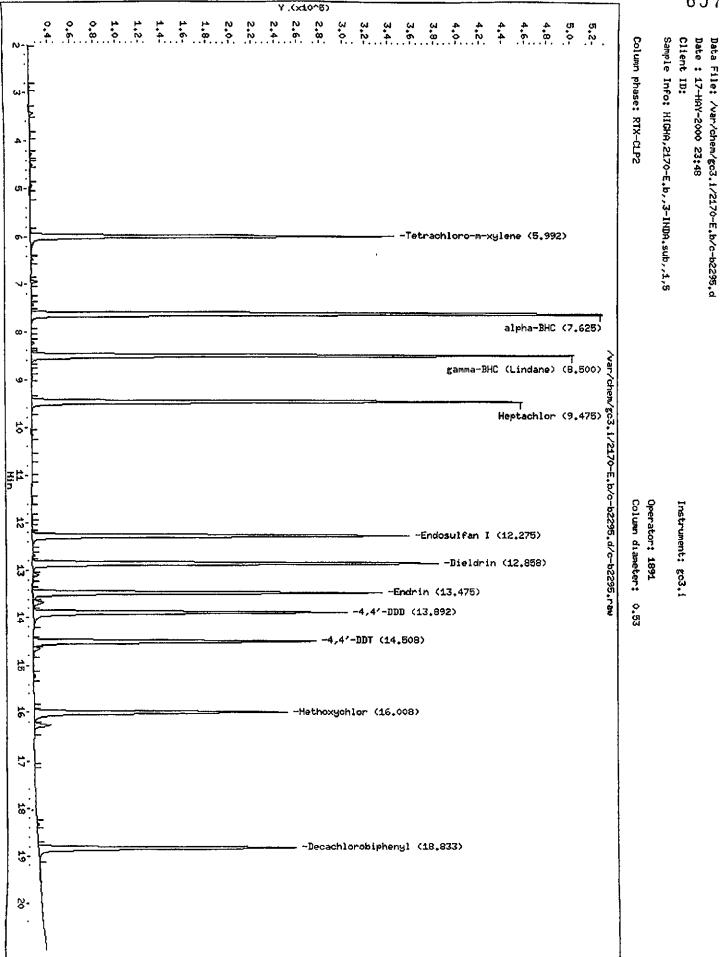
Cal Date : 17-MAY-2000 23:48 Calibration Sample, Level: 5 Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: 3-INDA.sub Integrator: Falcon

Target Version: 3.40

					AMOUNTS		
					CAL-AMT	ON-COL	
Compounds	RT	EXP RT	DLT RT	response	( ng)	( ng)	
	==			##=7=###	= 급 저 의 로 의 로	222224	
s 1 Tetrachloro-m-xylene	5.992	5.992	0.000	322080	0.10000	0.0783732	
S alpha-BHC	7.625	7.625	0.000	559179	0.10000	0.0877777	
6 gamma-BHC (Lindane)	8.500	8.508	-0.008	480537	0.10000	0.0892217	
10 Heptachlor	9.475	9.475	0 000	434641	0 10000	0.0821589	
15 Endosulfan I	12.275	12.275	0.000	334888	0 10000	0.0829622	
17 Dieldrin	12.858	12.858	0 000	359786	0.10000	0.0866623	
20 Endrin	13.475	13.475	0.000	310417	0 10000	0.0876900	
21 4,4'-DDD	13.892	13.892	0.000	278347	0.10000	0 0924258	
23 4,4'-DDT	14.508	14.508	0.000	250342	0.10000	0.0955424	
25 Methoxychlor	16.008	16.008	0.000	225075	0.20000	0.180058	
\$ 30 Decachlorobiphenyl	18.833	18.842	-0.009	228547	0.10000	0.0844898	



657 450 Data File: /var/chem/gc3.i/2170-E.b/c-b2296.d

Report Date: 18-May-2000 11:44

## STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2296.d

Lab Smp Id: LOWB

Inj Date : 18-MAY-2000 00:14 Operator : 1891 Inst ID: gc3.i

Smp Info : LOWB, 2170-E.b., 4-INDB. sub, ,1,1

Misc Info : 190-84-7

Comment : Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date : 18-May-2000 11:42 g Quant Type: ESTD

Cal File: c-b2296.d Cal Date : 18-MAY-2000 00:14

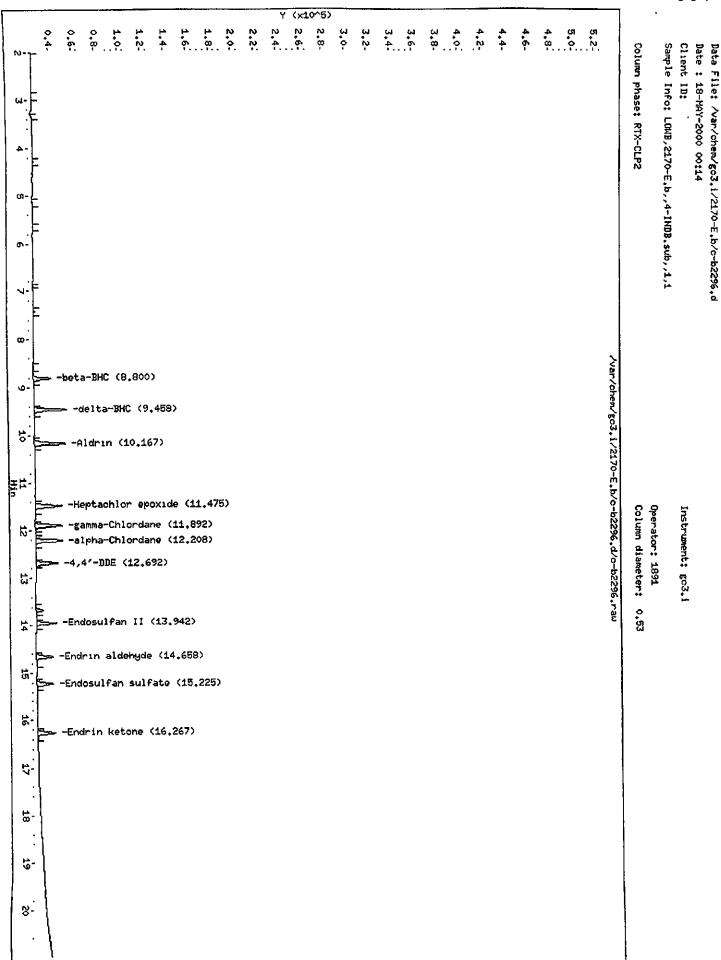
Calibration Sample, Level: 1 Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: 4-INDB.sub Integrator: Falcon

Target Version: 3.40

					AMOUNTS	
					CAL-AMT	ON-COL
Compounds	RT	BXP RT	DLT RT	RESPONSE	( ng)	( ng)
		ER###	DP=====	******	=======	******
11 Aldrin	10.167	10.167	0.000	27894	0.00500	0 00500000
7 beta-BHC	8.800	8.800	0.000	15896	0.00500	0.00500000
8 delta-BHC	9.458	9,458	0.000	28706	0.00500	0.00500000
12 Heptachlor epoxide	11.475	11 475	0.000	24110	0.00500	0.00500000
13 gamma-Chlordane	11 892	11 892	0 000	24762	0.00500	0.00500000
14 alpha-Chlordane	12.208	12 208	0.000	24400	0.00500	0.00500000
16 4,4'-DDB	12 692	12.692	0.000	20059	0.00500	0.00500000
22 Endosulfan II	13.942	13.942	0.000	18109	0.00500	0.00500000
24 Endrin aldehyde	14 658	14.658	0.000	15202	<b>0.00500</b>	0.00500000
26 Endosulfan sulfate	15.225	15.225	0.000	14592	0 00500	0.00500000
27 Endrin ketone	16.267	16.267	0.000	16130	0 00500	0.00500000



Data File: /var/chem/gc3.i/2170-E.b/c-b2297.d

Report Date: 18-May-2000 11:45

## STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2297.d

Lab Smp Id: MLOWB

Inj Date : 18-MAY-2000 00:40

Inst ID: gc3.i Operator: 1891

Smp Info : MLOWB, 2170-E.b, , 4-INDB. sub, , 1, 2

Misc Info : 190-84-8

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTB.m Meth Date : 18-May-2000 11:42 g Quar Quant Type: ESTD Cal File: c-b2297.d Cal Date : 18-MAY-2000 00:40

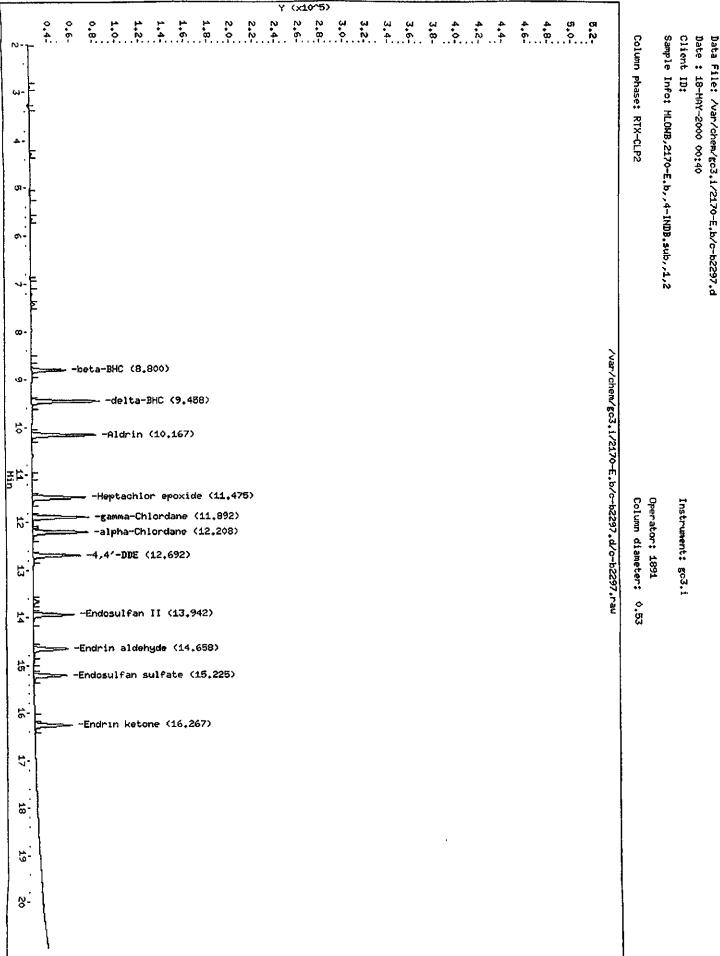
Calibration Sample, Level: 2 Als bottle: 1

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 4-INDB.sub

Target Version: 3.40

					AMOUNTS	
					CAL-AMT	ON-COL
Compounds	RT	EXP RT	DLT RT	rbsponsb	( ng)	( ng)
	20	35### <b>#</b>	*****	042222	***	3=====
11 Aldrin	10.167	10.167	0.000	56697	0 01000	0.0100808
7 beta-BHC	8.800	8.800	0 000	30704	0.01000	0.00982591
8 delta-BHC	9.458	9 458	0.000	59691	0.01000	0 0101946
12 Heptachlor epoxide	11.475	11.475	0.000	47916	0.01000	D 00996838
13 gamma-Chlordane	11.892	11 892	0 000	50094	0 01000	0 0100572
14 alpha-Chlordane	12.208	12.208	0.000	49070	0.01000	0.0100276
16 4,4'-DDE	12.692	12.692	0.000	42223	0.01000	0.0102556
22 Endosulfan II	13.942	13.942	0.000	36744	0 01000	0.0100721
24 Endrin aldehyde	14 658	14 658	0.000	30722	0 01000	0.0100520
26 Endosulfan sulfate	15 225	15.225	0.000	30063	0 01000	0.0101484
27 Endrin ketone	16 267	16.267	0.000	33517	0.01000	0.0101911



Data File: /var/chem/gc3.i/2170-E.b/c-b2298.d

Report Date: 18-May-2000 11:45

### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2298.d

Lab Smp Id: MEDB

Inj Date : 18-MAY-2000 01:06

Operator: 1891 Inst ID: gc3.i

Smp Info : MEDB, 2170-E.b, ,4-INDB.sub, ,1,3

Misc Info: 190-84-9

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date: 18-May-2000 11:42 g Quant Type: ESTD Cal Date: 18-MAY-2000 01:06 Cal File: c-b2298.d

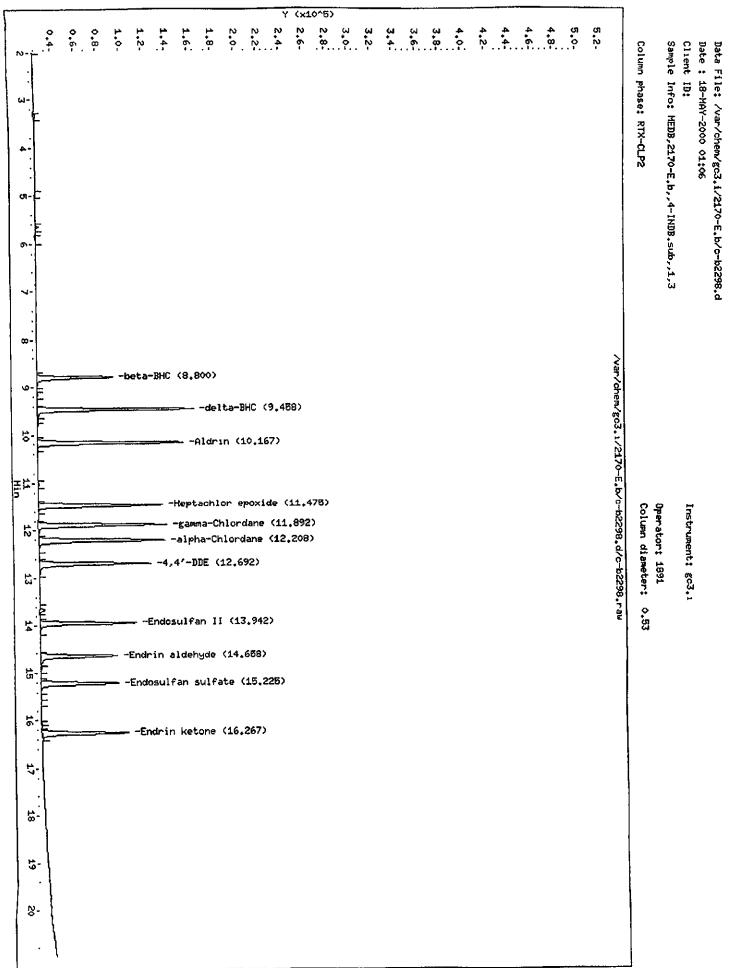
Als bottle: 1 Calibration Sample, Level: 3

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 4-INDB.sub

Target Version: 3.40

					AMOU	NTS
					CAL-AMT	ON-COL
Compounds	RT	EXP RT	DLT RT	response	( ng)	( ng)
	**	2020==	****	******	277722±	
11 Aldrin	10.167	10.167	0.000	127877	0.02500	0.0234442
7 beta-BHC	8.800	8.800	0.000	66768	0.02500	0.0224548
8 delta-BHC	9.458	9.458	0.000	137836	0.02500	0.0240080
12 Heptachlor epoxide	11.475	11.475	0.000	109575	0.02500	0.0234861
13 gamma-Chlordane	11 892	11.892	0.000	113875	0 02500	0.0235331
14 alpha-Chlordane	12.208	12.208	0.000	111661	0 02500	0.0235019
16 4,4'-DDE	12.692	12.692	0.000	98517	0.02500	0.0242757
22 Endosulfan II	13.942	13.942	0.000	84906	0 02500	0.0238222
24 Endrin aldehyde	14.658	14.658	0 000	67736	0.02500	0.0230341
26 Bndosulfan sulfate	15 225	15.225	0.000	69174	0.02500	0 0238760
27 Endrin ketone	16.267	16.267	0.000	76419	0.02500	0.0237955



Data File: /var/chem/gc3.i/2170-E.b/c-b2299.d

Report Date: 18-May-2000 11:45

### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-b2299.d

Lab Smp Id: MHIGHB

Inj Date : 18-MAY-2000 01:32

Operator: 1891 Inst ID: gc3.i

Smp Info : MHIGHB, 2170-E.b,, 4-INDB. sub,, 1,4

Misc Info : 190-84-10

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

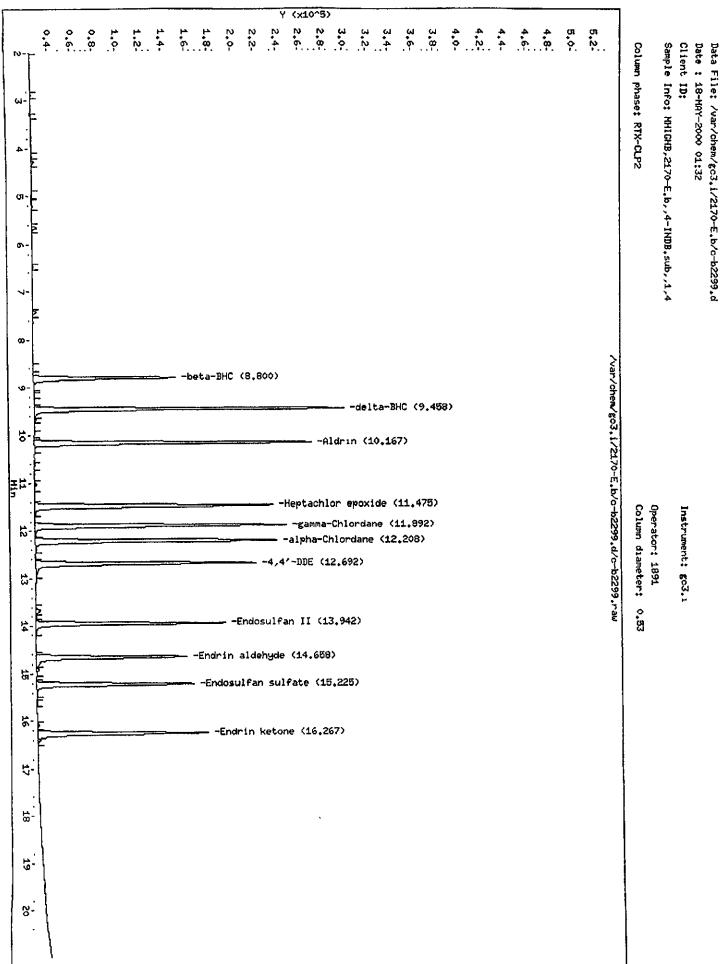
Meth Date: 18-May-2000 11:42 g Quant Type: ESTD Cal Date: 18-MAY-2000 01:32 Cal File: c-b2299.d

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 4-INDB.sub

Target Version: 3.40

					AMOU	NTS
					CAL-AMT	ON-COL
Compounds	RT	BXP RT	DLT RT	RESPONSE	( ng)	( ng)
	<b>48</b>	=====	****	****====	#24555	
11 Aldrin	10.167	10.167	0.000	245294	0.05000	0 0461307
7 beta-BHC	8.800	8.800	0.000	125530	0.05000	0.0439265
8 delta-BHC	9.458	9 458	0.000	274439	0.05000	0.0483326
12 Heptachlor epoxide	11 47S	11.475	0.000	210478	0.05000	0.0462432
13 gamma-Chlordane	11.892	11.892	0.000	222340	0.05000	0 0468983
14 alpha-Chlordane	12 208	12.208	0.000	213437	0.05000	0.0460933
16 4,4'-DDB	12 692	12.692	0.000	194889	0.05000	0.0485023
22 Endosulfan II	13.942	13.942	0.000	166887	0.05000	0 0475794
24 Endrin aldehyde	14.658	14.658	0.000	133001	0.05000	0.0463335
26 Endosulfan sulfate	15.225	15.225	0.000	139369	0.05000	0.0485647
27 Endrin ketone	16.267	16.267	0.000	150750	0.05000	0.0476700



Data File: /var/chem/gc3.i/2170-E.b/c-b2300.d

Report Date: 18-May-2000 11:45

### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2300.d

Lab Smp Id: HIGHB

Inj Date : 18-MAY-2000 01:57

Operator: 1891 Inst ID: gc3.i

Smp Info : HIGHB, 2170-E.b, , 4-INDB. sub, , 1, 5

Misc Info : 190-84-11

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date: 18-May-2000 11:42 g Quant Type: ESTD Cal Date: 18-MAY-2000 01:57 Cal File: c-b2300.d

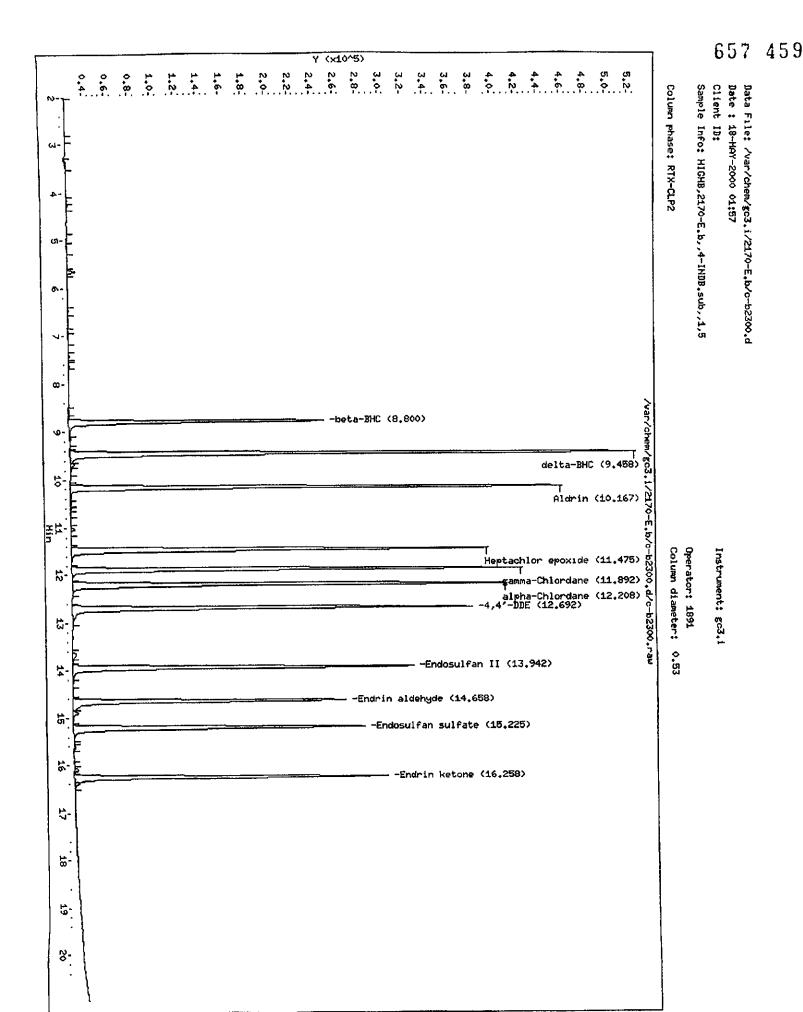
Als bottle: 1 Calibration Sample, Level: 5

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 4-INDB.sub

Target Version: 3.40

					AMOU	NTS
					CAL-AMT	ON-COL
Compounds	RT	BXP RT	DLT RT	response	( ng)	( ng)
	==	*****			P======	****
11 Aldrin	10.167	10.167	0.000	436171	0.10000	0.0850860
7 beta-BHC	8 800	8.800	0.000	225417	0.10000	0.0823586
8 delta-BHC	9.458	9.458	0.000	501642	0.10000	0.0904546
12 Heptachlor epoxide	11.475	11.475	0.000	370895	0.10000	0.0846208
13 gamma-Chlordane	11.892	11.892	0.000	400585	0 10000	0.0871995
14 alpha-Chlordane	12 208	12.208	0.000	386580	0 10000	0.0863365
16 4,4'-DDE	12 692	12.692	0.000	355841	0 10000	0.0906326
22 Endosulfan II	13.942	13.942	0.000	303765	0 10000	0.0889876
24 Endrin aldehyde	14.658	14.658	0 000	242240	0.10000	0 0871087
26 Endosulfan sulfate	15.225	15.225	0.000	258531	0.10000	0.0919100
27 Endrin ketone	16 258	16.267	-0.009	278391	0.10000	0.0901912



Data File: /var/chem/gc3.i/2170-E.b/c-b2301.d

Report Date: 18-May-2000 11:45

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2301.d

Lab Smp Id: 2ND A

Inj Date : 18-MAY-2000 02:23
Operator : 1891 Inst ID: gc3.i

Smp Info : 2ND A,2170-E.b,,INDA.sub,,2,3

Misc Info : 190-82-2

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Quant Type: ESTD Meth Date : 18-May-2000 11:42 g Cal File: c-b2300.d Cal Date : 18-MAY-2000 01:57

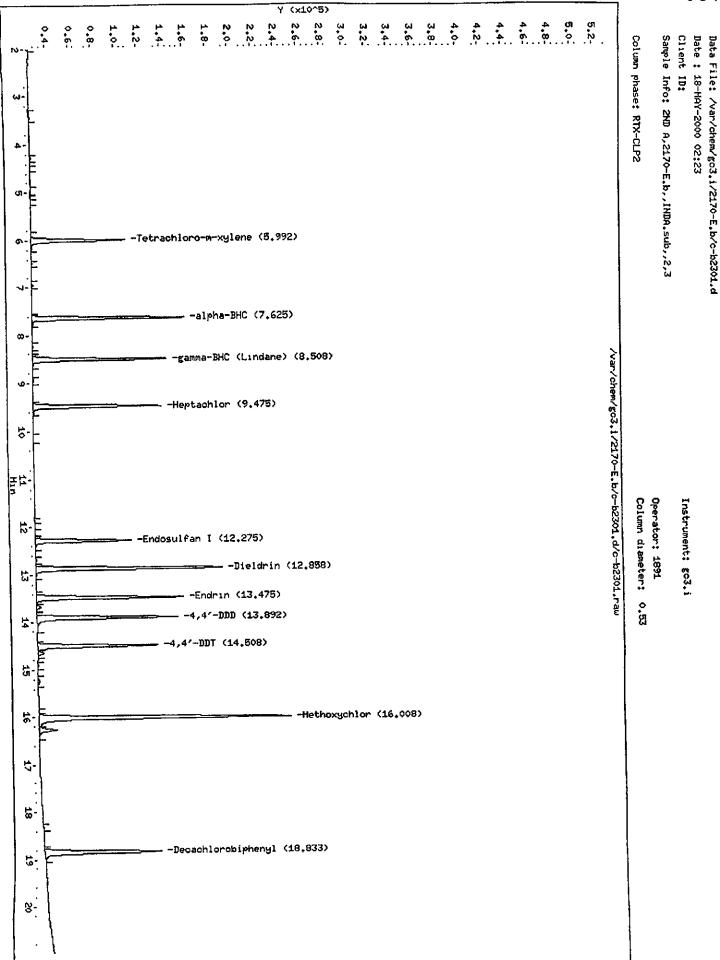
Continuing Calibration Sample Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: INDA.sub Integrator: Falcon

Target Version: 3.40

						MOU	nts
						CAL-AMT	ON-COL
Co	npounds	RT	BXP RT	DLT RT	RESPONSE	( ng)	( ng)
==		고목	EE227=	*****		网络沙尔萨罗斯	****
\$	1 Tetrachloro-m-xylene	5.992	5.992	0.000	83284	0 02500	0.0202659
•	5 alpha-BHC	7.625	7.625	0.000	133894	0 02500	0.0210182
	6 gamma-BHC (Lindane)	8.508	8.508	0.000	117671	0 02500	0.0218481
	10 Heptachlor	9.475	9.475	0.000	112288	0.02500	0.0212255
	15 Endosulfan I	12.275	12.275	0 000	84670	0.02500	0 0209754
	17 Dieldrin	12.658	12.858	0.000	164637	0.02500	0.0396564
	20 Bndrin	13.475	13.475	0.000	129700	0.02500	0 0366390
	21 4,4'-DDD	13.892	13.892	0.000	124532	0.02500	0.0413511
	23 4,4'-DDT	14.508	14.508	0.000	106368	0.02500	0.0405951
	25 Methoxychlor	16.008	16 008	0 000	222949	0.05000	0 178357
\$	30 Decachlorobiphenyl	18.833	18.842	-0.009	103128	0.02500	0.0381246



Data File: /var/chem/gc3.i/2170-E.b/c-b2302.d

Report Date: 18-May-2000 11:45

### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2302.d

Lab Smp Id: 2ND B

Inj Date : 18-MAY-2000 02:49

Operator: 1891 Inst ID: gc3.i

Smp Info : 2ND B,2170-E.b,,INDB.sub,,2,3

Misc Info : 190-82-5

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date: 18-May-2000 11:42 g Quant Type: ESTD Cal Date: 18-MAY-2000 01:57 Cal File: c-b2300.d

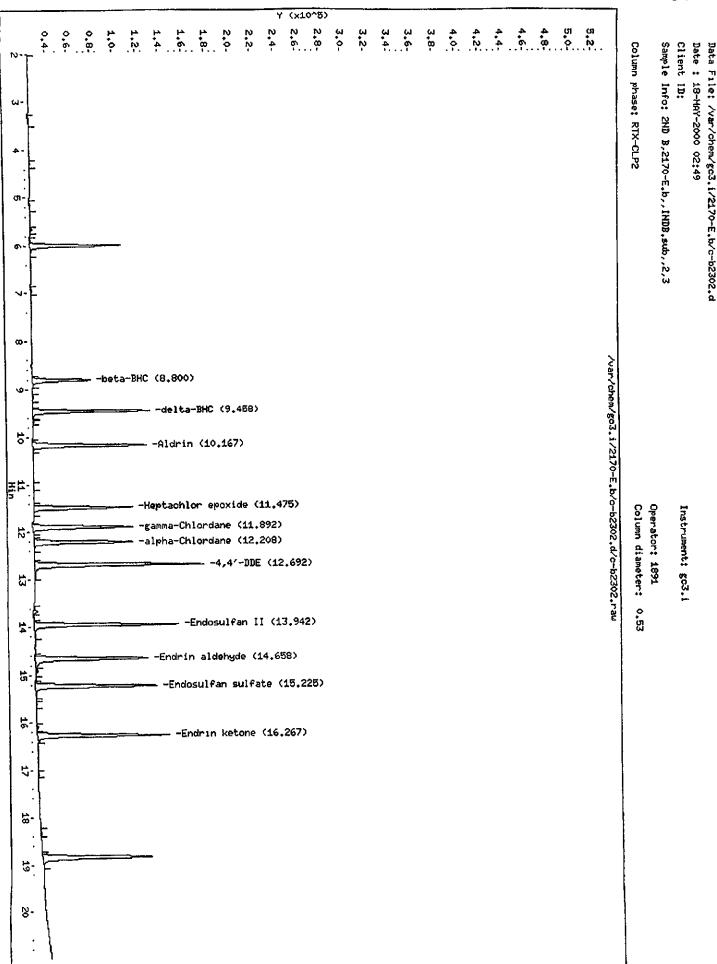
Als bottle: 1 Continuing Calibration Sample

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: INDB.sub

Target Version: 3.40

					AMOU	NTS
					CAL-AMT	ON-COL
Compounds	RT	BXP RT	DLT RT	response	( ng)	( ng)
== #2==================================	**	***	=====	*******	======	======
11 Aldrin	10.167	10 167	0 000	100971	0.02500	0.0196969
7 beta-BHC	8.800	8.800	0.000	51862	0.02500	0.0189484
8 delta-BHC	9.458	9.458	0.000	104007	0.02500	0.0187542
12 Heptachlor epoxide	11.475	11.475	0 000	87730	0.02500	0.0200159
13 gamma-Chlordane	11.892	11.892	0.000	88875	0.02500	0.0193463
14 alpha-Chlordane	12.208	12.208	0 000	87649	0.02500	0.0195750
16 4.4'-DDE	12.692	12.692	0.000	150009	0.02500	0.0382072
22 Endosulfan II	13.942	13.942	0.000	127424	0.02500	0.0373287
24 Endrin aldehyde	14.658	14.658	0.000	99733	0.02500	0.0358637
26 Endosulfan sulfate	15.225	15.225	0.000	107047	0.02500	0.0380561
27 Endrin ketone	16.267	16.267	0 000	116990	0.02500	0.0379016



Data File: /var/chem/gc3.i/2170-E.b/c-b2303.d

Report Date: 18-May-2000 11:45

## STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2303.d

Lab Smp Id: EVALB

Inj Date : 18-MAY-2000 03:15

Inst ID: gc3.i Operator : 1891

Smp Info : EVALB, 2170-E.b, , EVALBR. sub, , 3, 1

Misc Info : 190-88-8

Comment

: /var/chem/gc3.i/2170-E.b/PESTB.m Method

Quant Type: ESTD Meth Date : 18-May-2000 11:42 g Cal File: c-b2300.d Cal Date : 18-MAY-2000 01:57

QC Sample: PEM Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: EVALBR.sub Integrator: Falcon

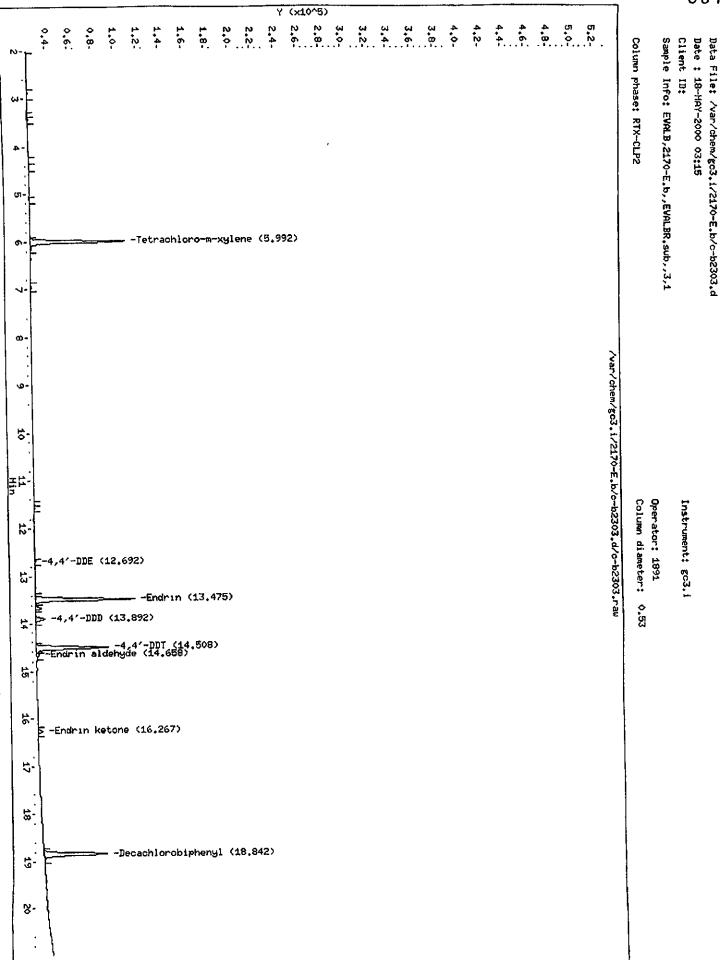
Target Version: 3.40

					CONCENT	rations
					ON-COLUMN	FINAL
Compounds	RT	EXP RT	DLT RT	response	( ng)	( ng)
	==	=====	===##=	==== <b>=</b> =	3#######	====#==
\$ 1 Tetrachloro-m-xylene	5.992	5.992	0 000	84240	0.02050	0.0204985(R)
16 4,4'-DDB	12.692	12.692	0.000	715	0.000182	0.000182110
20 Endrin	13.475	13.475	0.000	88469	0 02499	0.0249917
21 4,4'-DDD	13.892	13.892	0 000	8171	0.00271	0 00271320
23 4,4'-DDT	14.508	14.508	0.000	64749	0.02471	0.0247113
24 Endrin aldehyde	14.658	14 658	0.000	2023	0.000727	0.000727464
27 Endrin ketone	16.267	16.267	0 000	4449	0 00144	0.00144136
\$ 30 Decachlorobiphenyl	18.842	18.842	0.000	56206	0.02078	0.0207784(R)

# OC Flag Legend

R - Spike/Surrogate failed recovery limits.

00× 12.100 EXDRIN 6.500



Data File: /var/chem/gc3.i/2170-E.b/c-b2361.d

Report Date: 19-May-2000 11:25

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2361.d

Lab Smp Id: EVALB

Inj Date : 19-MAY-2000 08:32

Operator : 1891 Inst ID: gc3.i

Smp Info : EVALB, 2170-E.b, , EVALBR. sub, , 3, 1

Misc Info : 190-88-8

Comment

: /var/chem/gc3.i/2170-E.b/PESTB.m Method

Meth Date: 19-May-2000 11:24 eppinged Quant Type: ESTD Cal File: c-b2300.d Cal Date : 18-MAY-2000 01:57

QC Sample: PEM Als bottle: 1

Dil Factor: 1.00000 Compound Sublist: EVALBR.sub Integrator: Falcon

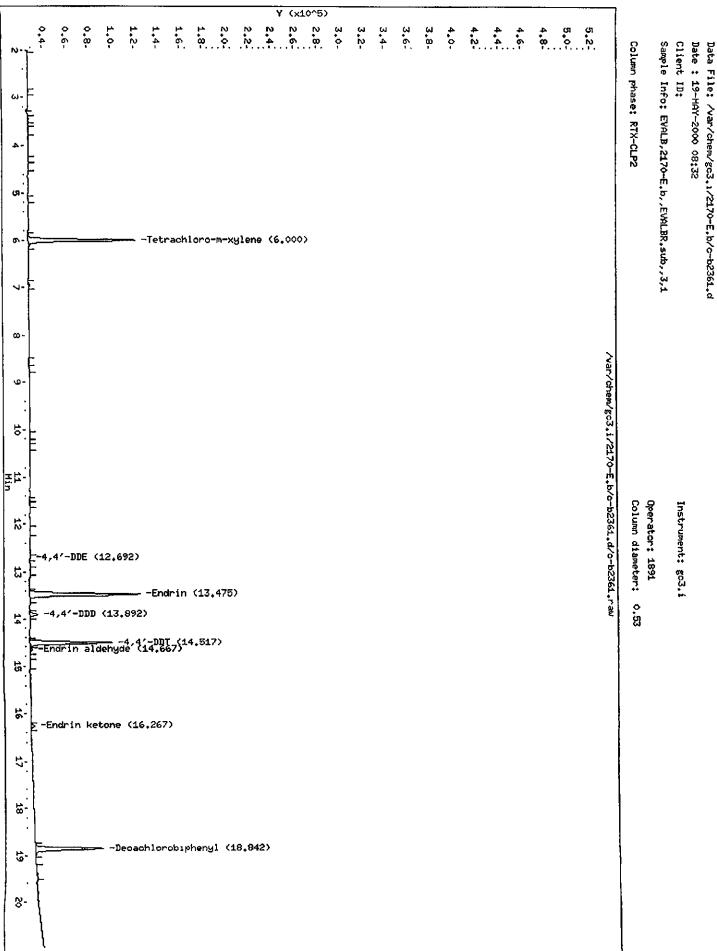
Target Version: 3.40

						CONCENTRATIONS		
						ON-COLUM	N PINAL	
Comp	ounds	R <b>T</b>	BXP RT	DLT RT	response	( ng)	( ng)	
		==	======	444005	855555		****	
\$	1 Tetrachloro-m-xylene	6.000	6.000	0.000	93867	0.02284	0.0228411(R)	
1	6 4,4'-DDB	12.692	12 692	0.000	769	0.000196	0.000195864	
2	0 Endrin	13.475	13.475	0.000	96844	0.02736	0.0273575	
2	1 4,4'-DDD	13.892	13.900	-0.008	7577	0.00252	0.00251596	
2	3 4,4'-DDT	14.517	14.517	0 000	72454	0.02765	0 0276519	
2	4 Endrin aldehyde	14.667	14.667	0.000	1660	0.000597	0.000596931	
2	7 Endrin ketone	16.267	16.267	0 000	3630	0.00118	0.00117602	
\$ 3	O Decachlorobiphenyl	18.842	18.842	0 000	59349	0.02194	0.0219403(R)	

## OC Flag Legend

R - Spike/Surrogate failed recovery limits

DOT 10:39.



Data File: /var/chem/gc3.i/2170-E.b/c-b2362.d

Report Date: 19-May-2000 11:23

### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2362.d

Lab Smp Id: MEDTOX

Inj Date : 19-MAY-2000 08:58

Operator: 1891 Inst ID: gc3.i

Smp Info : MEDTOX, 2170-E.b, ,1-TOX.sub, ,2,3

Misc Info : 190-84-13

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date: 19-May-2000 11:22 g Quant Type: ESTD Cal Date: 18-MAY-2000 01:57 Cal File: c-b2300.d

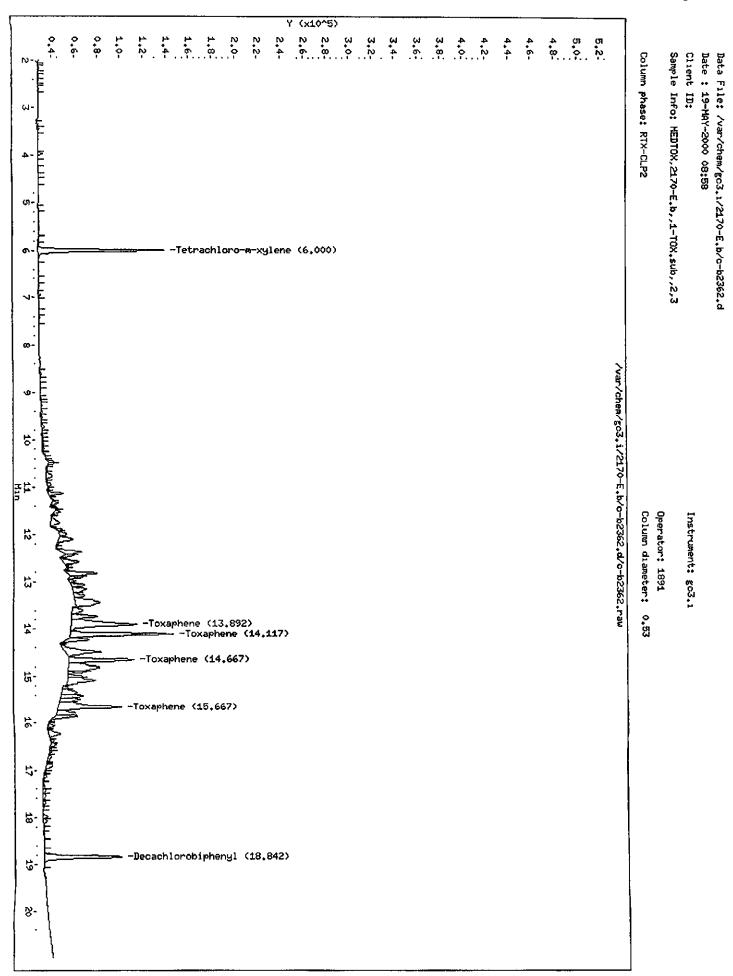
Als bottle: 1 Continuing Calibration Sample

Dil Factor: 1.00000 Integrator: Falcon

Integrator: Falcon Compound Sublist: 1-TOX.sub

Target Version: 3.40

					amounts		
					CAL+AMT	ON-COL	
Ça	mpounds	RT	EXP RT DLT RT	response	( ng)	( ng)	
==		==	****		=======		
	18 Toxaphene	13 892	13.892 0.000	5757 <del>9</del>	1.00000	1.08796	
\$	1 Tetrachloro-m-xylene	6 000	6 000 0,000	111015	0.02500	0.0270138	
\$	30 Decachlorobiphenyl	18.842	18.842 0.000	68483	0.02500	0.0253169	



Data File: /var/chem/gc3.i/2170-E.b/c-b2363.d

Report Date: 19-May-2000 11:23

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-b2363.d

Lab Smp Id: MEDCHLOR

Inj Date : 19-MAY-2000 09:24

Operator : 1891 Inst ID: gc3.i

Smp Info : MEDCHLOR, 2170-E.b,,2-CHLO.sub,,2,3 Misc Info : 190-85-10

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTB.m Meth Date : 19-May-2000 11:22 g Quar Cal Date : 18-MAY-2000 01:57 Cal Quant Type: ESTD Cal File: c-b2300.d

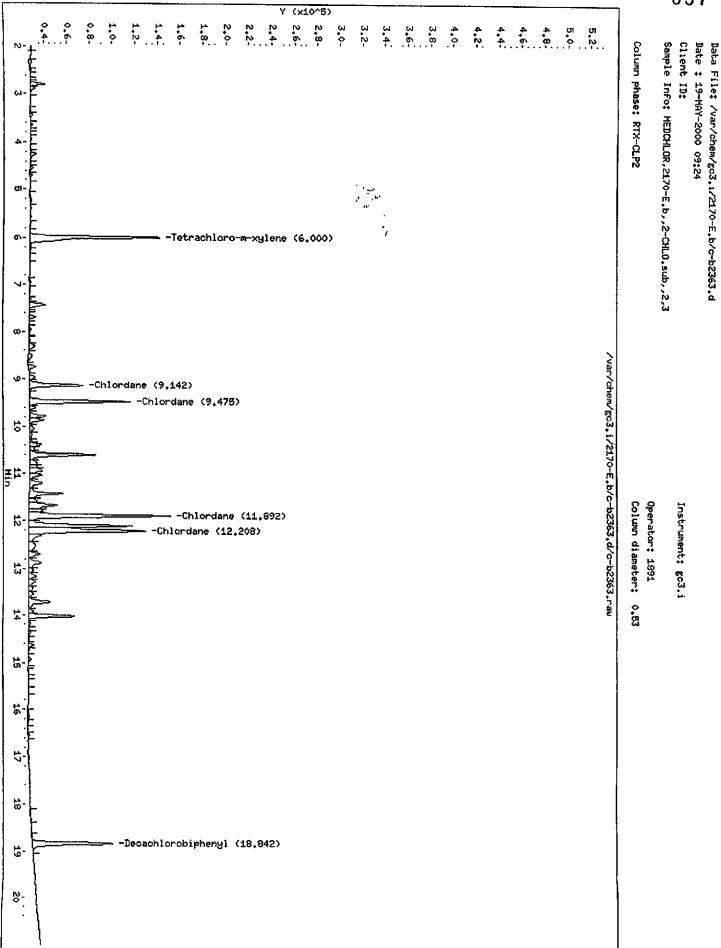
Als bottle: 1 Continuing Calibration Sample

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: 2-CHLO.sub

Target Version: 3.40

				DOMA	nts
				CAL-AMT	ON-COL
Compounds	RT	BXP RT DLT RT	RESPONSE	( ng)	( ng)
		22222 958888			======
9 Chlordane	9.142	9 142 0.000	47163	0 25000	0.300050
\$ 1 Tetrachloro-m-xylene	6 000	6 000 0.000	114138	0 02500	0 0277737
\$ 30 Decachlorobiphenyl	18 842	18 842 0.000	71224	0.02500	0 0263302



Data File: /var/chem/gc3.i/2170-E.b/c-b2364.d

Report Date: 19-May-2000 11:24

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2364.d

Lab Smp Id: MEDA

Inj Date : 19-MAY-2000 09:50

Operator : 1891 Smp Info : MEDA,2170-E.b,,INDA.sub,,2,3 Inst ID: gc3.i

Misc Info: 190-84-3

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTB.m Meth Date : 19-May-2000 11:22 g Quar Quant Type: ESTD Cal File: c-b2300.d Cal Date : 18-MAY-2000 01:57

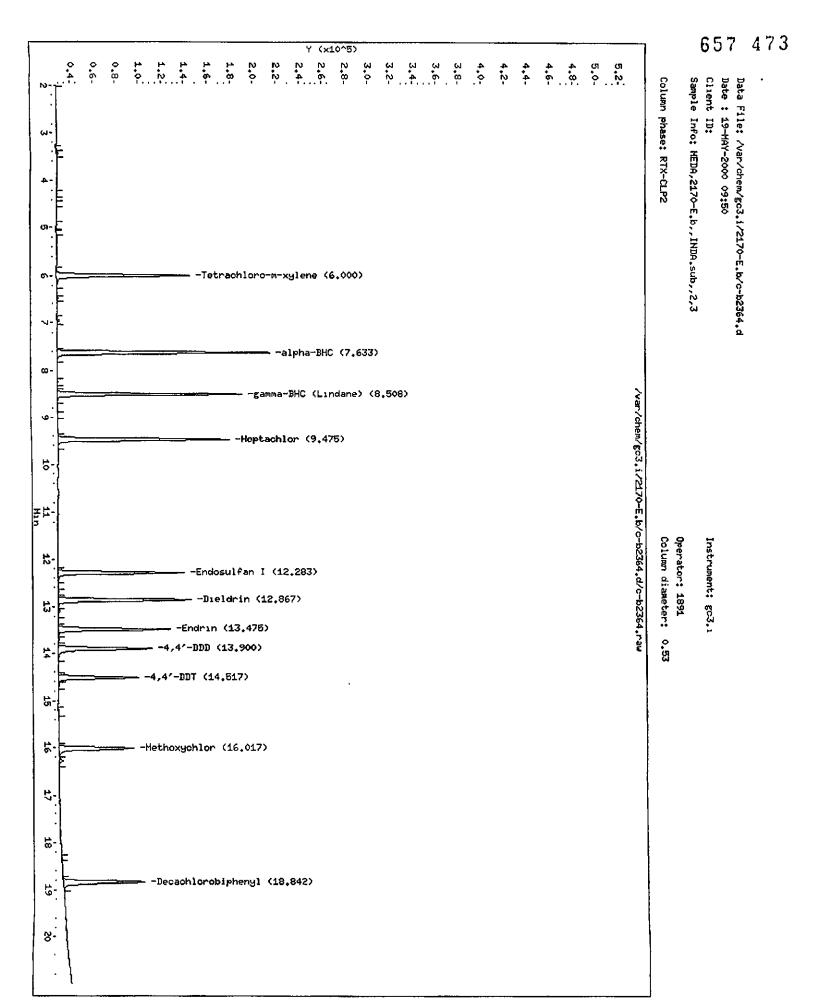
Continuing Calibration Sample Als bottle: 1

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: INDA.sub

Target Version: 3.40

					AMOU	nts
					CAL-AMT	ON-COL
Compounds	RT	BXP RT	DLT RT	response	(ng)	( ng)
	==	======			======	
\$ 1 Tetrachloro-m-xylene	6.000	6,000	0.000	117061	0 02500	0.0284850
5 alpha-BHC	7.633	7.633	0.000	185823	0.02500	0.0291698
6 gamma-BHC (Lindane)	8.508	8.508	0.000	161288	0.02500	0.0299465
10 Heptachlor	9.475	9.475	0.000	150514	0.02500	0.0284512
15 Endosulfan I	12.283	12 283	0.000	111488	0.02500	0.0276191
17 Dieldrin	12.867	12.867	0.000	117253	0.02500	0.0282429
20 Endrin	13.475	13.475	0.000	99123	0.02500	0.0280013
21 4,4'-DDD	13.900	13.900	0.000	62852	0.02500	0.0275112
23 4,4'-DDT	14.517	14.517	0.000	71033	0.02500	0.0271096
25 Methoxychlor	16.017	16.017	0.000	66694	0.05000	0.0533546
\$ 30 Decachlorobiphenyl	18.842	18.842	0.000	71717	0.02500	0.0265125



Data File: /var/chem/gc3.i/2170-E.b/c-b2365.d

Report Date: 19-May-2000 11:24

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2365.d

Lab Smp Id: MEDB

Inj Date : 19-MAY-2000 10:16

Operator: 1891 Inst ID: gc3.i

Smp Info : MEDB, 2170-E.b, , INDB. sub, , 2, 3

Misc Info : 190-84-9

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

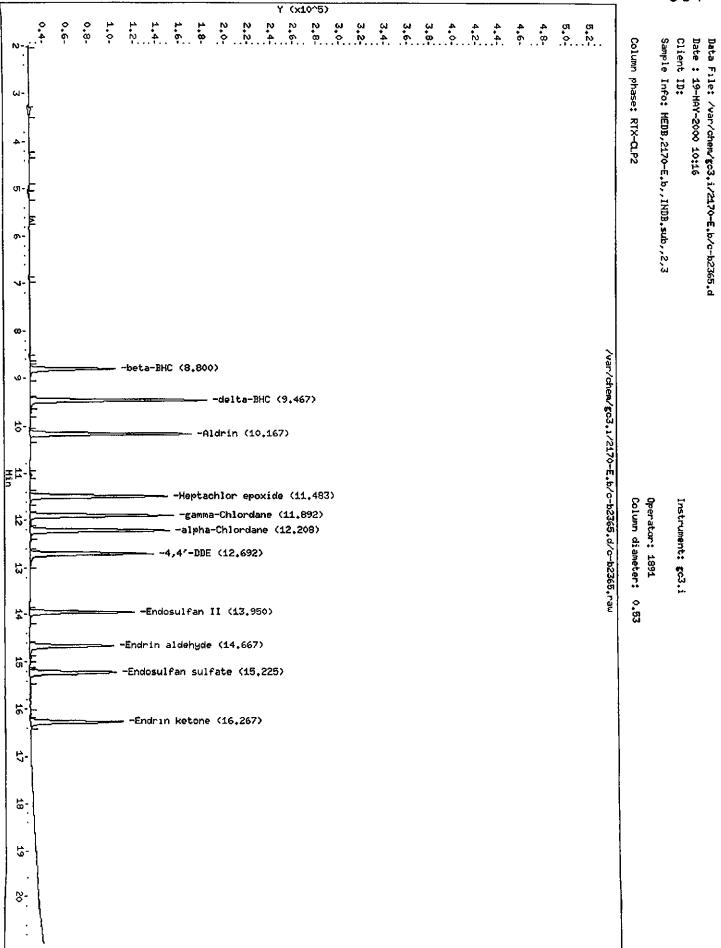
Meth Date: 19-May-2000 11:22 g Quant Type: ESTD Cal Date: 18-MAY-2000 01:57 Cal File: c-b2300.d

Als bottle: 1 Continuing Calibration Sample

Dil Factor: 1.00000 Integrator: Falcon Compound Sublist: INDB.sub

Target Version: 3.40

					AMOU	nts
					CAL-AMT	ON-COL
Compounds	RT	EXP RT	DLT RT	response	( ng)	( ng)
医亚巴亚古亚亚亚 经营票票票 医医毒素 医甲基甲甲甲二二烷	==		200000	*****		*****
11 Aldrin	10.167	10.167	0.000	142483	0 02500	0 0277949
7 beta-BHC	8.800	8.800	0.000	75023	0.02500	0 0274105
8 delta-BHC	9.467	9.467	0.000	155556	0.02500	0.0280494
12 Heptachlor epoxide	11.483	11.483	0.000	121281	0.02500	0.0276706
13 gamma-Chlordane	11.892	11.892	0.000	126693	0.02500	0.0275786
14 alpha-Chlordane	12.208	12.208	0.000	123299	0.02500	0.0275369
16 4,4'-DDE	12.692	12.692	0.000	108503	0.02500	0.0276357
22 Endosulfan II	13.950	13.950	0 000	92337	0.02500	0.0270500
24 Endrin aldehyde	14.667	14.667	0.000	74472	0.02500	0.0267799
26 Endosulfan sulfate	15.225	15.225	0.000	76466	0.02500	0.0271843
27 Endrin ketone	16.267	16.267	0.000	82040	0.02500	0.0265788



Data File: /var/chem/gc3.i/2170-E.b/c-b2374.d

Report Date: 19-May-2000 14:33

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2374.d

Lab Smp Id: MEDA

Inj Date : 19-MAY-2000 14:10

Operator: 1891 Inst ID: gc3.i

Smp Info : MEDA, 2170-E.b, , INDA. sub, , 2, 3

Misc Info: 190-84-3

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

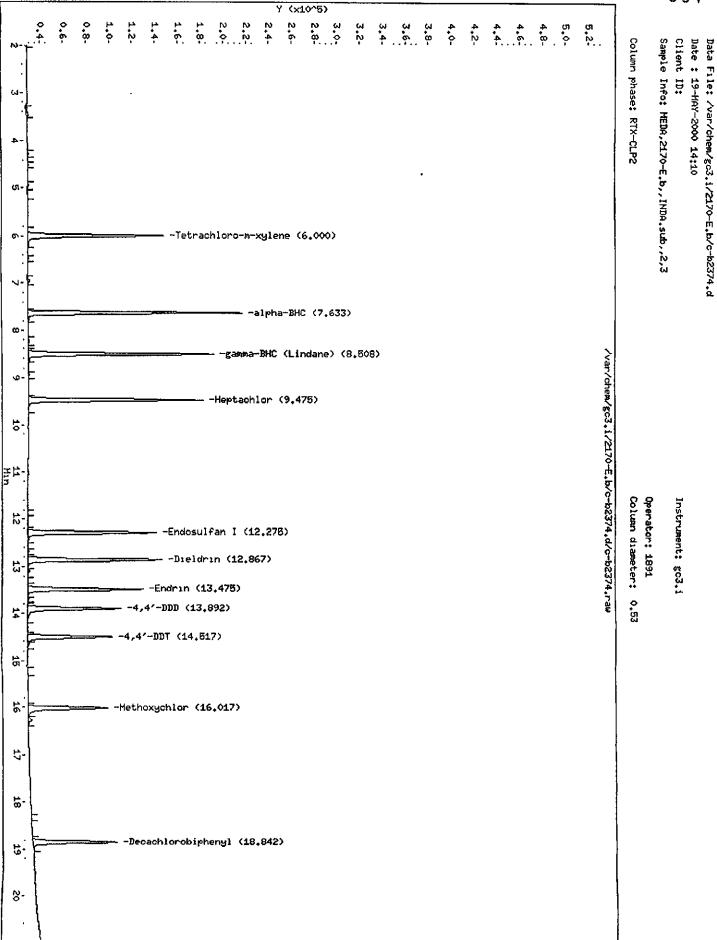
Meth Date: 19-May-2000 14:32 g Quant Type: ESTD Cal Date: 18-MAY-2000 01:57 Cal File: c-b2300.d

Als bottle: 1 Continuing Calibration Sample

Dil Factor: 1.00000
Integrator: Falcon
Compound Sublist: INDA.sub

Integrator: Falcon
Target Version: 3.40

					AMOU	NTS
					CAL-AMT	ON-COL
Compounds	RT	RXP RT	DLT RT	RESPONSE	( ng)	( ng)
· · · · · · · · · · · · · · · · · · ·	==	=====	======	******	****	******
\$ 1 Tetrachloro-m-xylene	6.000	6.000	0.000	119470	0 02500	0.0290712
5 alpha-BHC	7.633	7.633	0.000	188445	0 02500	0.0295814
6 gamma-BHC (Lindane)	8.508	8.508	0.000	162393	0.02500	0.0301516
10 Heptachlor	9.475	9.475	0.000	153991	0.02500	0.0291085
15 Endosulfan I	12.275	12,283	-0.008	113008	0.02500	0 0279956
17 Dieldrin	12.867	12.867	0.000	118409	0.02500	0 0285214
20 Endrin	13.475	13.475	0 000	101821	0.02500	0 0287635
21 4,4'-DDD	13 892	13.900	-0 008	82452	0 02500	0.0273784
23 4,4'-DDT	14.517	14.517	0.000	74192	0 02500	0.0283152
25 Methoxychlor	16.017	16.017	0.000	70066	0.05000	0.0560522
\$ 30 Decachlorobiphenyl	18.842	18.842	0 000	73722	0 02500	0.0272537



Data File: /var/chem/gc3.i/2170-E.b/c-b2375.d

Report Date: 19-May-2000 15:21

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-b2375.d

Lab Smp Id: MEDB Inj Date : 19-MAY-2000 14:36

Inst ID: gc3.i Operator : 1891

Smp Info : MEDB, 2170-E.b,, INDB. sub,, 2,3

Misc Info: 190-84-9

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Quant Type: ESTD Meth Date : 19-May-2000 15:21 g Cal File: c-b2300.d Cal Date : 18-MAY-2000 01:57

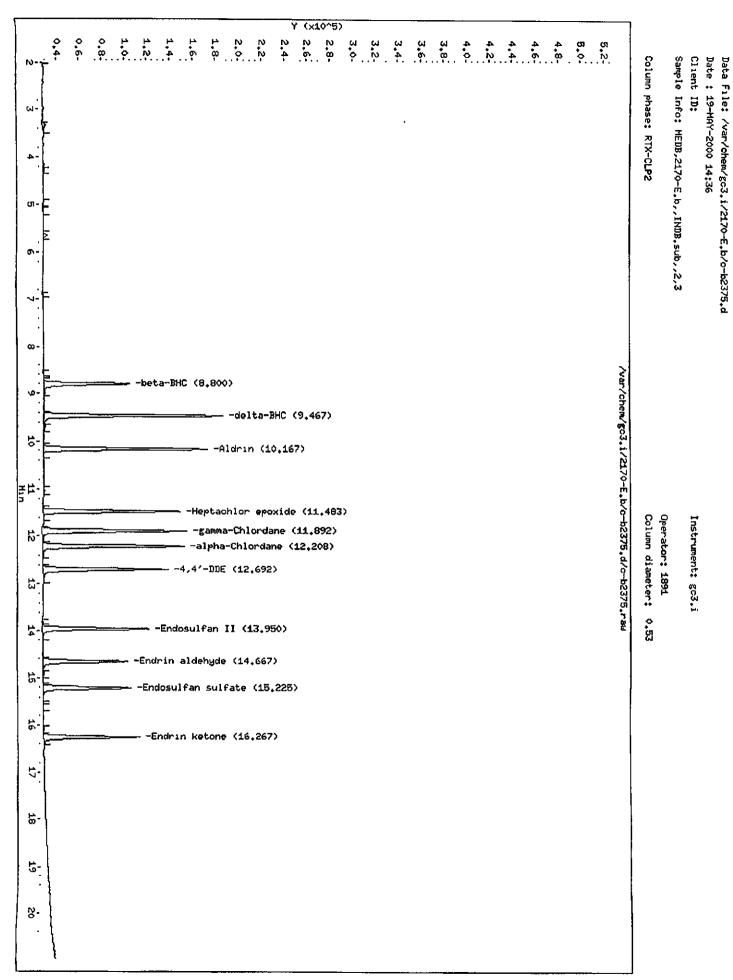
Continuing Calibration Sample Als bottle: 1

Dil Factor: 1.00000

Compound Sublist: INDB.sub Integrator: Falcon

Target Version: 3.40

					AMOU	NTS
					CAL-AMT	ON-COL
Compounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
	==		89555	****		======
11 Aldrin	10.167	10.167	0.000	144317	0.02500	0 0281526
7 beta-BHC	8.800	8.800	0.000	76218	0.02500	0 0278471
8 delta-BHC	9.467	9.467	0.000	157697	0.02500	0.0284354
12 Heptachlor epoxide	11 483	11.483	0.000	121252	0.02500	0.0276640
13 gamma-Chlordane	11 892	11 892	0.000	127769	0.02500	0.0278128
14 alpha-Chlordane	12.208	12,208	0.000	125020	0.02500	0 0279212
16 4,4'-DDB	12 692	12 692	0.000	110114	0 02500	0.0280460
22 Endosulfan II	13.950	13.950	0.000	94167	0.02500	0.0275861
24 Endrin aldehyde	14.667	14.667	0.000	75748	0 02500	0.0272387
26 Endosulfan sulfate	15.225	15.225	0.000	78306	0.02500	0.0278385
27 Endrin ketone	16.267	16.267	0.000	85148	0.02500	0.0275857



# PESTICIDE QC DATA

## UXB INTERNATIONAL METHOD BLANK COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Lab Sample ID: COE180000 460

Matrix: (soil/water) SOLID-WATER

Method: SW846 8081A

m5/23/00

Pesticides (8081A)

Sample WT/Vol: 100 / mL Work Order: DDE2E101 Dilution factor. 1

Date Received: 05/13/00 Date Extracted: 05/18/00 Date Analyzed: 05/19/00

QC Batch: 0139460

Client Sample Id: INTRA-LAB BLANK

### CONCENTRATION UNITS:

CAS NO.	COMPOUND (ug/L o	r ug/kg) mg/L	Q
57-74-9	Chlordane (technical)	0.0050	ן ט
72-20-8	Endrin	0.00050	וט
76-44-8	Heptachlor	0.00050	ן ט
1024-57-3	Heptachlor epoxide	0.00050	וט
58-89-9	Lindane	0.00050	[ <u>U</u> ]
72-43-5	Methoxychlor	0.0010	וט
8001-35-2	Toxaphene	0.020	ע

Data File: /var/chem/gc3.i/2170-E.b/c-a2372.d

Report Date: 19-May-2000 13:56

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2372.d Lab Smp Id: DDE2E101 Client Client Smp ID: PBLK

Inj Date : 19-MAY-2000 13:18

Operator : 1891 Inst ID: gc3.i

Smp Info : DDE2E101,2170-E.b,,PEST.sub,,3,

Misc Info: 130197BLK Comment: 40851960

: /var/chem/gc3.i/2170-E.b/PESTA.m Method

Meth Date: 19-May-2000 11:01 eppinged Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-a2300.d Als bottle: 1 QC Sample: BLANK

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF	1 000	Dilution Factor
۷t		Volume of final extract (uL)
Vo	100.000	Volume of sample extracted (mL)
Vi		Volume injected

CONCENTRATIONS

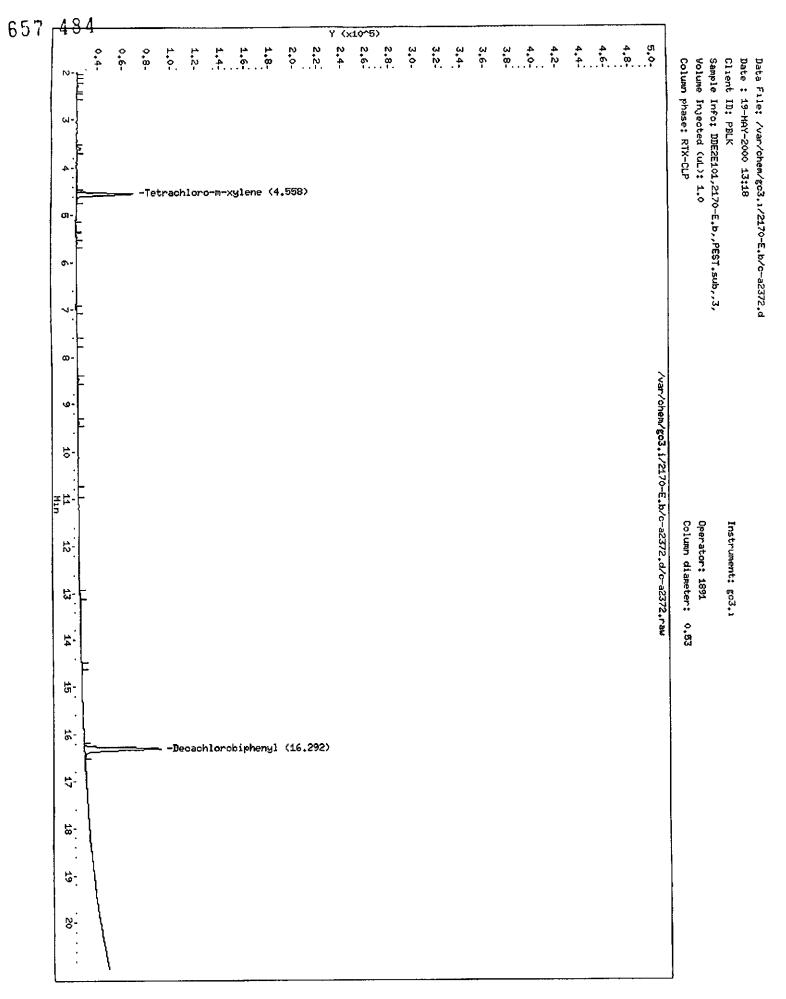
					cc	NCENTRA	TIONS
					ON-C	COLUMN	FINAL
unds	RT	EXP RT DLT	RT	response	(	ng)	( mg/L)
	==			******	===		*****
alpha-BHC	Com	pound Not D	etected				
HEXACHLOROBENZENE	Com	pound Not D	etected	•			
gamma-BHC (Lindane)	Com	pound Not D	etected				
beta-BHC	Com	pound Not E	etected				
Chlordane	Com	pound Not D	etected				
Heptachlor	Com	pound Not D	etected	•			
delta-BHC	Com	pound Not D	etected	•			
Aldrin	Com	pound Not D	etected	•			
Heptachlor epoxide	Com	pound Not D	etected				
gamma-Chlordane	Com	pound Not I	etected				
alpha-Chlordane	Com	pound Not E	etected	•			
Endosulfan I	Com	pound Not E	etect <b>e</b> d	•			
4,4'-DDE	Com	pound Not I	etected	•			
Dieldrin	Com	pound Not E	etected				
Endrin	Com	pound Not E	etected				
Toxaphene	Com	pound Not I	etected	•			
	alpha-BHC HEXACHLOROBENZENE gamma-BHC (Lindane) beta-BHC Chlordane Heptachlor delta-BHC Aldrin Heptachlor epoxide gamma-Chlordane alpha-Chlordane Endosulfan I 4,4'-DDE Dieldrin Endrin Toxaphene	alpha-BHC Com  HEXACHLOROBENZENE Com  gamma-BHC (Lindane) Com  beta-BHC Com  Chlordane Com  Heptachlor Com  delta-BHC Com  Heptachlor com  Heptachlor epoxide Com  gamma-Chlordane Com  alpha-Chlordane Com  endosulfan I Com  1,4'-DDE Com  Dieldrin Com  Endrin Com  Endrin Com	alpha-BHC Compound Not I Compound Not I Gamma-BHC (Lindane) Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compound Not I Compou	alpha-BHC Compound Not Detected HEXACHLOROBENZENE Compound Not Detected gamma-BHC (Lindane) Compound Not Detected beta-BHC Compound Not Detected Chlordane Compound Not Detected Heptachlor Compound Not Detected delta-BHC Compound Not Detected Aldrin Compound Not Detected gamma-Chlordane Compound Not Detected gamma-Chlordane Compound Not Detected alpha-Chlordane Compound Not Detected Endosulfan I Compound Not Detected Lindane Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Compound Not Detected Detection Detected Detection Detected Detection Detected Detection Detected Detection Detected Detection Detected Detection Detected Detection Detected Detection Detected Detection Detected Detection Detected Detection Detected Detection Detected Detection Detection Detection Detected Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Detection Dete	alpha-BHC Compound Not Detected.  HEXACHLOROBENZENE Compound Not Detected.  gamma-BHC (Lindane) Compound Not Detected.  beta-BHC Compound Not Detected.  Chlordane Compound Not Detected.  Heptachlor Compound Not Detected.  Aldrin Compound Not Detected.  Aldrin Compound Not Detected.  Heptachlor epoxide Compound Not Detected.  Heptachlor dane Compound Not Detected.  Gamma-Chlordane Compound Not Detected.  Endosulfan I Compound Not Detected.  Letter Compound Not Detected.  Letter Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.	alpha-BHC Compound Not Detected.  Chlordane Compound Not Detected.  Chlordane Compound Not Detected.  Chlordane Compound Not Detected.  Chlordane Compound Not Detected.  Aldrin Compound Not Detected.  Aldrin Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.  Compound Not Detected.	alpha-BHC Compound Not Detected.  HEXACHLOROBENZENE Compound Not Detected.  gamma-BHC (Lindane) Compound Not Detected.  beta-BHC Compound Not Detected.  Chlordane Compound Not Detected.  Heptachlor Compound Not Detected.  Aldrin Compound Not Detected.  Heptachlor epoxide Compound Not Detected.  Heptachlor epoxide Compound Not Detected.  gamma-Chlordane Compound Not Detected.  alpha-Chlordane Compound Not Detected.  Endosulfan I Compound Not Detected.  4,4'-DDE Compound Not Detected.  Endrin Compound Not Detected.  Endrin Compound Not Detected.

Data File: /var/chem/gc3.i/2170-E.b/c-a2372.d Report Date: 19-May-2000 13:56

					CONCENTR	ATIONS
					ON-COLUMN	FINAL
Compo	punds	RT	EXP RT DLT RT	RESPONSE	( ng)	( mg/L)
***	****************			*****		
21	4,4'-DDD	Co	mpound Not Detect	ed.		
22	Endosulfan II	Co	mpound Not Detect	eđ		
23	4,4'-DDT	Co	mpound Not Detect	ed		
24	Endrin aldehyde	Co	mpound Not Detect	ed.		
26	Endosulfan sulfate	Co	mpound Not Detect	ed.		
25	Methoxychlor	Co	mpound Not Detect	ed.		
27	Endrin ketone	Co	mpound Not Detect	ed.		
\$ 1	Tetrachloro-m-xylene	4.558	4.558 0.000	46643	0 01152	0.00115196 (AR)
\$ 30	Decachlorobiphenyl	16.292	16.300 -0 008	62757	0 01905	0.00190500 (aR)

## QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).
   R Spike/Surrogate failed recovery limits.



Data File: /var/chem/gc3.i/2170-E.b/c-b2372.d

Report Date: 19-May-2000 13:58

# STL-PITTSBURGH ,

Data file: /var/chem/gc3.i/2170-E.b/c-b2372.d

Lab Smp Id: DDE2E101 Client Smp ID: PBLK

Inj Date : 19-MAY-2000 13:18

Operator: 1891 In Smp Info: DDE2E101,2170-E.b,,PEST.sub,,3, Inst ID: gc3.i

Misc Info: 130194BLK

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date: 19-May-2000 11:24 eppinged Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-b2300.d Als bottle: 1 QC Sample: BLANK

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF Vt Vo Vi	10.000 100.000	Dilution Factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected

C	ONCENTRATIONS	

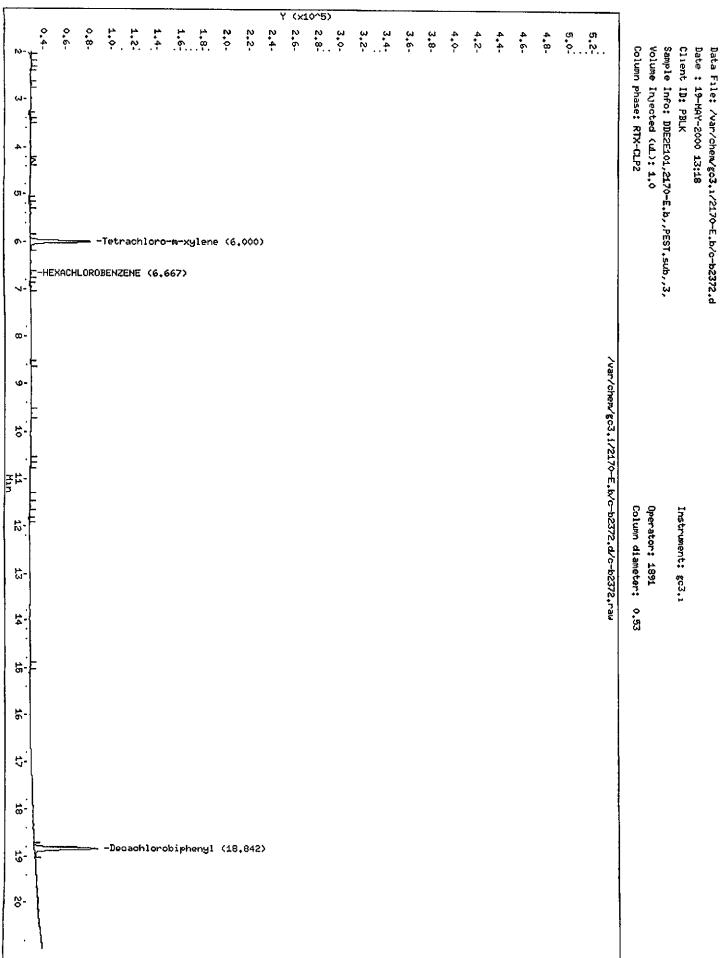
							071-0	COLUMN	PINAL
Compo	undo	RT	BXP RT	DLI	RT	response	(	ng)	(mg/L)
		-	905888	<b>t</b> res	ann	enenana	#5		
5	alpha-BHC	Com	pound N	ot I	otecte	d.			
4	HEXACHLOROBENZENE	6.667	6.650	0	017	458			(a)
6	gamma-BHC (Lindane)	Com	pound N	ot [	etecta	đ.			
7	beta-BHC	Com	pound N	ot I	etecte	đ.			
9	Chlordane	Com	pound N	ot [	etecte	d			
10	Heptachlor	Com	pound N	ot [	otecte	d			
8	delta-BHC	Com	pound N	ot I	etecte	đ.			
11	Aldrin	Com	pound N	ot I	etecte	d.			
12	Heptachlor epoxide	Com	pound N	ot [	etecte	d.			
13	gamma-Chlordane	Com	pound N	ot [	Datacta	d.			
14	alpha-Chlordane	Com	pound N	ot I	)etecte	đ.			
15	Endosulfan I	Соп	pound N	ot I	etecte	d.			
16	4,4'-DDB	Con	pound N	ot I	etecte	d.			
17	Dieldrin	Con	apound N	ot I	Detecte:	d.			
20	Bndrin	Con	pound N	ot I	etecte	d.			
18	Toxaphene	Com	pound N	ot I	etecte	d			

Data File: /var/chem/gc3.i/2170-E.b/c-b2372.d Report Date: 19-May-2000 13:58

		Concentrations
		ON-COLUMN FINAL
Compounds	RT EXP RT DLT RT RESPONS	BB ( ng) ( mg/L)
2617555555555555555		
21 4,4'-DDD	Compound Not Detected	
22 Endosulfan II	Compound Not Detected.	
23 4,4'-DDT	Compound Not Detected.	
24 Endrin aldehyde	Compound Not Detected.	
26 Endosulfan sulfate	Compound Not Detected.	
25 Methoxychlor	Compound Not Detected	
27 Endrin ketone	Compound Not Detected.	
\$ 1 Tetrachloro-m-xylene	6,000 6.000 0.000 5355	18 0.01382 0.00130228(aR)
\$ 30 Decachlorobiphenyl	18 842 18.842 0.000 5562	27 0.02086 0.00205643 (aR)

# QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).R - Spike/Surrogate failed recovery limits.



## UXB INTERNATIONAL CHECK SAMPLE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (SOIL/Water) SOLID WARK Lab Sample ID:COE180000 460 Method: SW846 8081A WS)33 50

Pesticides (8081A)

Date Received: 05/13/00 Sample WT/Vol: 100 / mL Work Order: DDE2E102 Date Extracted: 05/18/00 Date Analyzed: 05/19/00 Dilution factor: 1

QC Batch: 0139460

Client Sample Id: CHECK SAMPLE

## CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L Q
72-20-8	Endrin	0.00219
76-44-8	Heptachlor	0.00212
1024-57-3	Heptachlor epoxide	0.00219
58-89-9	Lindane	0.00189
72-43-5	Methoxychlor	0.00277

Data File: /var/chem/gc3.i/2170-E.b/c-a2373.d

Report Date: 19-May-2000 14:13

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2373.d

Lab Smp Id: DDE2E102 Client Smp ID: LCS

Inj Date : 19-MAY-2000 13:44

Operator : 1891 Inst ID: gc3.i

Smp Info : DDE2E102,2170-E.b,,PEST.sub,,3,
Misc Info : 130192LCS

Comment

: /var/chem/gc3.i/2170-E.b/PESTA.m Method

Meth Date: 19-May-2000 11:01 eppinged Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-a2300.d

Als bottle: 1 QC Sample: LCS

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF		Dilution Factor
Vt	10.000	Volume of final extract (uL)
Vo	100.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected

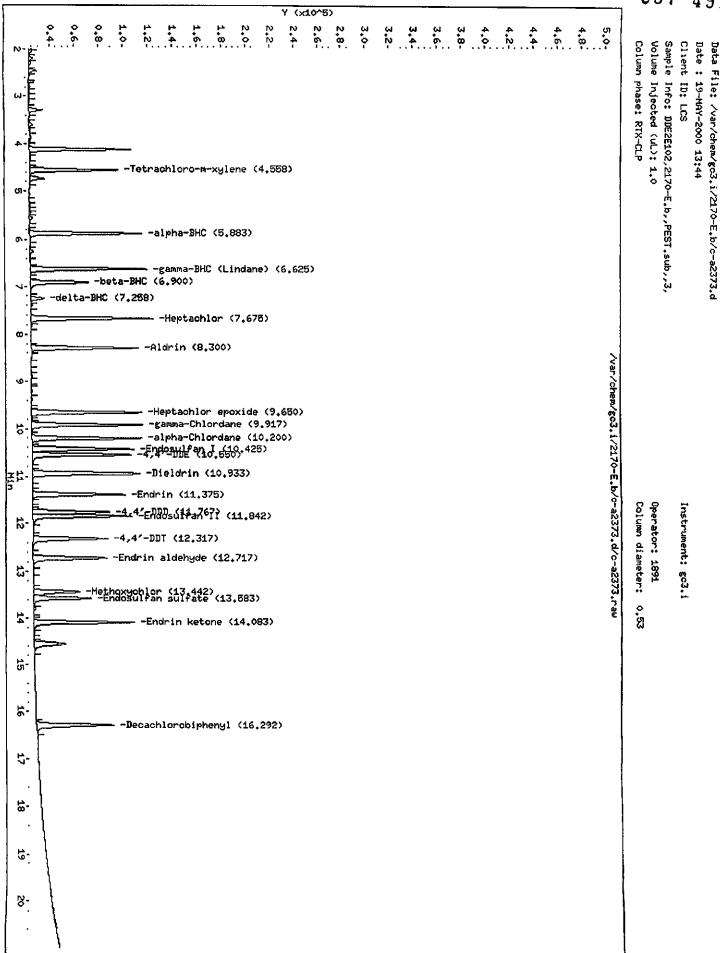
				CONCENTRA	tions
				ON-COLUMN	FINAL
Compounds	RT	EXP RT DLT RT	RESPONSE	( ng)	( mg/L)
CCDS # 부부부경제 COD SEE SEE SEE SE	85	2xcecs 23ee99		*****	=======
5 alpha-BHC	5 883	5.883 0 000	92466	0.01662 0	00166213(a)
4 HEXACHLOROBENZENE	Comp	pound Not Detect	eđ.		
6 gamma-BHC (Lindane)	6.625	6.625 0.000	96274	0.01887	00188671 (aR)
7 beta-BHC	6.900	6.908 -0.008	48715	0.02011 0	00201078(a)
9 Chlordane	Comp	pound Not Detect	ed.		
10 Heptachlor	7.675	7 675 0.000	101376	0.02115 0	00211506 (aR)
8 delta-BHC	7 258	7 258 0.000	11524	0.00245 0	000245204 (a)
11 Aldrin	8.300	8.300 0.000	88668	0.01989 0	00198905 (aR)
12 Heptachlor epoxide	9.650	9.650 0.000	91362	0 02189	.00218908(a)
13 gamma-Chlordane	9.917	9.925 -0.008	92010	0.02122 0	.00212158(a)
14 alpha-Chlordane	10.200	10.200 0.000	91125	0.02124 0	.00212446(a)
15 Endosulfan I	10.425	10.425 0.000	84658	0.02194 0	.00219413(a)
16 4,4'-DDE	10.550	10.550 0.000	81621	0.02241 0	00224116(a)
17 Dieldrin	10.933	10 942 -0.009	89238	0 02201 0	.00220109(aR)
20 Endrin	11.375	11.375 0.000	77463	0 02195	.00219482 (aR)
18 Toxaphene	Com	pound Not Detect	ed.		

Data File: /var/chem/gc3.i/2170-E.b/c-a2373.d Report Date: 19-May-2000 14:13

						CONCENTRATIONS			
						ON-	COLUMN		PINAL
Compounds		RT	EXP RT	DLT RT	RESPONSE	(	ng)		( mg/L)
	# # # # # # # # # # # # # # # # # # #	22			n=#500#=	==			2269633
21 4,4	- ססס	11.767	11.767	0.000	63681	0.	02182	0	00218243 (a)
22 Ende	osulfan II	11.842	11.850	-0.008	81978	0	02314	0	00231439(a)
23 4,4	-DDT	12.317	12.317	0.000	62411	٥.	02252	0	00225 <b>177(a</b> R)
24 End	rın aldehyde	12 717	12 717	0 000	61310	0	02087	0,	00208679(a)
26 Endo	sulfan sulfate	13.583	13.583	0.000	48290	0.	01576	٥.	00157623 (a)
25 Metl	noxychlor	13.442	13.450	-0.008	38614	0.	02772	0-	00277245 (a)
27 Endi	rin ketone	14.083	14.083	0 000	83156	0	02410	7	00240988(a)
\$ 1 Tet:	rachloro-m-xylene	4.558	4 558	0 000	73519	0	018عج	٥.	00181572 (aR)
\$ 30 Dec	achlorobiphenyl	16.292	16.300	-0.008	64216	ο.	01949	0	00194929 (aR)

# QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).
   R Spike/Surrogate failed recovery limits.



Data File: /var/chem/gc3.i/2170-E.b/c-b2373.d

Report Date: 19-May-2000 14:14

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-b2373.d

Lab Smp Id: DDE2E102 Client Smp ID: LCS

Inj Date : 19-MAY-2000 13:44

Operator: 1891 Inst ID: gc3.i

Smp Info : DDE2E102,2170-E.b,,PEST.sub,,3,

Misc Info: 130192LCS

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date: 19-May-2000 11:24 eppinged Quant Type: ESTD Cal Date: 18-MAY-2000 01:57 Cal File: c-b2300.d

Als bottle: 1 QC Sample: LCS

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF		Dilution Factor
٧t	10.000	Volume of final extract (uL)
Vo	100.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected

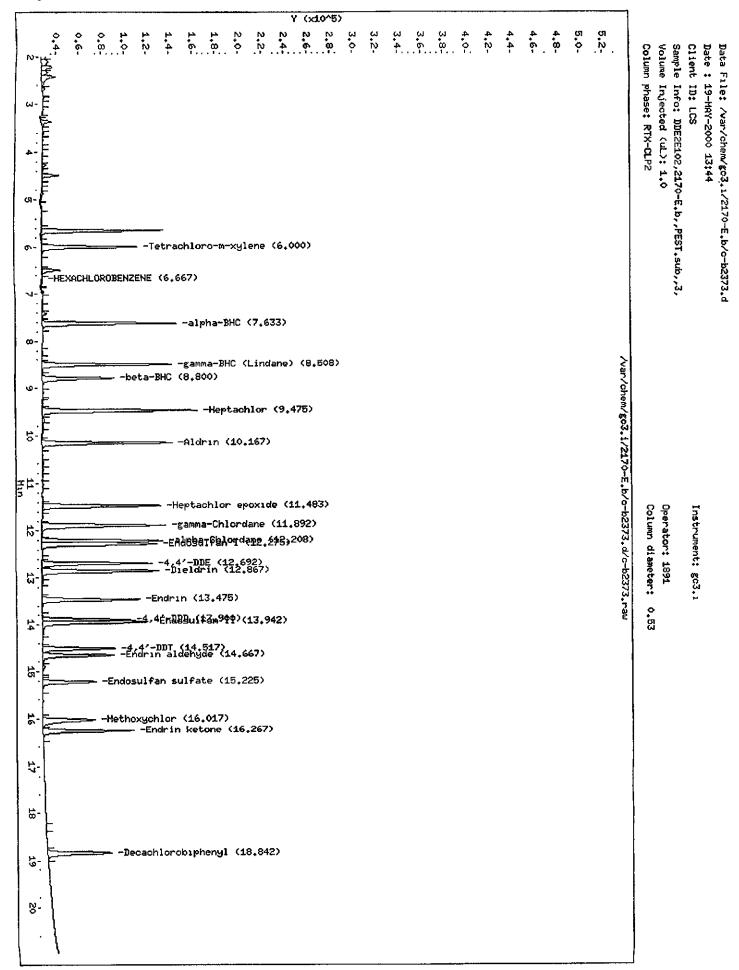
						Concentrations			
						on-column	FINAL		
Compo	undo	RT	BXP RT	DLT RT	RESPONSE	( ng)	(mg/L)		
===#=	d##############	a≠	220000	*****	=======	*****	******		
5	alpha-BHC	7 633	7.633	0.000	117972	0.01852	0.00185188(a)		
4	HEXACHLOROBENZENE	6 667	6.650	0 017	810		(a)		
6	gamma-BHC (Lindane)	8 508	8.508	0.000	114032	0.02117	0.00211724 (aR)		
7	beta-BHC	8.800	8.800	0.000	62664	0.02289	0.00228950(a)		
9	Chlordane	Con	pound No	ot Detect	ed.				
10	Heptachlor	9 475	9.475	0.000	135791	0 02567	0.00256682 (aR)		
8	delta-BHC	Com	pound No	ot Detect	ed.	_			
11	Aldrin	10.167	10.167	0.000	114348	0.02231	0 00223064(aR)		
12	Heptachlor epoxide	11 483	11.483	0.000	103705	0 02366	0.00236606(a)		
13	gamma-Chlordane	11.692	11.892	0.000	109170	0 02376	0 00237642(a)		
14	alpha-Chlordane	12.208	12.208	0.000	106404	0.02376	0.00237636(a)		
15	Endosulfan I	12.275	12.283	-0.008	100824	0.02498	0.00249772(a)		
16	4,4'-DDB	12.692	12.692	0.000	97024	0.02471	0.00247120(a)		
17	Dieldrin	12.867	12.867	0.000	102946	0.02480	0.00247968(aR)		
20	Endrin	13 475	13.475	0.000	86094	0.02432	0.00243208 (aR)		
18	Toxaphene	Com	pound No	ot Detect	eđ.				

Data File: /var/chem/gc3.i/2170-E.b/c-b2373.d Report Date: 19-May-2000 14:14

						CONCENTRATIONS
						on-column final
Comp	ounds	RT	BXP RT	DLT RT	rbsponsb	(ng) $(mg/L)$
	E########	25	*====	=====	#======	******
2	1 4,4'-DDD	13 900	13.900	0.000	77540	0.02575 0.00257473(a)
2	2 Endosulfan II	13.942	13.950	-0 008	92357	0 02706 0.00270559(a)
2	3 4,4'-DDT	14.517	14.517	0.000	64247	0.02452 0.00245197(aR)
2	4 Endrin aldehyde	14.667	14 667	0.000	62861	0.02260 0.00226046(a)
2	6 Bndosulfan sulfate	15.225	15.225	0 000	47106	0.01675 0.00167466(a)
2	5 Methoxychlor	16.017	16.017	0.000	46356	0 03708 0.00370844(a)
2	7 Endrin ketone	16.267	16.267	0 000	80005	0.02592 0.00259195(a)
\$	1 Tetrachloro-m-xylene	6.000	6.000	0.000	83297	0.02027 0.00202690 (aR)
\$ 3	0 Decachlorobiphenyl	18.842	18.842	0 000	S62S2	0 02080 0.00207954 (aR)

# QC Flag Legend

- a Target compound detected but, quantitated amount
   Below Limit Of Quantitation(BLOQ).
   R Spike/Surrogate failed recovery limits.



## 10A PESTICIDE IDENTIFICATION SUMMARY FOR SINGLE COMPONENT ANALYTES

LCS
-----

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT

Case No.: 40325

SDG No.: C0E130194

Lab Sample ID: DDE2E102

Date(s) Analyzed: 05/19/00 05/19/00

Instrument ID (1): GC3

Instrument ID (2): GC3

GC Column(1): RTX-CLP ID: 0.53 (mm) GC Column(2): RTX-CLP2 ID: 0.53 (mm)

	227	RT WINDOW				2-
ANALYTE	COL	R <b>T</b> =====	FROM	TO	CONCENTRATION	%D
alpha-BHC	1	5.88	5.81	5.91	0.00166213	
	2	7.63	7.56	7.66	0.00185188	11.4
beta-BHC	1	6.90	6.86	6.96	0.00201078	
	2	8.80	8.75	8.85	0.00228950	13.9
gamma-BHC (Lindane)	1	6.62	6.55	6.65	0.00188671	
	2	8.51	8.43	8.53	0.00211724	12.2
Heptachlor	1	7.67	7.60	7.70	0.00211506	
	2	9.47	9.40	9.50	0.00256682	21.4
Aldrin	1	8.30	8.25	8.35	0.00198905	
	2	10.17	10.12	10.22	0.00223064	12.1
Heptachlor epoxide	1	9.65	9.60	9.70	0.00218908	
	2	11.48	11.43	11.53	0.00236606	8.1
Endosulfan I	1	10.43	10.35	10.45	0.00219413	
	2	12.28	12.20	12.30	0.00249772	13.8
Dieldrin	1	10.93	10.86	10.96	0.00220109	
	2	12.87	12.79	12.89	0.00247968	12.6

page 1 of 3

FORM X PEST-1

OLM03.0

10A

EPA SAMPLE NO.

PESTICIDE IDENTIFICATION SUMMARY 657 496 FOR SINGLE COMPONENT ANALYTES

LCS

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: COE130194

Lab Sample ID: DDE2E102

Date(s) Analyzed: 05/19/00 05/19/00

Instrument ID (1): GC3

Instrument ID (2): GC3

GC Column(1): RTX-CLP ID: 0.53(mm) GC Column(2): RTX-CLP2 ID: 0.53(mm)

			RT WINDOW				
COL	RT	FROM	TO	CONCENTRATION	%D		
1	10.55	10.50	10.60	0.00224116			
2	12.69	12.64	12.74	0.00247120	10.3		
1	11.38	11.30	11.40	0.00219482			
2	13.47	13.40	13.50	0.00243208	10.8		
1	11.84	11.80	11.90	0.00231439			
2	13.94	13.90	14.00	0.00270559	16.9		
1	11.77	11.69	11.79	0.00218243			
2	13.90	13.82	13.93	0.00257473	18.0		
_	13.30	23.52					
1	13.58	13.53	13.63	0.00157623			
2	15.22	15.18	15.28	0.00167466	6.2		
_	13.22	13.20					
1	12.32	12.24	12.34	0.00225177			
2	14 52	14 44	14 54	0.00245197	8.9		
	14.52	14.44	14.54	0.00213137			
1	13.44	13.38	13.47	0.00277245			
					33.8		
4	16.02	15.94	16.04	0.003/0844	33.0		
1	14 09	14 03	14.13	0.00240988			
2	16.27	16.22	16.32	0.00259195	7.6		
	1 2 1 2 1 2 1 2	1 10.55 2 12.69 1 11.38 2 13.47 1 11.84 2 13.94 1 11.77 2 13.90 1 13.58 2 15.22 1 12.32 1 12.32 1 14.52 1 13.44 2 16.02 1 14.08	1       10.55       10.50         2       12.69       12.64         1       11.38       11.30         2       13.47       13.40         1       11.84       11.80         2       13.94       13.90         1       11.77       11.69         2       13.90       13.82         1       13.58       13.53         2       15.22       15.18         1       12.32       12.24         2       14.52       14.44         1       13.44       13.38         2       16.02       15.94         1       14.08       14.03	1       10.55       10.50       10.60         2       12.69       12.64       12.74         1       11.38       11.30       11.40         2       13.47       13.40       13.50         1       11.84       11.80       11.90         2       13.94       13.90       14.00         1       11.77       11.69       11.79         2       13.90       13.82       13.93         1       13.58       13.53       13.63         2       15.22       15.18       15.28         1       12.32       12.24       12.34         2       14.52       14.44       14.54         1       13.44       13.38       13.47         2       16.02       15.94       16.04         1       14.08       14.03       14.13	1       10.55       10.50       10.60       0.00224116         2       12.69       12.64       12.74       0.00247120         1       11.38       11.30       11.40       0.00219482         2       13.47       13.40       13.50       0.00243208         1       11.84       11.80       11.90       0.00231439         2       13.94       13.90       14.00       0.00270559         1       11.77       11.69       11.79       0.00218243         2       13.90       13.82       13.93       0.00257473         1       13.58       13.53       13.63       0.00157623         2       15.22       15.18       15.28       0.00167466         1       12.32       12.24       12.34       0.00225177         2       14.52       14.44       14.54       0.00245197         1       13.44       13.38       13.47       0.00277245         2       16.02       15.94       16.04       0.00370844         1       14.08       14.03       14.13       0.00240988		

page 2 of 3

FORM X PEST-1

OLM03.0

### 10A PESTICIDE IDENTIFICATION SUMMARY FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

LCS

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325

SDG No.: C0E130194

Lab Sample ID: DDE2E102

Date(s) Analyzed: 05/19/00 05/19/00

Instrument ID (1): GC3

Instrument ID (2): GC3

GC Column(1): RTX-CLP ID: 0.53(mm) GC Column(2): RTX-CLP2 ID: 0.53(mm)

	RT WINDOW						
ANALYTE	COL	RT	FROM	TO	CONCENTRATION	%D	
******************	===	=====	=====	=====	=======================================	=====	
Endrin aldehyde	1	12.72	12.67	12.77	0.00208679		
	2	14.67	14.62	14.72	0.00226046	8.3	
alpha-Chlordane	1	10.20	10.15	10.25	0.00212446		
	2	12.21	12.16	12.26	0.00237636	11.8	
gamma-Chlordane	1	9.92	9.88	9.97	0.00212158		
	2	11.89	11.84	11.94	0.00237642	12.0	
	1		<del></del>				
	2	<del></del>					
	1			•			
	2						
	1						
	2			<del></del> :			
	1						
	2						
	1		<del></del> .				
	2						

page 3 of 3

FORM X PEST-1

OLMO3.0

## UXB INTERNATIONAL MATRIX SPIKE COMPOUNDS

657 498

Lab Name. Severn Trent Laboratories, Inc SDG Number:

Matrix: (soil/water) SOLID WATER Lab Sample ID:COE130194 001

Method: SW846 8081A

un 5/03/00

Pesticides (8081A)

Sample WT/Vol: 100 / mL Work Order: DD6A411A Dilution factor: 1

Date Received · 05/13/00 Date Extracted: 05/18/00 Date Analyzed: 05/19/00

QC Batch: 0139460

Client Sample Id: DF/S1/0133/SDC/001 A

## CONCENTRATION UNITS ·

CAS NO.	COMPOUND (ug/L	or ug/kg) mg/L Q
72-20-8	Endrin	0.00212
76-44-8	Heptachlor	0.00236
1024-57-3	Heptachlor epoxide	0.00201
58-89-9	Lindane	0.00172
72-43-5	Methoxychlor	0.00289

Data File: /var/chem/gc3.i/2170-E.b/c-a2367.d

Report Date: 19-May-2000 12:39

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2367.d Lab Smp Id: DD6A411A Client Client Smp ID: DF/S1/0133/SDC/MS

Inj Date : 19-MAY-2000 11:08

Operator : 1891 Inst ID: qc3.i

Smp Info : DD6A411A,2170-E.b,,PEST.sub,,3,

Misc Info: 1301940018

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m Meth Date : 19-May-2000 11:01 eppinged Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-a2300.d

Als bottle: 1 OC Sample: MS

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF Vt Vo Vi	10.000	Dilution Factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected

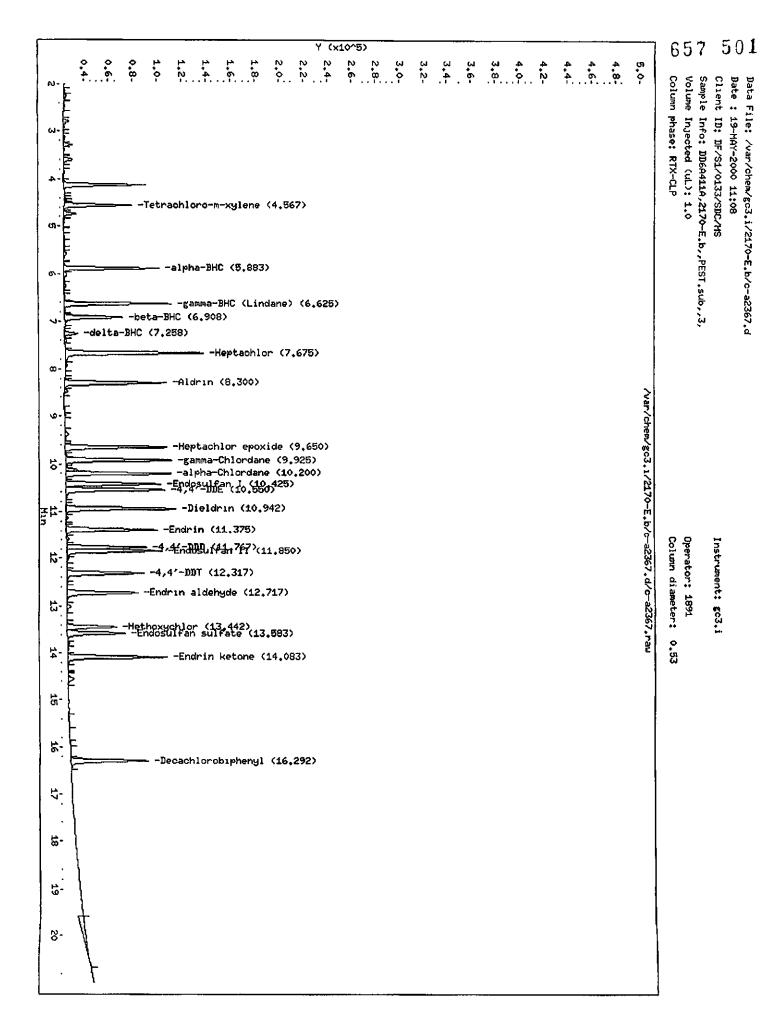
					CONCENT	RATIONS
					ON-COLUMN	FINAL
Compounds	RT	EXP RT D	LT RT	RESPONSE	( ng)	( mg/L)
					****	
5 alpha-BHC	5.883	5 883	0.000	78343	0.01408	0.00140826(a)
4 HEXACHLOROBENZENE	Соп	pound Not	Detect	ed.		
6 gamma-BHC (Lindane)	6.625	6.625	0.000	87566	0.01716	0.00171606 (aR)
7 beta-BHC	6.908	6 908	0.000	47433	0 01958	0.00195786(a)
9 Chlordane	Com	pound Not	Detect	ed.		
10 Heptachlor	7.675	7.675	0.000	113894	0 02376	0.00237622 (aR)
8 delta-BHC	7 258	7 258	0 000	11036	0 00235	0.000234820(a)
11 Aldrin	8.300	8.300	0.000	84159	0.01888	0.00188790(aR)
12 Heptachlor epoxide	9 650	9 650	0.000	83884	ر0.02010	0.00200991(a)
13 gamma-Chlordane	9 925	9 925	0 000	87619	0.02020	0.00202034(a)
14 alpha-Chlordane	10.200	10.200	0.000	87109	0.02031	0.00203083(a)
15 Endosulfan I	10.425	10.425	0.000	78573	0.02036	0.00203642(a)
16 4,4'-DDE	10 550	10.550	0 000	81784	0.02246	0.00224563(a)
17 Dieldrin	10.942	10.942	0.000	90402	0.02230	0 00222980(aR)
20 Endrin	11.375	11.375	0.000	74965	0.02124	0.00212405 (aR)
18 Toxaphene	Con	pound Not	Detect	ed.		

Data File: /var/chem/gc3.i/2170-E.b/c-a2367.d Report Date: 19-May-2000 12:39

						CONCENT	CATIONS
						ON-COLUMN	FINAL
Compo	unds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( mg/L)
52005		22		****			
21	4,4'-DDD	11.767	11 767	0 000	65696	0.02251	0.00225148(a)
22	Endosulfan II	11 850	11 850	0.000	78500	0.02216	0.00221620(a)
23	4,4'-DDT	12 317	12.317	0.000	63119	0.02277	0.00227731 (aR)
24	Endrin aldehyde	12.717	12.717	0 000	58410	0.01988	0.00198808(a)
26	Endosulfan sulfate	13.583	13 583	0 000	47876	0.01563	0 00156272(a)
25	Methoxychlor	13 442	13.450	-0.008	40276	0.02892	0.00289178(a)
27	Endrin ketone	14 083	14.083	0 000	82062	0.02378	0.00237817(a)
\$ 1	Tetrachloro-m-xylene	4 567	4.558	0 009	56113	سطِّ0.0138	0 00138584 (aR)
\$ 30	Decachlorobiphenyl	16.292	16 300	-0.008	63785	0.0193	0.00193620(aR)

## QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).R Spike/Surrogate failed recovery limits.



Data File: /var/chem/gc3.i/2170-E.b/c-b2367.d

Report Date: 19-May-2000 12:46

#### STL-PITTSBURGH

Data file : /var/chem/gc3.i/2170-E.b/c-b2367.d

Lab Smp Id: DD6A411A Client Smp ID: DF/S1/0133/SDC/MS

Inj Date : 19-MAY-2000 11:08

Operator : 1891 Inst ID: gc3.i

Smp Info : DD6A411A, 2170-E.b, , PEST. sub, , 3,

Misc Info: 130194001S

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date : 19-May-2000 11:24 eppinged Quant Type: ESTD Cal Date : 18-MAY-2000 01:57 Cal File: c-b2300.d

Als bottle: 1 QC Sample: MS

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF	1.000	Dilution Factor
۷t	10.000	Volume of final extract (uL)
Vo	100.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected

					CONCENT	RATIONS
					OM-COLUMN	FINAL
Compounds	RT	BXP RT	DLT RT	response	( ng)	( mg/L)
	==		sec===	******	- 10 11 11 11 11 11 11 11 11 11 11 11 11	======
5 alpha-BHC	7.633	7.633	0.000	103590	0.01626	0.00162612(a)
4 HEXACHLOROBENZENE	6.692	6 650	0.042	1327		(a)
6 gamma-BHC (Lindane)	8.508	8.508	0.000	104134	0.0193	0.00193346 (aR)
7 beta-BHC	8.800	8.800	0.000	59026	0 02157	0.00215658(a)
9 Chlordane	Com	pound N	ot Detect	ed.		
10 Heptachlor	9.475	9.475	0.000	124703	0.02357	0 00235722 (aR)
8 delta-BHC	Com	pound N	ot Detect	ed	_	
11 Aldrin	10 175	10.167	0.008	99958	0 01950	0.00194993 (aR)
12 Heptachlor epox1de	11 483	11.483	0.000	97087	0.02215	0.00221507(a)
13 gamma-Chlordane	11 892	11 892	0 000	101932	0.02219	0.00221886(a)
14 alpha-Chlordane	12 208	12 208	0.000	100679	0.02249	0.00224850(a)
15 Endosulfan I	12.283	12.283	0 000	93033	0.02305	0.00230472(a)
16 4,4'-DDB	12.692	12.692	0.000	96677	0.02462	0.00246236(a)
17 Dieldrin	12.867	12.867	0 000	103941	0.02504	0 00250364(aR)
20 Endrin	13.475	13.475	0 000	83874	0.02369	9.00236936 (aR)
18 Toxaphene	Соп	pound N	ot Detect	ed.	~	

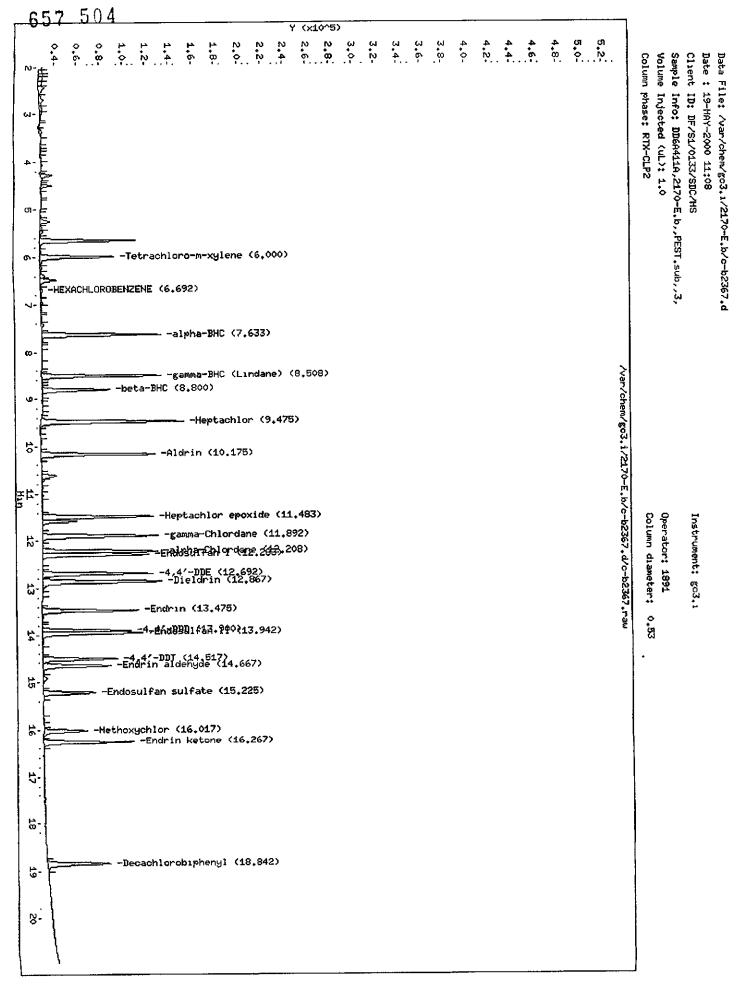
Data File: /var/chem/gc3.i/2170-E.b/c-b2367.d Report Date: 19-May-2000 12:46

							ATIONS
						on-column	FINAL
C	ompounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( mg/L)
=	*********	39.79					
	21 4,4'-DDD	13 900	13.900	0 000	77880	0.02586	0.00258602(a)
	22 Bndosulfan II	13.942	13.950	-0 008	87922	0.02576	0 00257566(a)
	23 4,4'-DDT	14.517	14.517	0.000	65556	0 02502	0.00250193 (aR)
	24 Endrin aldehyde	14.667	14.667	0.000	59377	0 02135	0.00213518(a)
	26 Bndosulfan sulfate	15.225	15.225	0.000	46000	0.01635	0.00163534(a)
	25 Methoxychlor	16.017	16.017	0.000	38052	0.03044	0.00304413(a)
	27 Endrin ketone	16.267	16 267	0 000	78718	0.02550	0.00255025(a)
\$	1 Tetrachloro-m-xylene	6.000	6.000	0 000	65108	0 01584	0.00159430(aR)
\$	30 Decachlorobiphenyl	18 842	18.842	0 000	54644	0.02020	0 00202009 (aR)

CONCENTRATIONS

# QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).
   R Spike/Surrogate failed recovery limits.



657 505

10A EPA SAMPLE NO PESTICIDE IDENTIFICATION SUMMARY

DF/S1/01

Lab Name: STL-PITTSBURGH

Contract:

33/SDC/MS

Lab Code: STLPIT Case No.: SAS No.: 40325

SDG No.: C0E130194

Lab Sample ID: DD6A411A

Date(s) Analyzed: 05/19/00 05/19/00

Instrument ID (1): GC3

Instrument ID (2): GC3

GC Column(1): RTX-CLP ID: 0.53 (mm) GC Column(2): RTX-CLP2 ID: 0.53 (mm)

FOR SINGLE COMPONENT ANALYTES

			RT W	MDOW		
ANALYTE	COL	RT	FROM	TO	CONCENTRATION	%D
	===	=====	=====	=====	=========	=====
alpha-BHC	1	5.88	5.81	5.91	0.00140826	
	2	7.63	7.56	7.66	0.00162612	15.5
beta-BHC	1	6.91	6.86	6.96	0.00195786	
	2	8.80	8.75	8.85	0.00215658	10.1
gamma-BHC (Lindane)	1	6.62	6.55	6.65	0.00171606	
	2	8.51	8.43	8.53	0.00193346	12.7
Heptachlor	1	7.67	7.60	7.70	0.00237622	
	2	9.47	9.40	9.50	0.00235722	0.8
Aldrin	1	8.30	8.25	8.35	0.00188790	
	2	10.18	10.12	10.22	0.00194993	3.3
Heptachlor epoxide	1	9.65	9.60	9.70	0.00200991	
	2	11.48	11.43	11.53	0.00221507	10.2
Endosulfan I	1	10.43	10.35	10.45	0.00203642	
	2	12.28	12.20	12.30	0.00230472	13.2
Dieldrin	1	10.94	10.86	10.96	0.00222980	
	2	12.87	12.79	12.89	0.00250364	12.3

page 1 of 3

FORM X PEST-1

10A PESTICIDE IDENTIFICATION SUMMARY FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

DF/S1/01 33/SDC/MS

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: COE130194

Date(s) Analyzed: 05/19/00 05/19/00

Lab Sample ID: DD6A411A

Instrument ID (2): GC3

Instrument ID (1): GC3

GC Column(1): RTX-CLP ID: 0.53(mm) GC Column(2): RTX-CLP2 ID: 0.53(mm)

			RT W	MDOW		
ANALYTE	COL	RT	FROM	TO	CONCENTRATION	%D
=======================================	===	=====	=====	=====	=======================================	=====
4,4'-DDE	1	10.55	10.50	10.60	0.00224563	
	2	12.69	12.64	12.74	0.00246236	9.6
Endrin	1	11.38	11.30	11.40	0.00212405	
	2	13.47	13.40	13.50	0.00236936	11.5
Endosulfan II	1	11.85	11.80	11.90	0.00221620	ļ
	2	13.94	13.90	14.00	0.00257566	16.2
4,4'-DDD	1	11.77	11.69	11.79	0.00225148	
	2	13.90	13.82	13.93	0.00258602	14.8
Endosulfan sulfate	1	13.58	13.53	13.63	0.00156272	
	2	15.22	15.18	15.28	0.00163534	4.6
4,4'-DDT	1	12.32	12.24	12.34	0.00227731	
	2	14.52	14.44	14.54	0.00250193	9.9
Methoxychlor	1	13.44	13.38	13.47	0.00289178	
	2	16.02	15.94	16.04	0.00304413	5.3
Endrin ketone	1	14.08	14.03	14.13	0.00237817	
	2	16.27	16.22	16.32	0.00255025	7.2

page 2 of 3

FORM X PEST-1

## 10A PESTICIDE IDENTIFICATION SUMMARY FOR SINGLE COMPONENT ANALYTES

DF/S1/01 33/SDC/MS

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.:

SAS No.: 40325

SDG No.: C0E130194

Lab Sample ID: DD6A411A

Date(s) Analyzed: 05/19/00 05/19/00

Instrument ID (1): GC3

Instrument ID (2): GC3

GC Column(1): RTX-CLP ID: 0.53(mm) GC Column(2): RTX-CLP2

ID: 0.53 (mm)

ANALYTE	CONCENTRATION	0.70				
MUCHANISH SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SERVICES SE	COL	RT	FROM	TO	CONCENTRATION	%D =====
Endrin aldehyde	1	12.72	12.67	12.77	0.00198808	
	2	14.67	14.62	14.72	0.00213518	7.4
alpha-Chlordane	1	10.20	10.15	10.25	0.00203083	
	2	12.21	12.16	12.26	0.00224850	10.7
gamma-Chlordane	1	9.93	9.88	9.97	0.00202034	
	2	11.89	11.84	11.94	0.00221886	9.8
	1			<del></del>	-	
	2			<del></del>		
	1		<del></del>			
	2	<del></del>	<del></del>			l
	1					
	2					
	1					
	2			<u></u>		
	1					
	2					

page 3 of 3

FORM X PEST-1

657 508

## UXB INTERNATIONAL MATRIX SPIKE DUPLICATE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WARL

Lab Sample ID:COE130194 001

Method: SW846 8081A Lm563 W

Pesticides (8081A)

Sample WT/Vol: 100 / mL Work Order: DD6A411C Dilution factor: 1

Date Received: 05/13/00 Date Extracted:05/18/00 Date Analyzed: 05/19/00

QC Batch: 0139460

Client Sample Id: DF/S1/0133/SDC/001 A

## CONCENTRATION UNITS:

CAS NO.	COMPOUND (	ug/L or ug/kg) mg/L Q
72-20-8	Endrin	0 00213
76-44-8	Heptachlor	0.00238
1024-57-3	Heptachlor epoxide	0.00199
58-89-9	Lindane	0.00172
72-43-5	Methoxychlor	0.00291

Data File: /var/chem/gc3.i/2170-E.b/c-a2368.d

Report Date: 19-May-2000 12:40

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-a2368.d

Lab Smp Id: DD6A411C Client Smp ID: DF/S1/0133/SDC/MSD

Inj Date : 19-MAY-2000 11:34 Operator : 1891 Smp Info : DD6A411C,2170-E.b,,PEST.sub,,3, Inst ID: gc3.i

Misc Info : 130194001D

Comment

Method : /var/chem/gc3.i/2170-E.b/PESTA.m

Meth Date: 19-May-2000 11:01 eppinged Quant Type: ESTD Cal Date: 18-MAY-2000 01:57 Cal File: c-a2300 QC Sample: MSD Cal File: c-a2300.d

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF Vt Vo Vi	10.000 100.000	Dilution Factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected

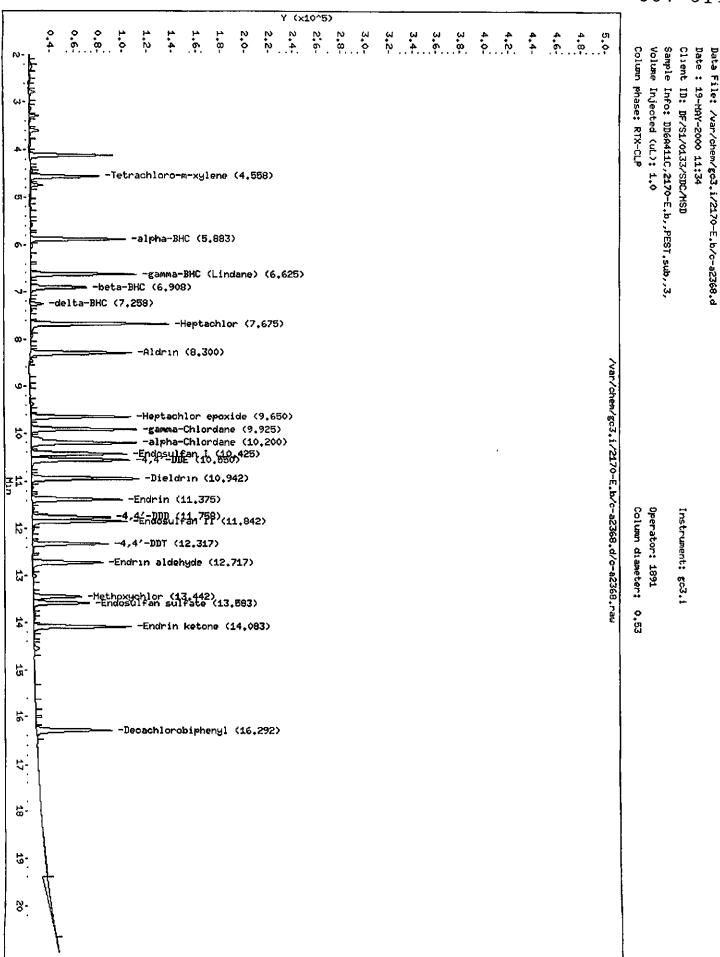
					CONCENT	RATIONS
					ON-COLUMN	FINAL
Compounds	RT	EXP RT D	LT RT	response	( ng)	( mg/L)
B. 10 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1				*******	****	
5 alpha-BHC	5.883	5.883	0.000	79047	0 01421	0.00142092(a)
4 HEXACHLOROBENZENE	Com	pound Not	Detecte	ed.		
6 gamma-BHC (Lindane)	6.625	6 625	0.000	87919	0.01723	0.00172298 (aR)
7 beta-BHC	6.908	6.908	0 000	46766	0.01930	0.00193033(a)
9 Chlordane	Com	pound Not	Detecte	ed.		
10 Heptachlor	7.675	7 675	0.000	114171	0.02382	0.00238200(aR)
8 delta-BHC	7.258	7.258	0 000	11043	0.00235	0.000234969(a)
11 Aldrin	8.300	8.300	0.000	84375	0.01893	0.0 <u>0189</u> 275(aR)
12 Heptachlor epoxide	9.650	9.650	0.000	83234	0.01994	0.00199433(a)
13 gamma-Chlordane	9.925	9.925	0.000	87283	0.02013	0 00201259(a)
14 alpha-Chlordane	10.200	10.200	0.000	87566	0.02041	0.00204148(a)
15 Endosulfan I	10 425	10 425	0.000	79279	0 02055	0.00205472(a)
16 4,4'-DDE	10 550	10.550	0.000	81886	0 02248	0.00224843(a)
17 Dieldrin	10 942	10 942	0.000	88925	0 02193	0.00219337(aR)
20 Endrin	11.375	11.375	0.000	75188	0 02130	0.00213037 (AR)
18 Toxaphene	Соп	pound Not	Detect	ed	•	

Data File: /var/chem/gc3.i/2170-E.b/c-a2368.d Report Date: 19-May-2000 12:40

						CONCENTI	RATIONS
						ON - COLUMN	FINAL
Compoun	da	RT	EXP RT	DLT RT	RESPONSE	( ng)	( mg/L)
3255582		12 22	225582	20222		****	== # = # = #
21 4	, 4 <sup>1</sup> - DDD	11 758	11 767	-0 009	65055	0 02230	0.00222952(a)
22 E	ndosulfan II	11.842	11 850	-0 008	78371	0.02213	0.00221256(a)
23 4	,4'-DDT	12 317	12 317	0.000	62682	0.02262	0.00226154(aR)
24 E	ndrin aldehyde	12.717	12.717	0.000	58199	0.01981	0.00198090(a)
26 E	ndosulfan sulfate	13 583	13 583	0.000	47034	0.01535	0.00153523(a)
25 M	ethoxychlor	13 442	13 450	-0 008	40595	0.02915	0.00291468(a)
27 E	ndrin ketone	14 083	14.083	0.000	80634	0.02337	0.00233679(a)
\$ 1 T	etrachloro-m-xylene	4 558	4.558	0.000	57052	0.01409	0.00140903 (aR)
\$ 30 D	ecachlorobiphenyl	16.292	16.300	-0.008	62801	0.01906	0.00190634 (aR)

## QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).
   R Spike/Surrogate failed recovery limits.



## 657 512

Data File: /var/chem/gc3.i/2170-E.b/c-b2368.d

Report Date: 19-May-2000 12:46

#### STL-PITTSBURGH

Data file: /var/chem/gc3.i/2170-E.b/c-b2368.d

Lab Smp Id: DD6A411C Client Smp ID: DF/S1/0133/SDC/MSD

Inj Date : 19-MAY-2000 11:34

Operator : 1891 Inst ID: gc3.i

Smp Info : DD6A411C, 2170-E.b, , PEST. sub, , 3,

Misc Info: 130194001D

Comment :

Method : /var/chem/gc3.i/2170-E.b/PESTB.m

Meth Date: 19-May-2000 11:24 eppinged Quant Type: ESTD Cal Date: 18-MAY-2000 01:57 Cal File: c-b2300.d

Als bottle: 1 QC Sample: MSD

Dil Factor: 1.00000
Integrator: Falcon Compound Sublist: PEST.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* (Vt/Vo)/Vi

Name	Value	Description
DF	1.000	Dilution Factor
۷t	10.000	Volume of final extract (uL)
Vo	100.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected

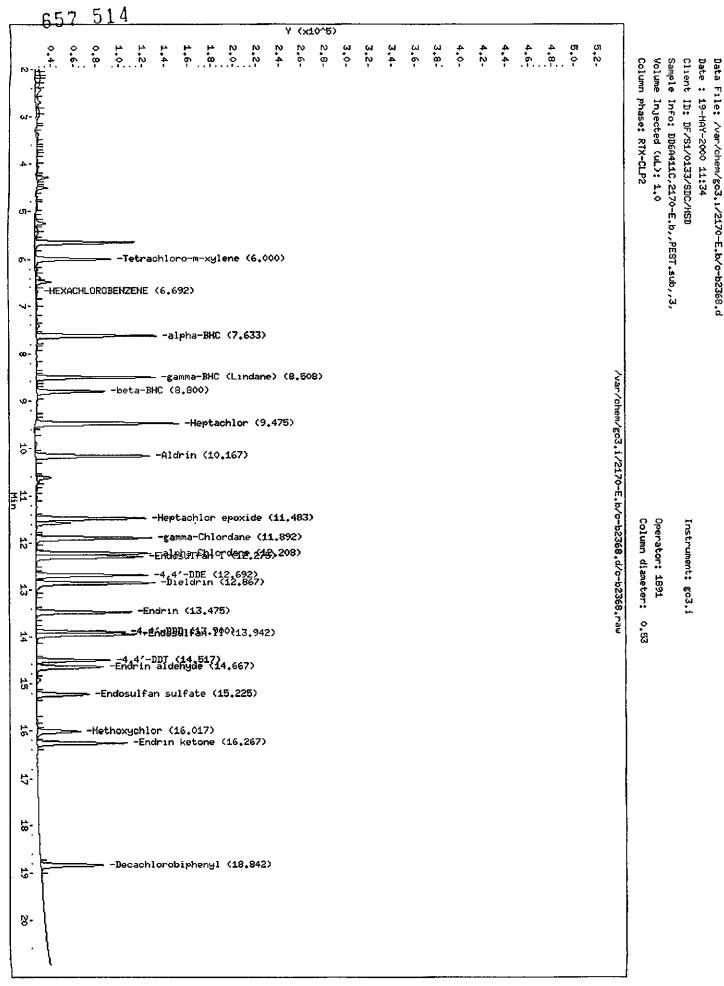
						CONCENT	RATIONS
						on-column	PINAL
Compou	ınds	RT	EXP RT	DLT RT	response	( ng)	( mg/L)
						======	22222
s	alpha-BHC	7.633	7.633	0 000	104056	0.01633	0.00163343(a)
4	HEXACHLOROBENZENS	6 692	6 650	0 042	1580		(a)
6	gamma-BHC (Lindane)	8.508	8.508	0.000	103697	0.01925	0.00192535 (aR)
7	beta-BHC	8.800	8.800	0.000	59482	0.02173	0.00217324(a)
9	Chlordane	Con	pound No	ot Detect	ed.		
10	Heptachlor	9.475	9.475	0.000	126057	0.02383	0.00238282 (aR)
8	delta-BHC	Con	pound No	ot Detect	ed.	_	
11	Aldrin	10 167	10 167	0.000	99731	0.01946	0.00194550(aR)
12	Heptachlor epoxide	11 483	11,483	0.000	96301	0 02197	0.00219714(a)
13	gamma-Chlordane	11 892	11 892	0.000	101686	0.02214	0.00221350(a)
14	alpha-Chlordane	12 208	12 208	0 000	100700	0 02249	0 00224897(a)
15	Endosulfan I	12.275	12 283	-0.00B	93587	0.02318	0 00231844(a)
16	4,4'-DDB	12.692	12.692	0 000	97818	0.02491	0.00249142(a)
17	Dieldrin	12.867	12.867	0 000	103369	0.02490	0.002489 <u>87(aR)</u>
20	Endrin	13.475	13.475	0.000	83511	سو 0.0235	0.00235911 (aR)
18	Toxaphene	Con	spound No	ot Detect	ed.	_	

Data File: /var/chem/gc3.i/2170-E.b/c-b2368.d Report Date: 19-May-2000 12:46

						CONCENT	RATIONS	
							ON-COLUMN	FINAL
Co	mpoı	ande	RT	EXP RT	DLT RT	response	( ng)	( mg/L)
==		************	==	=====			Charass	****
	21	4,4'-DDD	13.900	13.900	0.000	77560	0.02575	0 00257540(a)
	22	Bndosulfan II	13.942	13.950	-0 008	87926	0.02576	0 00257578(a)
	23	4,4'-DDT	14.517	14.517	0 000	65441	0.02498	0.00249754 (aR)
	24	Endrin aldehyde	14.667	14.667	0.000	58983	0.02121	0.00212101(a)
	26	Endosulfan sulfate	15.225	15.225	0.000	46383	0.01649	0.00164896(a)
	25	Methoxychlor	16 017	16.017	0.000	38247	0.03060	0.00305973(B)
	27	Bndrin ketone	16 267	16.267	0.000	79066	0.02562	0.00256153(a)
Ş	1	Tetrachloro-m-xylene	6.000	6.000	0.000	66649	0.01622	0 00162186 (aR)
\$	30	Decachlorobiphenyl	18 842	18.842	0.000	55151	0.02039	0.00203883 (aB)

## QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).R Spike/Surrogate failed recovery limits.



<u>657</u> 515 EPA SAMPLE NO.

## 10A PESTICIDE IDENTIFICATION SUMMARY FOR SINGLE COMPONENT ANALYTES

DF/S1/013 3/SDC/MSD

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.:

SAS No.: 40325

SDG No.: C0E130194

Lab Sample ID: DD6A411C

Date(s) Analyzed: 05/19/00 05/19/00

Instrument ID (1): GC3

Instrument ID (2): GC3

GC Column(1): RTX-CLP ID: 0.53(mm) GC Column(2): RTX-CLP2

ID: 0.53 (mm)

			RT W	NDOW		
ANALYTE	COL	RT	FROM	TO	CONCENTRATION	%D
**=====================================	===	=====	=====	=====	<b>#</b> #==========	=====
alpha-BHC	1	5.88	5.81	5.91	0.00142092	
	2	7.63	7.56	7.66	0.00163343	15.0
beta-BHC	1	6.91	6.86	6.96	0.00193033	
	2	8.80	8.75	8.85	0.00217324	12.6
gamma-BHC (Lindane)	1	6.62	6.55	6.65	0.00172298	
	2	8.51	8.43	8.53	0.00192535	11.7
Heptachlor	1	7.67	7.60	7.70	0.00238200	
	2	9.47	9.40	9.50	0.00238282	0.0
Aldrin	1	8.30	8.25	8.35	0.00189275	
	2	10.17	10.12	10.22	0.00194550	2.8
Heptachlor epoxide	1	9.65	9.60	9.70	0.00199433	
	2	11.48	11.43	11.53	0.00219714	10.2
Endosulfan I	1	10.43	10.35	10.45	0.00205472	
	2	12.28	12.20	12.30	0.00231844	12.8
Dieldrin	1	10.94	10.86	10.96	0.00219337	
	2	12.87	12.79	12.89	0.00248987	13.5

page 1 of 3

FORM X PEST-1

10A PESTICIDE IDENTIFICATION SUMMARY FOR SINGLE COMPONENT ANALYTES

EPA SAMPLE NO.

DF/S1/013 3/SDC/MSD

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: C0E130194

Lab Sample ID: DD6A411C

Date(s) Analyzed: 05/19/00 05/19/00

Instrument ID (1): GC3

Instrument ID (2): GC3

GC Column(1): RTX-CLP ID: 0.53(mm) GC Column(2): RTX-CLP2 ID: 0.53(mm)

			RT W	NDOW		
ANALYTE	COL	RT	FROM	TO	CONCENTRATION	%D
	===	=== <b>===</b>	=====	=====	=======================================	=====
4,4'-DDE	1	10.55	10.50	10.60	0.00224843	
	2	12.69	12.64	12.74	0.00249142	10.8
Endrin	1	11.38	11.30	11.40	0.00213037	
	2	13.47	13.40	13.50	0.00235911	10.7
Endosulfan II	1	11.84	11.80	11.90	0.00221256	
	2	13.94	13.90	14.00	0.00257578	16.4
4,4'-DDD	1	11.76	11.69	11.79	0.00222952	
	2	13.90	13.82	13.93	0.00257540	15.5
Endosulfan sulfate	1	13.58	13.53	13.63	0.00153523	į
	2	15.22	15.18	15.28	0.00164896	7.4
4,4'-DDT	1	12.32	12.24	12.34	0.00226154	
	2	14.52	14.44	14.54	0.00249754	10.4
Methoxychlor	1	13.44	13.38	13.47	0.00291468	
	2	16.02	15.94	16.04	0.00305973	5.0
Endrin ketone	1	14.08	14.03	14.13	0.00233679	
	2	16.27	16.22	16.32	0.00256153	9.6

page 2 of 3

FORM X PEST-1

#### 10A PESTICIDE IDENTIFICATION SUMMARY FOR SINGLE COMPONENT ANALYTES

DF/S1/013 3/SDC/MSD

Lab Name: STL-PITTSBURGH

Contract:

Lab Code: STLPIT

Case No.:

SAS No.: 40325

SDG No.: C0E130194

Lab Sample ID: DD6A411C

Date(s) Analyzed: 05/19/00 05/19/00

Instrument ID (1): GC3

Instrument ID (2): GC3

GC Column(1): RTX-CLP ID: 0.53(mm) GC Column(2): RTX-CLP2

ID: 0.53 (mm)

-	г		יונו ייינו	INDOW		<del></del> ,
ANALYTE	COL	RT	FROM	TO	CONCENTRATION	%D
	===	=====	88888	=====	=======================================	=====
Endrin aldehyde	1	12.72	12.67	12.77	0.00198090	
	2	14.67	14.62	14.72	0.00212101	7.1
alpha-Chlordane	1	10.20	10.15	10.25	0.00204148	
	2	12.21	12.16	12.26	0.00224897	10.2
gamma-Chlordane	1	9.93	9.88	9.97	0.00201259	
	2	11.89	11.84	11.94	0.00221350	10.0
	1					
	2					i
	1					
	2					
	1					
	2					
	1					
	2		<u> </u>			
	1				•	
	2					

page 3 of 3

FORM X PEST-1

# PESTICIDE MISCELLANEOUS

TCLP (Method 1311) Quanter 450 Wi.

Quanterra Incorporated 450 William Pitt Way Pittsburgh, Pennsylvania 15238 412/826-5477 FAX: 412/826-5571

**W**ua

Environmental Services

* = Sample determined to have free-liquid, % solid determination was performed  <5 = Extraction Fluid No. 1 57 mL Glacial Acetic Acid dil 500 mL + 64 3 mL of 1N Na OH dil to 1L (pH 4.93 ± 0.05)  >5 = Extraction Fluid No. 2 5.7 mL Glacial Acetic Acid dil to 1L (pH 2.88 ± 0.05)  \$\times 0.05 \tag{0.0 \tag{0.0} \tag{0.0} \tag{0.0}	5-100 12:15	(Riscord line number from above)  Date  Time  Availyst		8	73.16.00	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\			DICAG 4.59 1.79 4.50 1.70 4.50 1.70	9.44 1.73	Sample ID WgIVOI Init. pH Final pH Extract Fund to Acc  ODC A4 < 00, /8L C.   Q. C.	1	١ ٠	holling. Sant 5-16-00 creed by	GOUA-17	F- 412/826-5477 FAX: 412/826-5571
S-16-00   S'45   S-17-00   9:30   Agitation Apparatus RPMs: S-16-00   S'45   S-17-00   9:30   30±2   Sm.	5-16-0-12:00 1111-18	Location Date Time //// Analyst / Location	10	0	9 (3)	8.	50D6A4 V V V. 10.30	( 5))%,7	7	100, MY 100	to Acidity Sample ID WevDry Fluid Vol Tumbling (tor Metals)  1. DP-TCLP I-16-00 WevDry Fluid Vol Tumbling 1. 193		- 1 Solution Nur	5-17-00 Accument ARDS SN908	st # 8320701W Santal Number 038	

Sodium Sulfate Mfg. Conanterra N.\QA\LOGBOOKS\sepfun extr.doc May-00 Date Extraction Began Extraction Worksheet Analyst 14 3 Extract(s)

Record line number from above)

A A A & C COF Lot Number Aprove 30194 Date Completed Da DOIMS Sample ID 00/msp 003 004 LC5 5 स्वव 00 2756 Date Parameter 814 00000 Time **6** Client ID É Lot Number Extract(s) Received Method 3510C Hq6 B#0139460 10000h 1000 Sample Volume (mL) BiP OP-00-0035 0.0 Final Volume (ml.) δ F TAXANC Surrogate Number 8-06-06 Reviewed By 5-19.00 5-18-0 Solvent Lot ~44274 かれる Surrogate Volume (mL) 1.0m B.P. 03100 Time Extract(s) Relinquished Matrix Spike No. Solvent Mile SAK Clean up Method 90-85-8 8.0 NA Analysi STL Pittsburgh 450 William Pitt Way Pittsburgh, PA 15238 Matrix Spike Volume (mL) Ø 1.0m Page 12 of 80 FIRE KC Cleanup Date 0

Separatory Funnel

3200

657 521 Turbochrom Sequence File : H:\ACQUIRE\MET SEQ\2170-E.SEQ

Created by : DE11/02/98 Edited by : DE05/17/00

on: 5/17/00on: 5/17/0022:34

16:37

Description: QUANTERRA PGH 8081 RUN ON GC#3 CLP1/CLP2 COLUMNS

REVIEWED BY:

Number of Times Edited: 2

## Sequence File Header Information:

Number of Rows : 77

Instrument Type : 760 / 900 Series Intelligent Interface Injection Type : SINGLE

				Sequence Samp	le Descri	ptions -	Channel A					
Row	Туре	Sample Name	Sample Number	Study Name	Sample Amount	ISTD Amount	Sample Volume	Dil. Factor	Mult	Divisor	Addend	Norm. factor
1	Std Check	EVALB, 2170-E.b,	190-88-8	·	1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
2	Cal Replace	MEDTOX, 2170-E.b	190-84-13		1.000	1.000	1.000	1.000	1.000	1 000	0.000	100.000
3	Cal Replace	MEDCHLOR, 2170-E	190-85-10		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
4	Cal Replace	LOWAPPX9,2170-E	190-80-6		1.000	1 000	1.000	1.000	1.000	1.000	0.000	100.000
5	Cai.Replace	MLOWAPPX9,2170-	190-80-7		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
6	Cal.Replace	MEDAPPX9,2170-E	190-80-8		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
7	Cal:Replace	MHIGHAPPX9,2170	190-80-9		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
8	Cal Replace	HIGHAPPX9,2170-	190-80-10		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
9	Cal:Replace	LOWA, 2170-E.b.,	190-84-1	Pour outage	1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
10 11	Cal Replace	MLOWA, 2170-E.b.	190-84-2	1	1.000	1.000	1.000	1 000	1.000	1.000	0.000	100.000
11	Cal.Replace	MEDA, 2170-E.b,,	190-84-3		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
12	Cal:Replace	MHIGHA, 2170-E.b	190-84-4		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
13	Cal:Replace	HIGHA, 2170-E.b,	190-84-5		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
14	Cal:Replace	LOWB, 2170-E.b.,	190-84-7		1 000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
15	Cal Replace	MLOWB, 2170-E b,	190-84-8		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
16	Cal:Replace	MEDB, 2170-E.b.,	190-84-9		1.000	1.000	1.000	1.000	1.000	1.000	0 000	100.000
17 18	Cal:Replace	MHIGHB, 2170-E.b	190-84-10		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
19	Cal:Replace Std Check	HIGHB, 2170-E.b, 2ND A, 2170-E.b,	190-84-11 190-82-2		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
20	Std Check	2ND B,2170-E.b,	190-82-5		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
21	Std Check	EVALB, 2170-E.b,	190-88-8		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100 000
22	Sample	DD3QT10P,2170-E	110261005		1.000	1.000	1.000	1.000 1.000	1.000	1.000	0.000	100 000
23	Sample	DD3QV10P,2170-E	110126006		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
24	Sample	DD3QX124,2170-E	110261007		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100 000
25	Sample	DD3QX125,2170-E	1102610078		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000 100.000
26	Sample	DD3QX126,2170-E	110261007D.		1 000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
27	Sample	DD3R010P,2170-E	110261008		1 000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
<u>27</u>	Sample	DD5GR101,2170-E	110261BLK		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
- 29	Sample	DD5GR102,2170-E	110261LCS		1.000	1.000	1.000	1.000	1.000	1 000	0.000	100.000
30	Sample	DD11N10A, 2170-E	100168001		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
31	Sample	DD11V10A,2170-E	100168002	,	1.000	1.000	1.000	1.000	1.000	1 000	0.000	100 000
32	Sample	DD11X10A,2170-E	100168003		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
33	Sample	DD12310A, 2170-E	100168004 🗸		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100 000
34	Sample	DD12410A,2170-E	100168005		1.000	1.000	1.000	1.000	1.000	1 000	0.000	100 000
35	Sample	DD12510A,2170-E	100168006		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100 000
36	Sample	DD12810A,2170-E	100168008		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
37	Sample	DD12C10A, 2170-E	100168009		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
38	Sample	DD12D10A, 2170-E	100168010	,	1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
39	Sample Sample	DD1R5101,2170-E	100168BLK		1.000	1.000	1.000	1.000	1.000	1 000	0.000	100.000
40	Sample Sample	DD1R5102,2170-E DD1R5103,2170-E	100168LCS 100168LCD		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
41 42	Std Check	MEDAPPX9,2170-E	190-80-8		1.000	1.000	1.000	1.000	1.000	1 000	0.000	100 000
43	Std Check	MEDA, 2170-E.b.,	190-84-3		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
44		MEDB, 2170-E.b,,	190-84-9		1.000	1.000	1.000	1.000 1.000	1.000	1.000	0 000	100.000
45	Std Check	EVALB, 2170-E.b,	190-88-8		1.000	1.000	1.000	1 000	1.000	1.000	0.000	100.000
46	Sample	DD36010A,2170-E	110206001		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000 100.000
47	•	DD36810A,2170-E			1.000	1.000	1.000	1.000	1.000	1.000		100.000
48	Sample	DD36910A, 2170-E			1.000	1 000	1.000	1.000	1.000	1.000	0 000	100.000
49	Sample	DD36A10A, 2170-E	110206004		1.000	1 000	1.000	1.000	1.000	1.000	0.000	100 000
50	Sample	•	110206005		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
51	Sample	DD36D10A,2170-E	110206006		1.000	1 000	1.000	1.000	1.000	1.000	0.000	100 000
52	Sample		110206007		1.000	1.000	1 000	1.000	1.000	1.000	0 000	100.000
53	Sample	DD36G10A,2170-E	110206008		1 000	1.000	1 000	1.000	1.000	1.000	0 000	
54	Sample	DD36L10A,2170-E	110206009		1.000	1.000	1.000	1.000	1 000		0.000	
55	Sample	DD36N10A,2170-E	110206010		1.000	1.000	1.000	1.000	1.000	1.000	0.000	
56	Sample	DD36V10A,2170-E	110206011	,	1.000	1.000	1.000	1.000	1.000	1.000	0.000	
57	Sample		110206012		1.000	1.000	1 000	1.000	1.000	1.000	0.000	100.000
58	Sample	DD5Q31QA,2170-E	130132001	•	1.000	1.000	1.000	1.000	1 000	1.000	0.000	100 000
STE <sup>9</sup>	Firetsbur	gP050410A,2170-E	130132002		1.000	1 000	1 000	1.000	1.000	1 000	0.000	3°20°1°
		-										J = V =

65	57 522										
60	Sample	DD5Q510A,2170-E	130132003	1.000	1.000	1 000	1.000	1.000	1.000	0.000	100.000
61	Sample	DD5Q710A,2170-E	130132004	1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
62	Sample	DD5Q810A,2170-E	130132005	1,000	1.000	1.000	1.000	1.000	1.000	0.000	100 000
63	Sample	DD91V101,2170-E	130132BLK	1.000	1.000	1,000	1 000	1.000	1.000	0.000	100.000
64	Sample	DD91V102,2170-E	130132LCS	1 000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
65	Sample	DD91V103,2170-E	130132LCD	1 000	1.000	1.000	1.000	1 000	1.000	0.000	100 000
66	Std Check	MEDA, 2170-E.b.,	190-84-3	1,000	1.000	1.000	1.000	1.000	1.000	0.000	100 000
67	Std Check	MEDB, 2170-E b,	190-84-9	1,000	1.000	1.000	1.000	1.000	1.000	0.000	100 000
68	Std Check	EVALB, 2170-E b,	190-88-8	1 000	1.000	1.000	1.000	1.000	1 000	0.000	100.000
69	Sample	DD5QA10A,2170-E	130132006	1.000	1.000	1.000	1.000	1.000	1.000	0.000	100 000
70	Sample	DD5QC10A, 2170-E	130132007	1.000	1.000	1.000	1.000	1 000	1.000	0 000	100.000
71	Sample	DD5QE10A, 2170-E	130132008	1,000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
72	Sample Sample	DD50F10A, 2170-E	130132009	1,000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
73	Sample	DD5QH10A,2170-E	130132010	1 000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
74	Sample	DD5QK10A, 2170-E	130132011	1 000	1.000	1.000	1.000	1,000	1.000	0.000	100.000
75	Sample	DD5QM10A,2170-E	130132012	1.000	1.000	1 000	1.000	1.000	1.000	0.000	100.000
76	-	MEDA, 2170-E.b.,	190-84-3	1.000	1.000	1.000	1 000	1.000	1.000	0.000	100.000
	Std Check	MEDB, 2170-E.b,,	190-84-9	1.000	1.000	1.000	1,000	1.000	1.000	0.000	100 000
77	Std Check	MEDD, 2110-5.0) /	150 04 5		2.305						

								ocess Infor							
Row	Site	Rack	Vıal	Method	Process Method	Calib Method	Report Format	Raw File	Result File	File	Modified Raw File	Rpt	Level Name	Update RT	Out Dev
1	<b></b>	1	1	GEN3C	GEN3A	122190A	EVAL	C-A2281	C-A2281		C-A2281	-	_	-	LPT1.
2	_	1	2	GEN3C	GEN3A	122190A	TOX	C-A2282	C-A2282		C-A2282	Ŋ	MED	N	LPT1 · , LP
3	_	1	2	GEN3C	GEN3A	122190A	TOX	C-A2283	C-A2283		C-A2283	N	MED	N	LPT1:,LP
4	-	ī	4	GEN3C	GEN3A	122190A	INDA	C-A2284	C-A2284		C-A2284	N	LOW	N	LPT1:
5	_	1	5	GEN3C	GEN3A	122190A	INDA	C-A2285	C-A2285		C-A2285	N	MLOW	N	LPT1:
6	-	1	6	GEN3C	GEN3A	122190A	INDA	C-A2286	C-A2286		C-A2286	И	MLOW	N	LPT1:
7	-	1	7	GEN3C	GEN3A	122190A	INDA	C-A2287	C-A2287		C-A2287	N	MLOW	N N	LPT1:
8	-	1	8	GEN3C	GEN3A	122190A	INDA	C-A2288	C-A2288		C-A2288 C-A2289	N	MLOW LOW	N	LPT1: LPT1:
9	-	1	4	GEN3C	GEN3A	122190A	INDA	C-A2289 C-A2292	C-A2289 C-A2292		C-A2292	N	MLOW	N	LPT1:
10	-	1	5	GEN3C	GEN3A	122190A 122190A	INDA INDA	C-A2292	C-A2292		C-A2293	N	MLOW	N	LPT1:
11	-	1	6 7	GEN3C	GEN3A GEN3A	122190A 122190A	INDA	C-A2294	C-A2294		C-A2294	N	MLOW	N	LPT1:
12 13	-	l 1	8	GEN3C GEN3C	GEN3A	122190A	INDA	C-A2295	C-A2295		C-A2295	N	MLOW	N	LPT1:
14	_	i	9	GEN3C	GEN3A	122190A	INDA	C-A2296	C-A2296		C-A2296	N	LOW	N	LPT1:
15	_		10	GEN3C	GEN3A	122190A	INDA	C-A2297	C-A2297		C-A2297	N	MLOW	N	LPT1:
16	-	ì	11	GEN3C	GEN3A	122190A	INDA	C-A2298	C-A2298		C-A2298	N	MLOW	И	LPT1:
17	_	ī	12	GEN3C	GEN3A	122190A	INDA	C-A2299	C-A2299		C-A2299	N	MLOW	N	LPT1:
18	_	1	13	GEN3C	GEN3A	122190A	INDA	C-A2300	C-A2300		C-A2300	И	MLOW	N	LPT1:
19	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2301	C-A2301		C-A2301	-	**	-	LPT1:
20	-	1	24	GEN3C	GEN3A	122190A	INDA	C-A2302	C-A2302		C-A2302	-	-	-	LPT1:
21	-	1	1	GEN3C	GEN3A	122190A	EVAL	C-A2303	C-A2303		C-A2303	-	-	-	LPT1:
22	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2304	C-A2304		C-A2304	-		-	LPT1:
23	-	1	23	GEN3C	gen3a	122190A	INDA	C-A2305	C-A2305		C-A2305	-	-	-	LPT1:
24	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2306	C-A2306		C-A2306 C-A2307	-	_	-	LPT1: LPT1:
25	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2307 C-A2308	C-A2307 C-A2308		C-A2307	_	_	_	LPT1:
26		1	23	GEN3C	GEN3A	122190A 122190A	INDA	C-A2309	C-A2309		C-A2309	_	_	_	LPT1:
27	-	1	23	GEN3C	GEN3A GEN3A	122190A 122190A	INDA INDA	C-A2309	C-A2310		C-A2310	_	_	_	LPT1:
28 29	-	1	23 23	GEN3C GEN3C	GEN3A GEN3A	122190A	INDA	C-A2311	C-A2311		C-A2311	_	-	-	LPT1:
30	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2312	C-A2312		C-A2312	-	-	-	LPT1:
31	_	1	23	GEN3C	GEN3A	122190A	INDA	C-A2313	C-A2313		C-A2313	-	-	-	LPT1:
32		1	23	GEN3C	GEN3A	122190A	INDA	C-A2314	C-A2314		C-A2314	-	-	-	LPT1.
33		1	23	GEN3C	GEN3A	122190A	INDA	C-A2315	C-A2315		C-A2315	-	-	-	LPT1:
34		1	23	GEN3C	GEN3A	122190A	INDA	C-A2316	C-A2316		C-A2316	-	_	-	LPT1:
35	_	1	23	GEN3C	<b>GEN3A</b>	122190A	INDA	C-A2317	C-A2317		C-A2317	-	-	-	LPT1:
36	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2318	C-A2318		C-A2318	-	_	-	LPT1:
37		1	23	GEN3C	GEN3A	122190A	INDA	C-A2319	C-A2319		C-A2319	-	-	-	LPT1.
38		1	23	GEN3C	GEN3A	122190A	INDA	C-A2320	C-A2320		C-A2320 C <del>-</del> A2321	_	_	_	LPT1.
39		1	23	GEN3C	GEN3A	122190A	INDA	C-A2321 C-A2322	C-A2321 C-A2322		C-A2321	_	_	-	LPT1:
40		1	23	GEN3C	GENGA	122190A 122190A	INDA INDA	C-A2322	C-A2322		C-A2323	-	_	_	LPT1:
41		1 1	23 6	GEN3C GEN3C	gen3a gen3a	122190A	INDA	C-A2324	C-A2324		C-A2324	_	-	_	LPT1:
42 43		1	6	GEN3C	GEN3A	122190A	INDA	C-A2325	C-A2325		C-A2325	_	-	_	LPT1:
44		1	11	GEN3C	GEN3A	122190A	INDA	C-A2326	C-A2326		C-A2326	-	-	-	LPT1:
45		ī	1	GEN3C	GEN3A	122190A	EVAL	C-A2327	C-A2327		C-A2327	-	_	-	LPT1 ·
46		1	23	GEN3C	GEN3A	122190A	INDA	C-A2328	C-A2328		C-A2328	-	-	-	LPT1
47		1	23	GEN3C	GEN3A	122190A	INDA	C-A2329	C-A2329		C-A2329	-	-	-	LPT1.
48	_	1	23	GEN3C	GEN3A	122190A	INDA	C-A2330	C-A2330		C-A2330	-	-	-	LPT1:
49	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2331	C-A2331		C-A2331	-	-	-	LPT1:
50		1	23	GEN3C	GEN3A	122190A	INDA	C-A2332	C-A2332		C-A2332	-	-	-	LPT1:
51		1	23	GEN3C	GEN3A	122190A	INDA	C-A2333	C-A2333		C-A2333	-	-	-	LPT1: LPT1
52		1	23	GEN3C	GEN3A	122190A	INDA	C-A2334	C-A2334		C-A2334	_	-	_	LPT1:
53		ı	23	GEN3C	GEN3A	122190A	INDA	C-A2335	C-A2335		C-A2335 C-A2336		_	_	LPT1:
54		1	23	GEN3C	GEN3A	122190A	INDA	C-A2336 C-A2337	C-A2336 C-A2337		C-A2336 C-A2337	_	_	_	LPT1.
55		1	23	GEN3C	GEN3A	122190A 122190A	INDA INDA	C-A2338	C-A2337		C-A2338	_	_	-	LPT1:
56 57		1	23 23	GEN3C GEN3C	GEN3A GEN3A	122190A	INDA	C-A2339	C-A2339		C-A2339	-	_	-	LPT1:
58		1 1	23	GEN3C GEN3C	GEN 3A	122190A	INDA	C-A2340	C-A2340		C-A2340	-	-	-	LPT1:
58 58		1	23	GEN3C GEN3C	GENSA GENSA	122190A	INDA	C-A2341	C-A2341		C-A2341	-	-	-	LPT1:
5: 60		1	23	GEN3C	GEN3A	122190A		C-A2342	C-A2342		C-A2342	-	-	-	LPT1:
۷.	١ _	1	23	дензс Э <b>ст</b> изс	GEN3A	122190A		C-A2343	C-A2343		C-A2343	-	-	-	lpT1. 3≥2102
						122190A		C-A2344	C-A2344		C-A2344		-		

63	-	1	23	GEN3C	GEN3A	122190a	INDA	C-A2345	C-A2345	C-A2345	_		657	523
		-	23								-	-	00 -	TIT!
64	-	1		GEN3C	gen3a	122190A	INDA	C-A2346	C-A2346	C-A2346	_	-	-	LPT1:
65	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2347	C-A2347	C-A2347	-	-	-	LPT1:
66	-	1	6	GEN3C	GEN3A	122190A	INDA	C-A2348	C-A2348	C-A2348	_	_		LPT1:
67	-	1	11	GEN3C	GEN3A	122190A	INDA	C-A2349	C-A2349	C-A2349	-	-	-	LPT1:
68	-	1	1	GEN3C	gen3a	122190A	EVAL	C-A2350	C-A2350	C-A2350	-	-	-	LPT1
69	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2351	C-A2351	C-A2351	-	_	•	LPT1:
70	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2352	C-A2352	C-A2352	-	-	-	LPT1.
71	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2353	C-A2353	C-A2353	-	-	_	LPT1.
72	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2354	C-A2354	C-A2354	-	-	_	LPT1
73	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2355	C-A2355	C-A2355	_	-	_	LPT1:
74	-	1	23	GEN3C	GEN3A	122190A	INDA	C-A2356	C-A2356	C-A2356	_	••	_	LPT1:
75	•	1	23	GEN3C	GEN3A	122190A	INDA	C-A2357	C-A2357	C-A2357	-	-	_	LPT1:
76	-	1	6	GEN3C	GEN3A	122190A	INDA	C-A2358	C-A2358	C-A2358	-	-	_	LPT1.
77	-	1	11	GEN3C	GEN3A	122190A	INDA	C-A2359	C-A2359	C-A2359	-	-	-	LPT1:

Turbochrom Sequence File: H:\ACQUIRE\MET\_SEQ\2190-E.SEQ

Created by : DE11/02/98 on : 5/19/00 10:00 Edited by : DE05/19/00 on : 5/19/00 10:07

Description: QUANTERRA PGH 8081 RUN ON GC#3 CLP1/CLP2 COLUMNS

REVIEWED BY:

Number of Times Edited: 1

## Sequence File Header Information:

Number of Rows : 15

Instrument Type : 760 / 900 Series Intelligent Interface
Injection Type : SINGLE .

				Sequence Samp	ole Descri	ptions -	Channel A					-
Row	Туре	Sample	Sample	Study Name	Sample	ISTD	Sample	Dil.	Mult	Divisor	Addend	Norm.
	-28-	Name	Number		Amount	Amount	Volume	Factor				factor
						1.000	1,000	1.000	1.000	1.000	0,000	100 000
1	Std Check	EVALB,2170-E.b,	190-88-8		1.000				1.000	1.000	0.000	100 000
2	Std Check	MEDTOX,2170-E.b	190-84-13		1.000	1.000	1.000	1.000				
3	Std Check	MEDCHLOR, 2170-E	190-85-10		1.000	1.000	1.000	1.000	1.000	1.000	0 000	100.000
Ā	Std Check	MEDA, 2170-E.b.,	190-84-3		1.000	1.000	1.000	1.000	1 000	1.000	0.000	100.000
	Std Check	MEDB. 2170-E.b.,	190-84-9		1.000	1.000	1,000	1.000	1.000	1.000	0 000	100.000
<del>- }</del>		DD6A4103,2170-E	130194001		1.000	1,000	1.000	1.000	1.000	1.000	0.000	100.000
ఫ	Sample		-		1.000	1.000	1.000	1 000	1.000	1.000	0 000	100.000
7	Sample	DD6A411A, 2170-E	130194001S			1.000	1,000	1 000	1,000		0.000	100.000
8	Şample	DD6A411C,2170-E	130194001D		1 000				1.000		0.000	100.000
9	Sample	DD6A5103,2170-E	130194002		1.000	1.000	1.000	1.000				
<u>10</u>	Sample	DD6A6103,2170-E	130194003		1.000	1 000	1.000	1.000	1.000		0.000	100.000
11	Sample	DD6A7103.2170-E	130192004		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100 000
12	Sample	DDE2E101,2170-E	130192BLK		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100 000
.13	•	DDE2E102,2170-E	130192LCS		1.000	1.000	1,000	1.000	1.000	1.000	0.000	100 000
	Sample	•	190-84-3		1.000	1.000		1.000	1.000	1.000	0.000	100.000
14	Std Check	MEDA, 2170-E.b.,				1.000		1.000	1.000		0.000	
15	Std Check	MEDB,2170-E b,,	190-84-9		1.000	1.000	1 000	1.000	1.000	1.000	0.000	100.000

Row	Site	Rack	Vıal	Inst Method	Process Method	Se Calib Method	equence Pro Report Format	ocess Infor Raw File	mation - ( Result File	Channel A Baseline File	Modified Raw File	Cal Rpt	Level Name	Update RT	Out Dev
1		1	1	GEN3C	GEN3A	122190A	EVAL	C-A2361	C-A2361		C-A2361	-	-	_	LPT1:
2	_	ī	2	GEN3C	GEN3A	122190A	TOX	C-A2362	C-A2362		C-A2362	-	-	-	LPT1:,LP
3		ī	2	GEN3C	GEN3A	122190A	TOX	C-A2363	C-A2363		C-A2363	-	-	-	LPT1:,LP
4	_	ī	6	GEN3C	GEN3A	122190A	INDA	C-A2364	C-A2364		C-A2364	-	-	-	LPT1:
5	-	ī	11	GEN3C	GEN3A	122190A	INDA	C-A2365	C-A2365		C-A2365	-	-	-	LPT1:
6	_	1	23	GEN3C	GEN3A	122190A	INDA	C-A2366	C-A2366		C-A2366	-	-	-	LPT1:
7	_	1	23	GEN3C	GEN3A	122190A	INDA	C-A2367	C-A2367		C-A2367	-	-	-	LPT1:
8	_	1	23	GEN3C	GEN3A	122190A	INDA	C-A2368	C-A2368		C-A2368	-	-	-	LPT1:
9	_	ī	23	GEN3C	GEN3A	122190A	INDA	C-A2369	C-A2369		C-A2369	-	-	-	LPT1.
10		ī	23	GEN3C	GEN3A	122190A	INDA	C-A2370	C-A2370		C-A2370	-	-	-	LPT1:
11	_	1	23	GEN3C	GEN3A	122190A	INDA	C-A2371	C-A2371		C-A2371	-	-	-	LPT1:
12	_	1	23	GEN3C	GEN3A	122190A	INDA	C-A2372	C-A2372		C-A2372	-	-	-	LPT1:
13	_	1	23	GEN3C	GEN3A	122190A	INDA	C-A2373	C-A2373		C-A2373	-	-	_	LPT1:
14		ī	6	GEN3C	GEN3A	122190A	INDA	C-A2374	C-A2374		C-A2374	-	-	-	LPT1:
15		ī	11	GEN3C	GEN3A	122190A	INDA	C-A2375	C-A2375		C-A2375	-	-	-	LPT1:

657 525 .

Turbochrom Sequence File : H:\ACQUIRE\MET SEQ\2170-E.SEQ

Created by: DE11/02/98 on: 5/17/00 16:37 Edited by: DE05/17/00 on: 5/17/00 22:34

Description: QUANTERRA PGH 8081 RUN ON GC#3 CLP1/CLP2 COLUMNS

REVIEWED BY:

Number of Times Edited: 2

## 110673

## Sequence File Header Information:

Number of Rows : 77

Instrument Type : 760 / 900 Series Intelligent Interface
Injection Type : SINGLE

Row	Туре	Sample Name	Sample Number	Sequence Sampl Study Name	le Descri Sample Amount	ptions - ISTD Amount	Channel B Sample Volume	Dil. Factor	Mult	Divisor	Addend	Norm. factor
1	Std Check	EVALB, 2170-E.b,	190-88-8		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
2	Cal·Replace	MEDTOX, 2170-E.b	190-84-13		1.000	1.000	1.000	1 000	333.000	1.000	0.000	100.000
3	Cal:Replace	MEDCHLOR, 2170-E	190-85-10		1.000	1.000	1.000	1 000	333.000	1.000	0.000	100.000
4	Cal:Replace	Lowappx9,2170-E	190-80-6		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
5	Cal:Replace	MLOWAPPX9,2170-	190-80-7		1.000	1.000	1.000	1.000	1 000	1.000	0.000	100 000
6	Cal:Replace	MEDAPPX9,2170-E	190-80-8		1.000	1 000	1.000	1.000	1.000	1.000	0 000	100 000
7	Cal Replace	MHIGHAPPX9,2170	190-80-9		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
9	Cai:Replace	HIGHAPPX9,2170-	190-80-10	0 04.4	1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
9	Cal:Replace	LOWA, 2170-E.b.,	190-84-1	Power Outrige	1.000	1.000	1.000	1.000 1.000	1.000 1.000		0.000	100 000 100.000
10	Cal:Replace	MLOWA, 2170-E.b,	190-84-2	•	1.000	1.000	1.000	1.000	1.000		0.000	
11	Cal:Replace	MEDA, 2170-E.b.,	190-84-4		1.000	1.000	1.000	1.000	1.000		0.000	100.000
12	Cal:Replace	MHIGHA, 2170-E.b HIGHA, 2170-E.b,	190-84-5		1.000	1.000	1.000	1 000	1.000		0.000	
13 14	Cal.Replace Cal:Replace	LOWB, 2170-E.b,	190-84-7		1.000	1.000	1.000	1.000	1.000		0.000	
15	Cal:Replace	MLOWB, 2170-E.b.	190-84-8		1.000	1.000	1.000	1.000	1 000		0.000	
16	Cal:Replace	MEDB, 2170-E.b.	190-84-9		1 000	1.000		1.000	1.000		0 000	
17	Cal Replace	MHIGHB, 2170-E.b	190-84-10		1 000	1 000		1.000	1.000		0 000	
18	Cal:Replace	HIGHB, 2170-E b,	190-84-11		1 000	1.000		1.000	1.000	1.000	0.000	100 000
19	Std Check	2ND A, 2170-E.b.	190-82-2		1.000	1.000	1.000	1 000	1.000	1.000	0.000	100 000
20	Std Check	2ND B, 2170-E.b,	190-82-5		1.000	1.000	1.000	1.000	1.000	1.000	0 000	100 000
21	Std Check	EVALB, 2170-E.b,	190-88-8		1,000	1 000	1.000	1.000	1.000		0.000	
22	Sample	DD3QT10P,2170-E	110261005		1.000	1.000		1.000	1.000		0.000	
23	Sample	DD3QV10P,2170-E	110126006	/	1.000	1.000		1.000	1.000		0 000	
24	Sample	DD3QX124,2170-E	110261007		1.000	1.000		1.000	1.000		0.000	
25	Sample	DD30X125,2170~E	110261007		1.000	1.000		1.000	1.000		0 000	
26	Sample	DD3QX126,2170-E	110261007		1.000	1.000		1.000	1.000		0.000	
27	Sample	DD3R010P,2170-E	110261008		1 000	1.000		1,000	1.000		0.000	
28	Sample	DD5GR101,2170-E	110261BLK		1.000	1.000			1.000		0.000	
29 30	Sample	DD5GR102,2170-E DD11N10A,2170-E	110261LCS 100168001		1.000	1.000			1.000		0.000	
31	Sample Sample	DD11V10A, 2170-E	100168002	•	1.000	1.000			1.000		0.000	
32	Sample	DD11X10A,2170-E	100168003	/	1.000	1.000			1.000		0 000	
33	Sample	DD12310A,2170-E	100168004		1.000	1.000			1.000		0 000	
34	Sample	DD12410A, 2170-E	100168005		1 000	1.000			1.000		0 000	100.000
35		DD12510A, 2170-E	100168006		1.000	1.000	1.000	1 000	1.000	1.000	0 000	100.000
36		DD12810A,2170-E	100168008		1.000	1.000	1.000	1.000	1.000	1.000	0.000	
37	Samole	DD12C10A,2170-E	100168009	_	1.000	1.000			1.000		0.000	
38	Sample	DD12D10A,2170-E	100168010		1 000	1.000			1.000		0.000	
39	•	DD1R5101,2170-E	100168BLK		1 000	1.000			1.000		0.000	
40	•	DD1R5102,2170-E	100168LCS		1.000	1.000			1.000		0 000	
41	•	DD1R5103,2170-E	100168LCD	1	1.000	1.000			1.000		0.000	
42		MEDAPPX9,2170-E	190-80-8		1.000	1.000			1.000		0.000	
43		MEDA, 2170-E.b,, MEDB, 2170-E.b,,	190-84-3 190-84-9		1.000				1.000		0.000	
44		EVALB, 2170-E.b,	190-88-8		1.000				1.000		0.000	
46		DD36010A,2170-E			1.000				1.000		0.000	
47		DD36810A,2170-E			1.000				1.000		0.000	
48		DD36910A,2170-E			1.000				1.000		0 000	100.000
49		DD36A10A,2170-E			1.000				1 000		0.000	100.000
50	•	DD36C10A, 2170-E			1.000				1.00	1.000	0.000	100 000
51	-	D036D10A,2170-E			1 000		1.000	1.000	1 00			
52		DD36F10A,2170-E	110206007		1.000				1.00			
53	Sample	DD36G10A,2170-E			1 000							
54	-	DD36L10A,2170-E			1.000							
55	-	DD36N10A,2170-E			1.000				1.00			
56	•	DD36V10A, 2170-E			1.000							
57	•	DD36W10A,2170-E			1.000							
58	Sample	DD5Q310A,2170-E	130132007		1.000							
STLP2	rittsbur	gh <sup>p050410A, 2170-E</sup>	130132004	•	1.000	. 1.00	1.000	. 1,000		- *.000		° 3205°

526	1 Sar 2 Sar 3 Sar 4 Sar 5 Sar	mple mple mple mple mple mple mple mple	•	OD5Q510A, DD5Q710A, DD5Q810A, DD91V101, DD91V102, DD91V103, MEDA, 2170	2170-E 1 2170-E 1 2170-E 1 2170-E 1 2170-E 1	130132003 130132004 130132005 130132BLK 130132LCS 130132LCD 190-84-3		1 000 1.000 1.000 1.000 1.000 1.000	1,000 1,000 1,000 1,000 1,000 1,000	1.000 1.000 1.000 1.000 1.000 1.000	1 000 1 000 1.000 1.000 1.000 1.000	1.000 1.000 1.000 1.000 1.000 1.000	1.000 1.000 1.000 1.000 1.000 1.000	0.000 0.000 0.000 0.000 0.000	100 000 100 000 100 000 100 000 100 000 100 000 100 000
<b>€</b> €		i Check i Check		MEDB, 2170- EVALB, 217	• •	.90-84 <b>-</b> 9 .90-88 <b>-</b> 8		1.000 1.000	1.000 1.000	1.000	1.000	1.000	1.000	0 000	100 000
6 7		mple mple		DD5QA10A,		130132006		1.000 1.000	1.000 1.000	1.000 1.000	1,000	1.000	1.000		100.000 100.000
	i San	nple		DDSQE10A,	2170-E 1	130132008 🖊	•	1.000	1.000	1.000	1.000	1 000 1 000	1.000		100.000
7		nple aple		DD5QF10A, DD5QH10A,		L30132009 / L30132010 /		1.000 1.000	1.000	1.000 1.000	1.000 1.000	1,000	1.000		100.000
7	-	nple		DD5QK10A,		130132011		1.000	1.000	1.000	1.000	1.000	1.000		100.000
7		nple		DD5QM10A,		130132012		1.000	1.000 1.000	1.000 1.000	1.000 1.000	1.000	1.000		100.000
7 7		i Checi i Checi		MEDB, 2170		190-84 <b>-</b> 3 190-84 <b>-</b> 9		1.000		1.000	1.000	1.000	1.000		100.000
,	, 500	ı Çileci	•	11000,02,0											
							-	cocess Infor			Madi Evad	C-1	T area l	[[adato	Out
Row	Site	Rack	Vial	Inst Method	Process Method	Calıb Method	Report Format	Raw File	Result File	Baseline File	Modified Raw File	Cal Rpt	Level Name	Update RT	Dev
								C-B2281	C-B2281		C-B2281				LPT1
	1 - 2 -	1	1 2	GEN3C GEN3C	GEN3B GEN3B	122190A 122190A	EVAL TOX	C-B2282	C-B2282		C-B2282	И	MED	N	LPT1.,LP
	3 -	1	2	GEN3C	GEN3B	122190A	TOX	C-82283	C-B2283		C-B2283	N	MED	N	LPT1:,LP
	4 -	1	4	GEN3C	GEN3B	122190A	INDA	C-82284	C-B2284		C-B2284	N	LOW	N	LPT1.
	5 -	1	5	GEN3C	GEN3B	122190A	INDA	C-B2285 C-B2286	C-B2285 C-B2286		C-B2285 C-B2286	N	MLOW MLOW	N N	LPT1. LPT1:
	6 - 7 -	1 1	5 7	GEN3C GEN3C	GEN3B GEN3B	122190A 122190A	INDA INDA	C-B2287	C-B2287		C-B2287	Ŋ	MLOW	N	LPT1.
	8 -	1	8	GEN3C	GEN3B	122190A	INDA	C-B2288	C-B2288		C-B2288	N	MLOW	Ŋ	LPT1:
	9 -	1	4	GEN3C	GEN3B	122190A	INDA	C-B2289	C-B2289		C-B2289	N	TOM	N	LPT1:
	0 -	1	5	GEN3C	GEN3B	122190A	INDA	C-B2292	C-B2292		C~B2292 C-B2293	N N	MTOM WTOM	N	LPT1: LPT1:
	1 -	1	6 7	GEN3C GEN3C	GEN3B GEN3B	122190A 122190A	INDA INDA	C-B2293 C-B2294	C-B2293 C-B2294		C-82294	И	MLOM	N	LPT1:
	.2 -	1	8	GEN3C GEN3C	GEN3B	122190A	INDA	C-B2295	C-B2295		C-B2295	N	MLOW	N	LPT1:
	4 -	1	9	GEN3C	GEN3B	122190A	INDA	C-B2296	C-B2296		C-B2296	N	LOW	N	LPT1:
	.5 -	1	10	GEN3C	GEN3B	122190A	INDA	C-B2297	C-B2297		C-B2297	N N	MLOW	N N	LPT1: LPT1:
	.6 -	1	11	GEN3C	GEN3B	122190A 122190A	INDA INDA	C-B2298 C-B2299	C-B2298 C-B2299		C-B2298 C-B2299	N	MLOW MLOW	N	LPT1:
	.7 -	1 1	12 13	GEN3C GEN3C	GEN3B GEN3B	122190A 122190A	INDA	C-B2299	C-B2300		C-B2300	N	MLOW	N	LPT1:
	9 -	ī	23	GEN3C	GEN3B	122190A	INDA	C-B2301	C-B2301		C-B2301	-	-	-	LPT1:
	20 -	1	24	GEN3C	GEN3B	122190A	INDA	C-B2302	C-B2302		C-B2302	-	-	-	LPT1: LPT1:
	21 -	1	1	GEN3C	GEN38	122190A 122190A	EVAL INDA	C-B2303 C-B2304	C-B2303 C-B2304		C-B2303 C-B2304	-	-	-	LPT1:
	?2 - ?3 -	1 1	23 23	GEN3C GEN3C	GEN3B GEN3B	122190A 122190A	INDA	C-B2305	C-B2305		C-B2305	_	-	-	LPT1.
	24 -	î	23	GEN3C	GEN3B	122190A	INDA	C-B2306	C-B2306		C-B2306	-	-	-	LPT1:
	25 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2307	C-B2307		C-B2307 C-B2308	_	_	-	LPT1: LPT1:
	26 <b>-</b> 27	1 1	23 23	GEN3C GEN3C	GEN3B GEN3B	122190A 122190A	INDA INDA	C-B2308 C-B2309	C-B2308 C-B2309		C-B2309	<del></del>	-	*	LPT1:
	27 28 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2310	C-B2310		C-B2310	-	-	-	LPT1:
	29 -	ī	23	GEN3C	GEN3B	122190A	INDA	C-B2311	C-B2311		C-B2311	-	-	-	LPT1.
:	30 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2312	C-B2312 C-B2313		C-B2312 C-B2313	-	-	-	LPT1: LPT1:
	31 - 32 -	1	23 23	GEN3C GEN3C	GEN3B GEN3B	122190A 122190A	INDA INDA	C-B2313 C-B2314	C-B2313 C-B2314		C-B2313	-	-	-	LPT1.
	33 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2315	C-B2315		C-B2315	-	-	-	LPT1.
	34 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2316	C-B2316		C-B2316	-	-	-	LPTI
	35 -	1	23	GEN3C	GEN3B	122190A	INDA INDA	C-B2317 C-B2318	C-B2317 C-B2318		C-82317 C-B2318	-	-	-	LPT1: LPT1:
	36 ~ 37 ~	1	23 23	GEN3C GEN3C	GEN3B GEN3B	122190A 122190A	INDA	C-B2319	C-B2310		C-B2319	-	-	-	LPT1:
	30 -	•	23	GENISC	GENISE	122190A	TNOA	C=B2320	C-B2320		C-82320	_	_	-	LPT1:

2 -	1	2	GEN3C	GEN3B	122190A	TOX	C-B2282	C-B2282	C-82282	64	MED	IN .	LPTI.,LP
3	1	2	GEN3C	GEN3B	122190A	TOX	C-82283	C-B2283	C-B2283	И	MED	Ŋ	LPT1:,LP
							C-B2284	C-B2284	C-B2284	N	LOW	N	LPT1 ·
4 -	1	4	GEN3C	GEN3B	122190A	INDA							
5 -	1	5	GEN3C	GEN3B	122190A	INDA	C-B2285	C-B2285	C-B2285	N	MLOW	N	LPT1.
6 -	ī	6	GEN3C	GEN3B	122190A	INDA	C-B2286	C-B2286	C-B2286	И	MLOW	N	LPT1:
					1001000		C-B2287	C-B2287	C-B2287	N	MLOW	N	LPT1 ·
7 -	1	7	GEN3C	GEN3B	122190A	INDA							
8 -	1	8	GEN3C	GEN3B	122190A	INDA	C-B2288	C-B2288	C-B2288	N	MLOW	И	LPT1:
				GEN3B	122190A	INDA	C-B2289	C-B2289	C-B2289	N	LOW	N	LPT1:
9 -	1	4	GEN3C								MLOW	N	LPT1:
10 -	1	5	GEN3C	GEN3B	122190A	INDA	C-B2292	C-B2292	C~B2292	N			
11 -	1	6	GEN3C	GEN3B	122190A	INDA	C-B2293	C-B2293	C-B2293	N	MLOW	N	LPT1:
					122190A	INDA	C-B2294	C-B2294	C-B2294	N	MLOW	N	LPT1:
12 -	1	7	GEN3C	GEN3B									
13 -	1	8	GEN3C	GEN3B	122190A	INDA	C-B2295	C-B2295	C-B2295	N	MLOW	N	LPT1:
		9	GEN3C	GEN3B	122190A	INDA	C-B2296	C-B2296	C-B2296	N	LOW	N	LPT1:
14 -	1								C-B2297	N	MLOW	N	LPT1:
15 -	1	10	GEN3C	GEN3B	122190A	INDA	C-B2297	C-B2297					
16 -	1	11	GEN3C	GEN3B	122190A	INDA	C-B2298	C-B2298	C-B2298	Ŋ	MLOW	N	LPT1:
10 -							C-B2299	C-B2299	C-B2299	N	MLOW	N	LPT1:
17 -	1	12	GEN3C	GEN3B	122190A	INDA							
18 -	1	13	GEN3C	GEN3B	122190A	INDA	C~B2300	C-B2300	C-B2300	N	MLOW	N	LPT1:
			GEN3C	GEN3B	122190A	INDA	C-B2301	C-B2301	C-B2301	_	-	-	LPT1:
19 -	1	23								-		_	LPT1:
20 -	1	24	GEN3C	GEN3B	122190A	INDA	C-B2302	C-B2302	C-B2302				
21 -	1	1	GEN3C	GEN38	122190A	EVAL	C-B2303	C-B2303	C-B2303	-	_	-	LPT1:
							C-B2304	C-B2304	C-B2304	-	-	-	LPT1:
22 -	1	23	GEN3C	GEN3B	122190A	INDA							
23 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2305	C-B2305	C-B2305	-	-	-	LPT1.
				GEN3B	122190A	INDA	C-B2306	C-B2306	C-B2306	-		-	LPT1:
24 -	1	23	GEN3C									_	LPT1:
25 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2307	C-B2307	C-B2307	-	-		
26 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2308	C-B2308	C-B2308			-	LPT1:
					122190A		C-B2309	C-B2309	C-B2309	-	-		LPT1:
27 -	1	23	GEN3C	GEN3B		INDA							
28 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2310	C-B2310	C-B2310	-	-	-	LPT1:
		23	GEN3C	GEN3B	122190A	INDA	C-B2311	C-B2311	C-B2311	_	-	-	LPT1.
	1									_		-	LPT1:
30 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2312	C-B2312	C-B2312		-		
31 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2313	C-B2313	C-B2313	-	-	-	LPT1:
				GEN3B	122190A	INDA	C-B2314	C-B2314	C-B2314	-	_	_	LPT1.
32 -	1	23	GEN3C									_	LPT1.
33 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2315	C-B2315	C-82315	-	*	-	
34 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2316	C-B2316	C-B2316	-	-	-	LPTI
							C-B2317	C-B2317	C-82317	_	_	_	LPT1:
35 -	1	23	GEN3C	GEN3B	122190A	INDA					=		
36 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2318	C-B2318	C-B2318	-	-	-	LPT1:
				GEN3B	122190A	INDA	C-B2319	C-B2319	C-B2319	-	-	_	LPT1:
37 -	1	23	GEN3C									_	LPT1:
38 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2320	C-B2320	C-82320	-	_		
39 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2321	C-B2321	C-B2321	-	-	-	LPT1:
								C-B2322	C-B2322	_	_	_	LPT1:
40 -	1	23	GEN3C	GEN3B	122190A	INDA	C-82322						
41 -	1	23	GEN3C	GEN3B	1221 <del>9</del> 0A	INDA	C-B2323	C-B2323	C-82323	-	-	-	LPT1:
				GEN3B	122190A	INDA	C-B2324	C-B2324	C-B2324	_	-	-	LPT1:
42 -	1	6	GEN3C		1221304						_	_	LPT1:
43 -	1	6	GEN3C	GEN3B	122190A	INDA	C-B2325	C-B2325	C-B2325	-	-	_	
44 -	1	11	GEN3C	GEN3B	122190A	INDA	C-B2326	C-B2326	Ç-B2326	-	-	-	LPT1:
							C-B2327	C-B2327	C-B2327	_	_	-	LPT1:
45	1	1	GEN3C	GEN3B	122190A	EVAL							
46 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2328	C-B2328	C-B2328	-	-	-	LPT1:
	1	23	GEN3C	GEN3B	122190A	INDA	C-B2329	C-B2329	C-B2329	_	-	-	LPT1:
							C-B2330	C-B2330	C-B2330	_	_	_	LPT1:
48 -	1	23	GEN3C	GEN3B	122190A	INDA							
49 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2331	C-B2331	C-B2331	-	-	-	LPT1:
			GEN3C	GEN3B	122190A	INDA	C-B2332	C-B2332	C-82332	-	-	_	LPT1:
50 <del>-</del>	1	23							C-B2333	_		_	LPT1:
51 <b>-</b>	1	23	GEN3C	GEN3B	122190A	INDA	C-B2333	С-В2333					
52 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2334	C-B2334	C-B2334	-	-	-	LPT1:
							C-B2335	C-82335	C-B2335	-	_	_	LPT1:
53 -	1	23	GEN3C	GEN3B	122190A	INDA							
54 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2336	C-B2336	C-B2336	-	-	-	LPT1
					122190A	INDA	C-B2337	C-B2337	C-B2337	-	-	-	LPT1
55 -	1	23	GEN3C	GEN3B						_	_	_	LPT1
56 -	1	23	GEN3C	GEN38	122190A	INDA	C-B2338	C-B2338	C-B2338	-	_	-	
57 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2339	C-B2339	C-B2339	-	-	-	LPT1:
							C-B2340	C-B2340	C-82340	_	-	-	LPT1 ·
58 <b>-</b>	1	23	GEN3C	GEN3B	122190A	INDA						_	
59 -	1	23	GEN3C	GEN3B	122190A	INDA	C-B2341	C-B2341	C-B2341	_	-	-	LPT1
60 -	ī	23	GEN3C	GEN3B	122190A	INDA	C-B2342	C-B2342	C-B2342	-	-	-	LPT1:
							C-B2343	C-B2343	C-B2343	-	_	-	LPT1 ·
61 -	1	23	GEN3C	GEN3B	122190A	INDA					_	_	
STE? Pi	++10	2h23	~⊶¢€изс	GEN3B	122190A	INDA	C-B2344	C-B2344	C-B2344		_	_	¥206
CITE ET		لمابدر	- 911										J_ 0 0

													,		
63	-	1	23	GEN3C	GEN3B	122190A	INDA	C-B2345	C-B2345	C-B2345	_	_		J)P <b>_</b> 1	
64	-	1	23	GEN3C	GEN 3B	122190A	INDA	C-B2346	C-B2346	C-B2346	-	_		LPT1:	
65	-	ı	23	GEN3C	GEN3B	122190A	INDA	C-B2347	C-B2347	C-B2347	_	-	_	LPT1:	
66	-	1	6	GEN3C	GEN3B	122190A	INDA	C-B2348	C-B2348	C-B2348	-	-	-	r.prr.	
67	-	1	11	GEN3C	GEN3B	122190A	INDA	C-B2349	C-B2349	C-B2349	-	_	_	LPT1.	S
68	-	1	1	GEN3C	GEN3B	122190A	EVAL	C-B2350	C-B2350	C-B2350	_	_	_	LPT1:	G
69	-	1	23	GEN3C	GEN3B	122190A	INDA	C-B2351	C-B2351	C-B2351	-	-	_	LPT1:	~
70	-	1	23	GEN3C	GEN3B	122190A	INDA	C-B2352	C~B2352	C-B2352	-	-	-	LPT1:	
71	-	1	23	GEN3C	GEN3B	122190A	INDA	C-B2353	C-B2353	C-B2353	-	-	_		Ċ
72	-	1	23	GEN3C	GEN3B	122190A	INDA	C-B2354	C-B2354	C~B2354	-	-	-	LPT1.	
73	-	1	23	GEN3C	GEN3B	122190A	INDA	C-B2355	C-B2355	C-B2355	_	-	_		-3
74	-	1	23	GEN3C	GEN3B	122190A	INDA	C-82356	C-B2356	C-B2356	_	-	_	LPT1.	~
75	-	1	23	GEN3C	GEN3B	122190A	INDA	C-82357	C-B2357	C-B2357	-	-	-	LPT1:	
76	-	1	6	GEN3C	GEN3B	122190A	INDA	C-B2358	C-B2358	C-B2358	_	-	-	LPT1:	
77	-	1	11	GEN3C	GEN3B	122190A	INDA	C-B2359	C-B2359	C-B2359	-	~	-	LPT1:	

Turbochrom Sequence File : H:\ACQUIRE\MET\_SEQ\2190-E.SEQ

Created by : DE11/02/98 on : 5/19/00 10:00 Edited by : DE05/19/00 on : 5/19/00 10:07

Description : QUANTERRA PGH 8081 RUN ON GC#3 CLP1/CLP2 CQLUMNS

REVIEWED BY:

Number of Times Edited: 1

## Sequence File Header Information:

Number of Rows : 15

Instrument Type : 760 / 900 Series Intelligent Interface
Injection Type : SINGLE

				Sequence Samp	ole Descri	ptions -	Channel B	1				
Row	Type	Sample	Sample	Study Name	Sample	ISTD	Sample	Díl.	Mult	Divisor	Addend	Norm.
	-76-	Name	Number	-	Amount	Amount	Volume	Factor				factor
									1 000		0.000	100.000
1	Std Check	EVALB, 2170-E.b,	190-88 <b>-</b> 8		1.000	1.000	1.000	1.000	1.000			
2	Std Check	MEDTOX, 2170-E b	190-84-13		1.000	1.000	1.000	1.000	333.000	1 000	0.000	100 000
3	Std Check	MEDCHLOR, 2170-E	190-85-10	FAILS 1	1.000	1.000	1.000	1 000	333.000	1.000	0.000	100.000
4	Std Check	MEDA. 2170-E b	190-84-3		1.000	1.000	1.000	1 000	1.000	1.000	0.000	100.000
5	Std Check	MEDB, 2170-E b,	190-84-9		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
6	_	DD6A4103,2170-E	130194001		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
<u> </u>		DD6A411A, 2170-E	1301940018	:	1,000	1.000	1.000	1.000	1.000	1.000	0 000	100 000
,	Sample	DD6A411C,2170-E	1301940010		1.000	1.000	1.000	1.000	1.000	1.000	0.000	100.000
8	Sample		130194001	,	1 000	1.000	1.000	1,000	1.000	1.000	0.000	100 000
9	Sample	DD6A5103,2170-E					1.000	1.000	1.000	1.000	0.000	100 000
10	Sample	DD6A6103,2170-E	130194003		1.000	1 000					0.000	100.000
11	Sample	DD6A7103,2170-E	130192004		1.000	1 000	1.000	1.000	1.000	1.000		
12	Sample	DDE2E101,2170-E	130192BLK		1.000	1.000	1.000	1 000	1.000	1.000	0.000	100.000
13	Sample	DDE2E102,2170-E	130192LCS		1.000	1.000	1.000	1.000	1 000	1.000	0.000	100.000
14	Std Check	MEDA, 2170-E b,,	190-84-3		1.000	1.000	1.000	1.000	1.000	1.000	0 000	100.000
15	Std Check	MEDB, 2170-E.b.,	190-84-9		1 000	1 000	1.000	1.000	1.000	1.000	0.000	100.000

Row	Site	Rack	Vıal	Inst Method	Process Method	Se Calib Method	equence Pr Report Format	ocess Info Raw File	rmation - ( Result File	Channel B Baseline File	Modified Raw File	Cal Rpt		Update RT	Out Dev
1				GEN3C	GEN3B	122190A	EVAL	C-B2361	С-в2361		C-B2361	-	-	-	LPT1:
2	_	î	2	GEN3C	GEN3B	122190A	TOX	C-B2362	C-B2362		C-B2362	-	-	-	LPT1 . LP
3	_	3	2	GEN3C	GEN3B	122190A	TOX	C-B2363	C-B2363		C-B2363	-	-	-	LPT1.,LP
4	-	ī	6	GEN3C	GEN3B	122190A	INDA	C-B2364	C-B2364		C-B2364	-	-	-	LPT1:
٠,	-	1	11	GEN3C	GEN3B	122190A	INDA	C-B2365	C-B2365		C-B2365	-	-	_	LPT1.
6		i	23	GEN3C	GEN3B	122190A	INDA	C-B2366	C-B2366		C-B2366	-	-	_	LPT1 ·
7	_	1	23	GEN3C	GEN3B	122190A	INDA	C-B2367	C-B2367		C-B2367	-	-	-	LPT1.
9	_	ī	23	GEN3C	GEN3B	122190A	INDA	C-B2368	C-B2368		C-B2368	-	-	-	LPT1
9		ī	23	GEN3C	GEN3B	122190A	INDA	C-B2369	C-B2369		C-B2369	-	-	-	LPT1
10		ī	23	GEN3C	GEN3B	122190A	INDA	C-B2370	C-B2370		C-B2370	-	-	_	LPT1:
11	_	ī	23	GEN3C	GEN3B	122190A	INDA	C-B2371	C-B2371		C-B2371	-	_	-	LPT1
12		î	23	GEN3C	GEN3B	122190A	INDA	C-B2372	C-B2372		C-B2372	-	-	-	LPT1:
13	_	1	23	GEN3C	GEN3B	122190A	INDA	C-B2373	C-B2373		C-B2373	-	-	-	LPT1:
14	-	1	6	GEN3C	GEN3B	122190A	INDA	C-B2374	C-B2374		C-B2374	_	-	-	LPT1.
15		1	11	GEN3C	GEN3B	122190A	INDA	C-B2375	C-B2375		C-B2375	-	-	_	LPT1:

PSRO24 5/18/00 11:43:53 MT SAMPLE CUSTODIAN REMOVAL REQUEST

PAGE 001

REQUESTED BY: PINOB

METHOD: QJ Pesticides (8081A)

STORAGE LOCATION	N WORK ORDER #	PICKED CNTR#	CONTROL #	CLIENT #	ANALYSIS	LOTID	SMP#	<u>sfx</u>	MATRIX DESCRIPTION	QTY <u>RCV</u>	' QT <u>'D RE</u>	
2F CLP1 50	DD6A4-1-03		234746	399411	A-36-QJ	COE130194	001	R C SOLID		0	3	1
2F CLP1	DD6A5-1-03	<del></del>	234747	399411	A-36-QJ	COE130194	002	SOLID		0	3	1
2F CLP1	DD6A6-1-03		234748	399411	A-36-QJ	COE130194	003	SOLID		0	3	1
2F CLP1	DD6A7-1-03		234749	399411	A-36-QJ	COE130194	004	SOLID		0	3	1

Brian a. Prio	Bian a. Priso	5-18-00 1500 5-18-00 1800

## HERBICIDE DATA

HERBICIDE QC SUMMARY

SW846 8151A SURROGATE RECOVERY

657 532

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

QESSDG:

Lot #: C0E130194

CLIENT ID.	SRG01	TOT OUT
)=====================================	=   ======	======
01 DF/S1/0133/SDC/001 A	51	00
02 DF/S1/0133/SDC/001 B	57	00
03 DF/S1/0133/SDC/001_C	56	<u>00</u>
04 DF/S1/0133/SDC/001 D	66	00
05 METHOD BLK. DDE4W101	78	00
06 LCS DDE4W102	89	00
07 DF/S1/0133/SDC/001 A D	86	00
08 DF/S1/0133/SDC/001 A S	82	00

SURROGATES SRG01 = DCAA QC LIMITS ( 42-125)

<sup>#</sup> Column to be used to flag recovery values

<sup>\*</sup> Values outside of required QC Limits

D System monitoring Compound diluted out

## SW846 8151A CHECK SAMPLE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Lot #: C0E180000

WO #: DDE4W102 BATCH: 0139489

T'	SPIKE	SAMPLE		QC	[
	ADDED	CONCENT.	ક	LIMITS	1
COMPOUND	(mg/L)	(mg/L)	REC	REC	QUAL
		=======================================	====		
2,4-D	0.160	0.133	83	28- 136	Ì
2,4,5-TP (Silvex)	0.0400	0.0323	81	50- 128	

* Values outside	of QC limits		
Spike Recovery:	0 out of	2 outside limits	ı
COMMENTS:			

NOTES (S):

657 534

SW846 8151A MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A

Lot #: C0E130194

WO #: DD6A411D BATCH: 0139489

	SPIKE ADDED	SAMPLE CONCENT.	MS CONCENT.	MS %	LIMITS	
COMPOUND	(mg/L )	(mg/L )	(mg/L)	REC	REC	LAUQ
=======================================			=======			=======
2,4-D	0.160	ND	0.133	83_	<u> 35- 133</u>	l1
2,4,5-TP (Silvex)	0.0400	ND	0.0322	80	50- 131	

N	$\mathbf{O}$	'ES	1	21	
- 13	~1	. DJ			

# Column to be used to flag recovery and RPD values with an asterisk * Values outside of QC limits
RPD: 0 out of 0 outside limits Spike Recovery: 0 out of 2 outside limits
COMMENTS:

#### SW846 8151A MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: Severn Trent Laboratories, Inc. Client: UXB INTERNATIONAL

Lab Code: QESPIT

SDG No:

Matrix Spike ID: DF/S1/0133/SDC/001 A

Lot #: C0E130194

WO #: DD6A411E BATCH: 0139489

     COMPOUND	SPIKE ADDED (mg/L )	MSD CONCENT. (mg/L )	MSD % REC	% RPD	QC RPD	LIMITS REC	QUAL
77	0.160 0.0400	0.144 0.0339	90	8.4 5.2	20 20	- <u>!</u>	=====================================

NOTES (S):			

RPD: 0 out of 2 outside limits Spike Recovery: \_\_\_0 out of \_\_\_2 outside limits

COMMENTS:

<sup>#</sup> Column to be used to flag recovery and RPD values with an asterisk

<sup>\*</sup> Values outside of QC limits

#### SW846 8151A METHOD BLANK SUMMARY

BLANK WORKORDER NO.
DDE4W101

Lab Name: Severn Trent Laboratories, Inc.

Lab Code: QESPIT SDG Number:

Lab File ID: a-a20536. Lot Number: COE130194

Matrix: SOLID WATER Extraction Method:

Un 5 boloo Date Extracted: 05/18/00

Date Analyzed(1): 05/20/00 Date Analyzed(2): N/A

Time Analyzed(1): 14:27 Time Analyzed(2): N/A

Instrument ID(1): A/B Instrument ID(2): N/A

GC Column(1): DB5/DB1701 ID: 053 GC Column(2): N/A ID: N/A

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS, AND MSD:

Ī		SAMPLE	DATE	DATE
į	CLIENT ID.	WORK ORDER #	ANALYZED(1)	ANALYZED(2)
1		=======================================	=======================================	=========
01	CHECK SAMPLE	DDE4W102 C	05/20/00	N/A
02	DF/S1/0133/SDC/001 A	DD6A4104	05/20/00	<u>N/A</u>
03	DF/S1/0133/SDC/001 A	DD6A411D S	05/20/00	_N/A
04	DF/S1/0133/SDC/001 A	DD6A411E D	05/20/00	N/A
05	DF/S1/0133/SDC/001 B	DD6A5104	05/20/00	N/A
06	DF/S1/0133/SDC/001 C	DD6A6104	05/20/00	_N/A
07	DF/S1/0133/SDC/001 D	DD6A7104	05/20/00	N/A
08				
09				i
10				<u></u>
11				l <u></u> _
12				
13			1	<u> </u>
14			<u> </u>	<u> </u>
15				
16		l <u> </u>		
17				
18				ll
19			1	<u></u>
20	i	1	1	

COMMENTS:		
	 FORM IV	

# HERBICIDE SAMPLE DATA

657 538

#### UXB INTERNATIONAL

SDG Number: Lab Name: Severn Trent Laboratories, Inc.

Matrix: (soil/water) SOLID WATEL

Method: SW846 8151A un skalos Herbicides (8151A)

Sample WT/Vol: 100 / mL Work Order: DD6A4104 Dilution factor: 1

Moisture %:10

Date Received: 05/13/00 Date Extracted: 05/18/00 Date Analyzed: 05/20/00

Lab Sample ID: C0E130194 001

QC Batch: 0139489

Client Sample Id: DF/S1/0133/SDC/001 A

#### CONCENTRATION UNITS:

	CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L Q	
Ī	94-75-7	2,4-D	0.040	<u>u</u>
j	93-72-1	2,4,5-TP (Silvex)	0.010	U

Data File: /var/chem/gc1.i/2180.b/a-a20530.d

Report Date: 22-May-2000 08:54

# STL-PITTSBURGH

Data file : /var/chem/gc1.i/2180.b/a-a20530.d

Lab Smp Id: DD6A4104 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 20-MAY-2000 11:33 Operator : 01797 Smp Info : DD6A4104,2180.b

Inst ID: gcl.i

Misc Info : 130194001

Comment

Method : /var/chem/gcl.i/2180.b/LONGH.m

Meth Date: 22-May-2000 08:08 g Quant Type: ESTD Cal Date : 18-MAY-2000 12:08 Als bottle: 3 Dil Factor: 1.00000 Cal File: a-a20468.d

Integrator: Falcon Compound Sublist: all.sub

Target Version: 3.40

## Concentration Formula: Amt \* DF \* 20\*Vt/Vo/Vi

Name	Value	Description
DF	1.000	Dilution Factor
Vt	10.000	Volume of final extract (uL)
Vo		Volume of sample extracted (mL)
Vi		Volume injected

		CONCENTRATIONS
		ON-COLUMN FINAL
Compounds	RT EXP RT DLT RT RESPONSE	( ng) ( mg/L)
<b>以中华的大学的国际企业企业企业企业的企业</b>	医名 医多色素素 医自己性红 拉拉拉克拉拉斯	***
1 DALAPON	Compound Not Detected.	-
\$ 2 DCAA	10.600 10.594 0.006 7760966	0 02558 0 05116(a)
3 MCPP	Compound Not Detected.	
4 DICAMBA	Compound Not Detected.	
5 MCPA	Compound Not Detected.	
6 DICHLOROPROP	Compound Not Detected.	
7 2,4-D	Compound Not Detected.	
8 PENTACHLOROPHENOL	Compound Not Detected.	
9 2,4,5-TP(SILVEX)	Compound Not Detected.	
10 2,4,5-T	Compound Not Detected.	
11 DINOSEB	Compound Not Detected.	
12 2,4-DB	18 551 18 579 -0.028 52472	0 00133 0.002662(a)

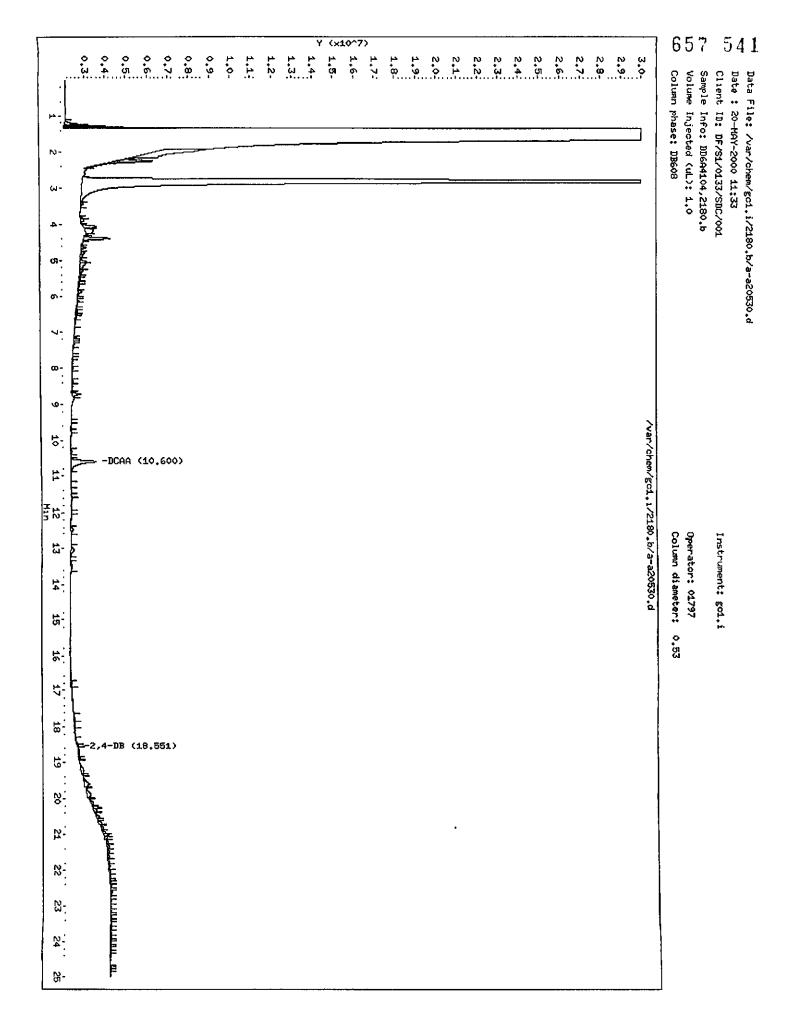
STL Pittsburgh 4010

657 539

Data File: /var/chem/gc1.i/2180.b/a-a20530.d Report Date: 22-May-2000 08:54

QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).



### UXB INTERNATIONAL

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) COLID WATEL Lab Sample ID:COE130194 002

Method: SW846 8151A um 5/22/00

Herbicides (8151A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A5104 Date Extracted:05/18/00 Dulution factor: 1 Date Analyzed: 05/20/00

Moisture %:8.6

QC Batch: 0139489

Client Sample Id: DF/S1/0133/SDC/001 B

# CONCENTRATION UNITS:

	CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L	Q
_	94-75-7	2,4-D	0.040	<u>  U                                   </u>
	93-72-1	2,4,5-TP (Silvex)	0.010	<u>                                     </u>

Data File: /var/chem/gc1.i/2180.b/a-a20533.d

Report Date: 22-May-2000 08:55

# STL-PITTSBURGH

CONCENTRATIONS ON 007 1701 77177

Data file : /var/chem/gc1.i/2180.b/a-a20533.d Lab Smp Id: DD6A5104 Clier Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 20-MAY-2000 13:00

Inst ID: gcl.i Operator: 01797

Smp Info : DD6A5104,2180.b

Misc Info: 130194002

Comment

Method : /var/chem/gcl.i/2180.b/LONGH.m Meth Date : 22-May-2000 08:08 g Qu Cal Date : 18-MAY-2000 12:08 Cal Quant Type: ESTD Cal File: a-a20468.d

Als bottle: 6

Dil Factor: 1.00000

Compound Sublist: all.sub Integrator: Falcon

Target Version: 3.40

Concentration Formula: Amt \* DF \* 20\*Vt/Vo/Vi

Name	Value	Description
DF Vt Vo Vi	10.000 100.000	Dilution Factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected

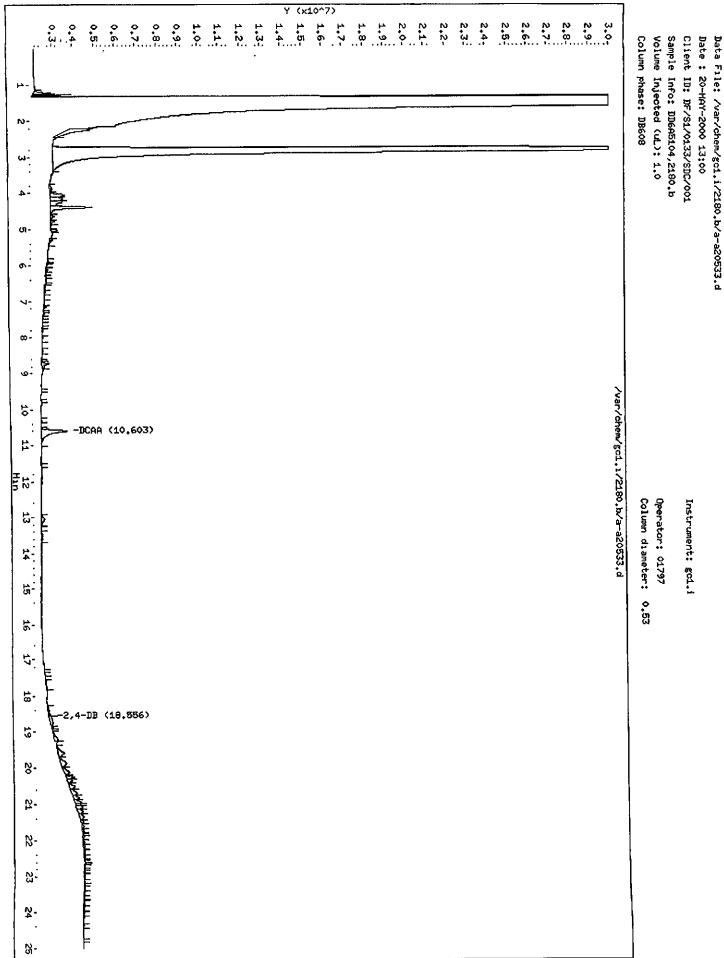
				ON-COTOWN	FINAL
Compounds	RT	EXP RT DLT R	T RESPONSE	( ng)	( mg/L)
*****************	==	=#### DDE38			******
1 DALAPON	Con	pound Not Det	ected.		
\$ 2 DCAA	10 603	10.594 0.0	09 8702403	0.02868	0 05737(a)

3	МСРР	Compound	Not	Detected.
4	DICAMBA	Compound	Not	Detected.
5	MCPA	Compound	Not	Detected.
6	DICHLOROPROP	Compound	Not	Detected.
7	2,4-D	Compound	Not	Detected.
8	PENTACHLOROPHENOL	Compound	Not	Detected.
9	2,4,5-TP(SILVEX)	Compound	Not	Detected
10	2,4,5-T	Compound	Not	Detected
11	DINOSEB	Compound	Not	Detected.

12 2,4-DB 18 556 18 579 -0 023 108784 0.00276 0 005519(a)

657 544 Data File: /var/chem/gc1.i/2180.b/a-a20533.d Report Date: 22-May-2000 08:55

# QC Flag Legend



### UXB INTERNATIONAL

SDG Number: Lab Name: Severn Trent Laboratories, Inc.

Matrix: (soil/water) SOLID WATER Lab Sample ID: C0E130194 003

Method: SW846 8151A um 5/00/00 Herbicides (8151A)

Date Received: 05/13/00 Sample WT/Vol: 100 / mL Date Extracted:05/18/00 Work Order: DD6A6104 Date Analyzed: 05/20/00 Dilution factor: 1

Moisture %:10

QC Batch: 0139489

Client Sample Id: DF/S1/0133/SDC/001 C

# CONCENTRATION UNITS:

	CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L (	3
Ī	94-75-7	2,4-D	0.040	<u> </u>
i	93-72-1	2,4,5-TP_(Silvex)	0.010	U

Data File: /var/chem/gc1.i/2180.b/a-a20534.d

Report Date: 22-May-2000 08:55

#### STL-PITTSBURGH

Data file : /var/chem/gc1.i/2180.b/a-a20534.d

Lab Smp Id: DD6A6104 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 20-MAY-2000 13:29

Operator: 01797 Inst ID: gcl.i

Smp Info : DD6A6104,2180.b

Misc Info : 130194003

Comment

12 2,4-DB

Method : /var/chem/gcl.i/2180.b/LONGH.m

Meth Date: 22-May-2000 08:08 g Quant Type: ESTD Cal Date: 18-MAY-2000 12:08 Cal File: a-a20468.d

Als bottle: 7

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: all.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* 20\*Vt/Vo/Vi

Name	Value	Description
DF	1.000	Dilution Factor
Vt		Volume of final extract (uL)
Vo	100.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected

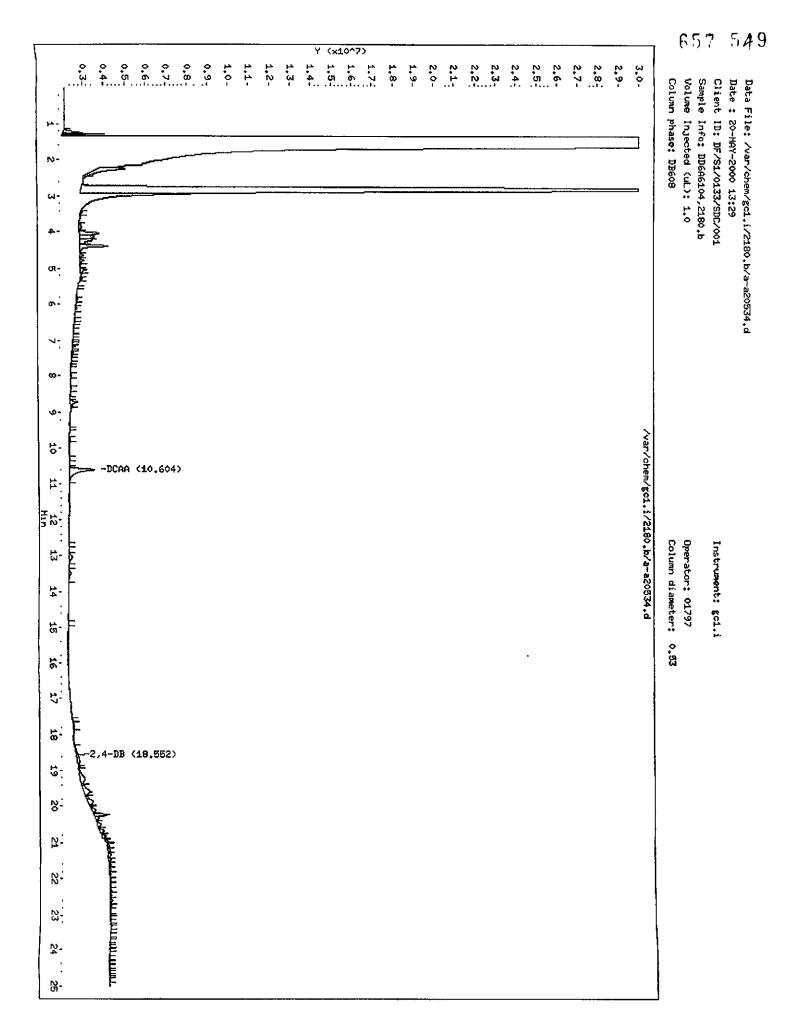
#### CONCENTRATIONS ON-COLUMN FINAL EXP RT DLT RT RESPONSE (ng) ( mg/L) Compounds \*\*\*\* 222222 -----1 DALAPON Compound Not Detected 0 02806 0.05612(a) 10 604 10 594 0 010 8513361 2 DCAA 3 MCPP Compound Not Detected. 4 DICAMBA Compound Not Detected. Compound Not Detected. 5 MCPA 6 DICHLOROPROP Compound Not Detected Compound Not Detected. 7 2,4-D Compound Not Detected 8 PENTACHLOROPHENOL 9 2,4,5-TP(SILVEX) Compound Not Detected. 10 2,4,5-T Compound Not Detected. 11 DINOSEB Compound Not Detected

STL Pittsburgh 4018

18.552 18.579 -0.027 108134 0.00274 0.005486(a)

Data File: /var/chem/gc1.i/2180.b/a-a20534.d Report Date: 22-May-2000 08:55

# QC Flag Legend



# UXB INTERNATIONAL

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATER Lab Sample ID:COE130194 004

Herbicides (8151A) LMS/22/00 Method: SW846 8151A

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Date Extracted: 05/18/00 Work Order: DD6A7104 Date Analyzed: 05/20/00 Dilution factor: 1

Moisture %:11

QC Batch: 0139489

Client Sample Id: DF/S1/0133/SDC/001 D

# CONCENTRATION UNITS:

	CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L	<u>}</u>
1	94-75-7	2,4-D	0.040	ט
i	93-72-1	2,4,5-TP (Silvex)	0.010	U

Data File: /var/chem/gc1.i/2180.b/a-a20535.d

Report Date: 22-May-2000 08:55

# STL-PITTSBURGH

Data file : /var/chem/gc1.i/2180.b/a-a20535.d

Lab Smp Id: DD6A7104 Client Smp ID: DF/S1/0133/SDC/001

Inj Date : 20-MAY-2000 13:58 Operator : 01797 Smp Info : DD6A7104,2180.b Inst ID: qc1.i

Misc Info: 130194004

Comment

Method : /var/chem/gc1.i/2180.b/LONGH.m

Meth Date : 22-May-2000 08:08 g Quant Type: ESTD

Cal Date : 18-MAY-2000 12:08 Cal File: a-a2046

Als bottle: 8 Cal File: a-a20468.d

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: all.sub

Target Version: 3.40

# Concentration Formula: Amt \* DF \* 20\*Vt/Vo/Vi

Name	Value	Description
DF	1 000	Dilubias Darks
		Dilution Factor
Vt		Volume of final extract (uL)
Vo	100.000	Volume of sample extracted (mL)
Vi	1.000	Volume injected

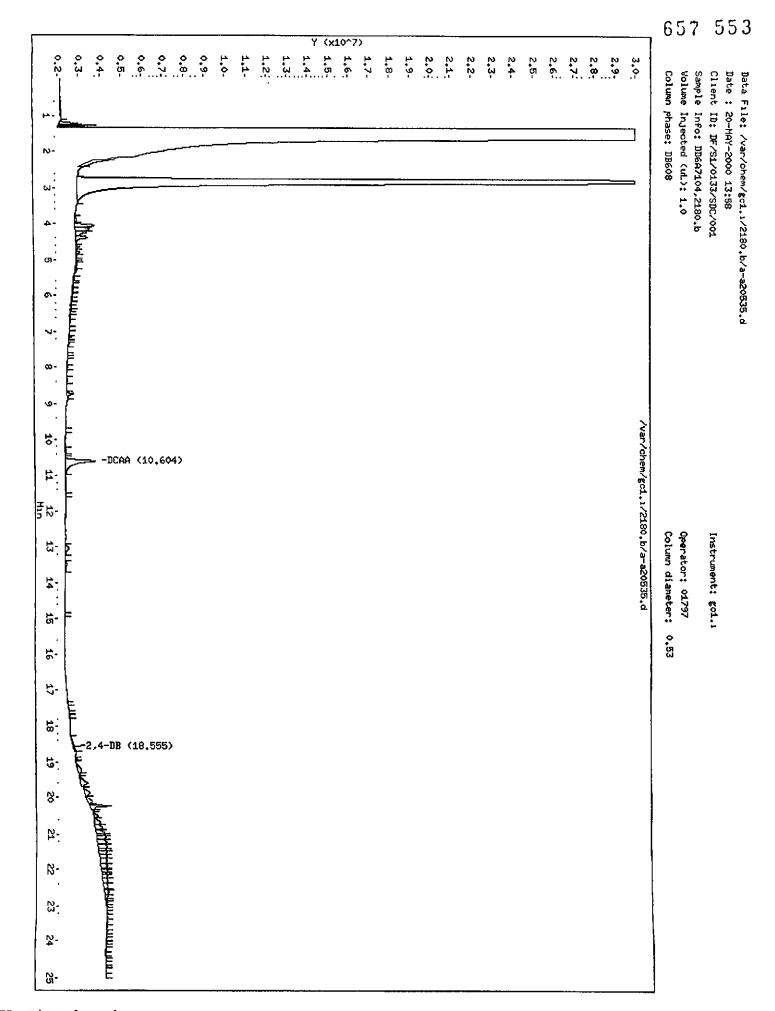
# CONCENTRATIONS

						ON-COLUMN	FINAL
Compo	ounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	(mg/L)
	化过滤性 化铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁铁				*******		******
:	DALAPON	Cor	mpound No	t Detecte	d.		
\$ :	2 DCAA	10.604	10.594	0.010	9970107	0.03286 /	0.06572(a)
;	3 MCPP	Cor	pound No	t Detecte	d	`	
•	1 DICAMBA	Cor	npound No	t Detecte	đ		
!	5 MCPA	Cor	pound No	t Detecte	:đ		
	5 DICHLOROPROP	Cor	mpound No	t Detecte	d.		
	7 2,4-D	Cor	mpound No	t Detecte	d.		
	9 PENTACHLOROPHENOL	Cor	mpound No	t Detecte	d.		
,	9 2,4,5-TP(SILVEX)	Cor	npound No	t Datecte	:d		
10	D 2,4,5-T	Cor	mpound No	t Detecte	eđ.		
1	1 DINOSEB	Cor	mpound No	t Detecte	d.		
1	2 2,4-DB	18.555	18.579	-0 024	77324	0.00196	0.003923(a)

Data File: /var/chem/gc1.i/2180.b/a-a20535.d Report Date: 22-May-2000 08:55

QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).



# HERBICIDE CALIBRATION DATA

Report Date : 20-May-2000 11:44



# STL-PITTSBURGH

# COMPOUND LISTING

Method file

: /var/chem/gcl.i/2180.b/LONGH.m : ESTD Quant Method Target Version : 3.40 Number of Cpnds : 12 Last Update : 20-May-2000 11:44

: GC MULTI COMP Data Type

Global Integrator : Falcon

Values Chromat Events \_\_\_\_ 3608.000000 Initial:Start Threshold 1804.000000 Initial: End Threshold Initial:Area Threshold 36080.000000 1.000000 Initial:P-P Resolution 10.000000 Initial:Bunch Factor ON Initial:Negative Peaks

0.200000 Initial: Tension

Compound	RT	RT Window	RF
1 DALAPON \$ 2 DCAA 3 MCPP 4 DICAMBA 5 MCPA 6 DICHLOROPROP 7 2,4-D 8 PENTACHLOROPHENOL 9 2,4,5-TP(SILVEX) 10 2,4,5-T 11 DINOSEB 12 2,4-DB	11.212 12.001 12.836 14.348 15.580 16.791 17.910	10.966-11.106 11.142-11.282 11.931-12.071 12.766-12.906 14.278-14.418 15.510-15.650 16.721-16.861 17.840-17.980 18.009-18.149	4.747e+05 2.235e+08 5.619e+05 5.838e+07 4.745e+07 5.548e+08 2.818e+08 3.299e+08 3.108e+08

Report Date : 18-May-2000 12:35



# STL-PITTSBURGH

# INITIAL CALIBRATION DATA

Start Cal Date : 18-MAY-2000 10:12 End Cal Date : 18-MAY-2000 12:08 Quant Method : ESTD

Quant Method : ESTD
Origin : Disabled
Target Version : 3.40
Integrator : Falcon

Method file : /var/chem/gc1.i/2180.b/LONGH.m Cal Date : 18-May-2000 12:35 eppinged

Curve Type : Average

Calibration File Names:

Level 1: /var/chem/gcl.i/2180.b/a-a20464.d Level 2: /var/chem/gcl.i/2180.b/a-a20465.d Level 3: /var/chem/gcl.i/2180.b/a-a20466.d Level 4: /var/chem/gcl.i/2180.b/a-a20467.d Level 5: /var/chem/gcl.i/2180.b/a-a20468.d

Compound  1 DALAPON 3 MCPP 4 DICAMBA 5 MCPA 6 DICHLOROPROP 7 2,4-D 8 PENTACHLOROPHENOL 9 2,4,5-TP(SILVEX) 10 2,4,5-T 11 DINOSEB 12 2,4-DB	0.00500   0.01000   0.02500   0.05000   0 10000
\$ 2 DCAA	348787512 331743765 305258919 281785053 249384444 303391938  13 031

4027

Data File: /var/chem/gc1.i/2180.b/a-a20529.d

Report Date: 20-May-2000 11:44

# STL-PITTSBURGH

# CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gcl.i Lab File ID: a-a20529.d

Analysis Type:

Lab Sample ID: Mherb Quant Type: ESTD

Injection Date: 20-MAY-2000 11:04

Init. Calibration Date(s): 05/18/0 05/18/0
Init. Calibration Times: 10:12 12:08

Method File: /var/chem/gc1.i/2180.b/LONGH.m

		l	1	MIN	XAM
(	COMPOUND	RRF	RFO	RRF	%D   %D
	1 DALAPON	89103950.173	87265056.	948 0 010	2.1 15 0
ş	2 DCAA	303391938.48	4 29887207	9 906 0 010	1.5  15 0
	3 MCPP	474681.118	438552	230 0.010	7.6 15.0
	4 DICAMBA	223500149.45	8 22152550	5 882 0 010	0 9  15.0
l	5 MCPA	561852.499	5  514815.	537 0 010	8.4 15.0
l	6 DICHLOROPROP	58376589.252	2 57607653.	302   0.010	1.3  15.0
i	7 2,4-D	47446236.733	45952173	913 0.010	3.1 15 0
1	8 PENTACHLOROPHENOL	554813477.97	72 55380336	33.459 0.010	0.2  15.0
ł	9 2,4,5-TP(SILVEX)	281846101.32	29 27081090	00 474   0.010	3.9  15.0
ļ	10 2,4,5-T	329934786.86	55 31931919	94 313 0 010	3.2  15.0
}	11 DINOSEB	310771875.03	71 30943598	34.252   0.010	0.4  15.0
1	12 2,4-DB	39423470.393	7 37884142	.012[0.010]	3 9 15 0
1		1			1

Data File: /var/chem/gc1.i/2180.b/a-a20550.d Report Date: 22-May-2000 08:07

### STL-PITTSBURGH

# CONTINUING CALIBRATION COMPOUNDS

Instrument ID: gcl.i

Lab File ID: a-a20550.d Analysis Type:

Lab Sample ID: Mherb

Injection Date: 20-MAY-2000 21:15

Init. Calibration Date(s): 05/18/0 05/18/0 Init. Calibration Times: 10:12 12:08 Method File: /var/chem/gcl.i/2180.b/LONGH.m

Quant Type: ESTD

		l	1		MIN	1	MAX
	COMPOUND	RRF	l	RF0	RRF	%D	<b>%</b> D
<b></b>	· 中华代表学院社会教育主题的的名词复数中华学校中央公司和中华中国	*==   ==========				-	=====
	1 DALAPON	89103950.1	73   92	052095.6	72 0 01	0  -3.3	15.0
,	2 DCAA	303391938	484   3	24467021	7 027 0.	010  -6	9j 15
	3 MCPP	474681 1	.18	455284.6	524   0.01	0 4.1	15.0
	4 DICAMBA	223500149.	458 2	4167604	7.059 0.	010  -8	.1 15
	S MCPA	561852.4	95	533000.	701 0.01	0  5.1	15 0
	6 DICHLOROPROP	58376589.2	52   62	090188.	579 0 01	0  -6.4	15.0
	7 2,4-D	47446236.7	733   49	971974.	148 0.01	0  -5.3	15 0
	8 PENTACHLOROPHENOL	554813477.	972 5	8634859	226 0.	010  -5	.7  15
	9 2,4,5-TP(SILVEX)	281846101	329 2	29072454	9.763 0.	010  -3	.2  15
	10 2,4,5-T	329934786	865 3	4572431	2 796 0.	010  -4	.8  15
	11 DINOSEB	310771875	071	32734653	5 433 0	010) -5	.3 15
	12 2,4-DB	39423470.3	397   39	9901372.	781 0 <b>0</b> 1	0  -1.2	15.0
						1	1

4029

Lab Name: STL-PITTSBURGH Contract:

Lab Code: STLPIT Case No.: SAS No.: 40325 SDG No.: C0E130194

GC Column: DB608 ID: 0.53 (mm) Init. Calib. Date(s): 05/18/00 05/18/00

Instrument ID: GC1

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS, SAMPLES, AND STANDARDS IS GIVEN BELOW:

	MEAN SURROC S1 : 10.59	SATE RT FROM 1	INITIAL CALI	BRATION		
	EPA SAMPLE NO.	LAB SAMPLE ID	DATE	TIME	S1 RT #	RT #
					π	
10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 27	DF/S1/0133/S DF/S1/0133/S DF/S1/0133/S DF/S1/0133/S DF/S1/0133/S DF/S1/0133/S	LAB SAMPLE ID ====================================	DATE ANALYZED  05/18/00 05/18/00 05/18/00 05/18/00 05/18/00 05/20/00 05/20/00 05/20/00 05/20/00 05/20/00 05/20/00 05/20/00	TIME ANALYZED  1012 1041 1110 1139 1208 1104 1133 1202 1231 1300 1329 1358 1427 1456 2115	S1 RT # ======= 10.60 10.60 10.60 10.59 10.60 10.60 10.60 10.60 10.60 10.60	RT #
28						
29						
30				<u> </u>	]	
31 32						

S1 = DCAA

QC LIMITS (+/- 0.07 MINUTES)

# Column used to flag retention time values with an asterisk.

\* Values outside of QC limits.

page 1 of 1

FORM VIII PEST

OLM03.0

Data File: /var/chem/gc1.i/2180.b/a-a20464.d

Report Date: 18-May-2000 12:35

### STL-PITTSBURGH

Data file : /var/chem/gcl.i/2180.b/a-a20464.d Lab Smp Id: Lherb

Inj Date : 18-MAY-2000 10:12

Inst ID: gcl.i Operator : 01797

Smp Info : Lherb, 2180.b Misc Info : 190-80-1

Comment

: /var/chem/gc1.i/2180.b/LONGH.m Method

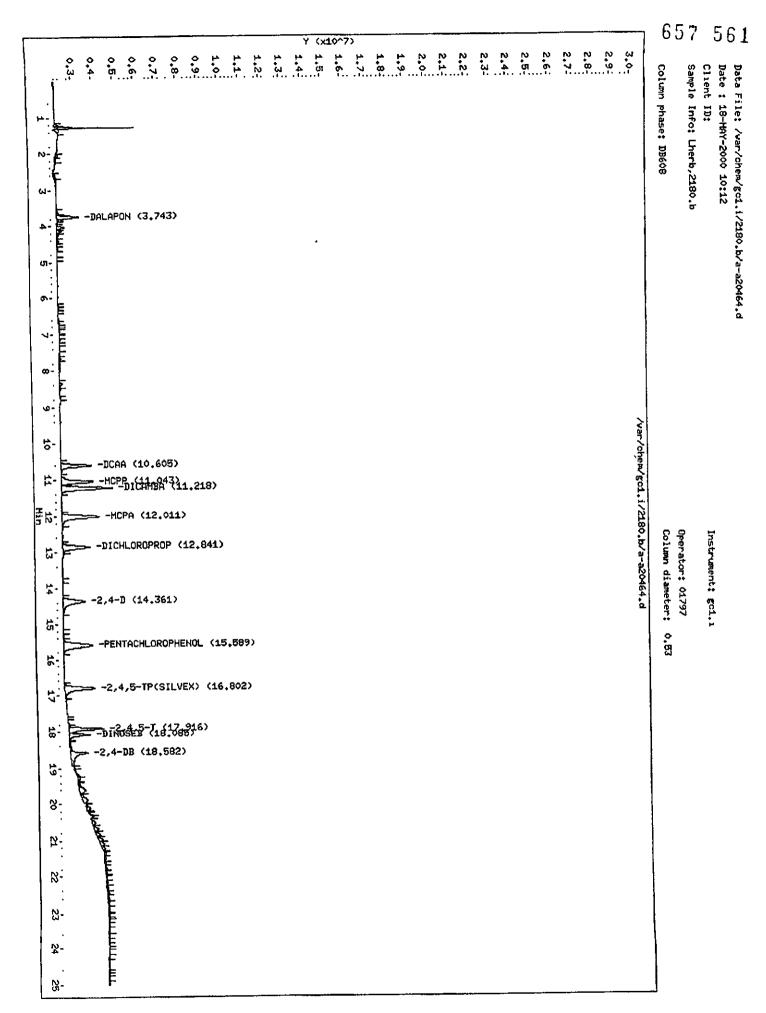
Meth Date: 18-May-2000 12:35 eppinged Quant Type: ESTD Cal File: a-a20464.d Cal Date : 18-MAY-2000 10:12

Calibration Sample, Level: 1

Als bottle: 2 Dil Factor: 1.00000 Integrator: Falcon Compound Sublist: 1-all.sub

Target Version: 3.40

						AMOUN	TS
						CAL-AMT	ON-COL
Соп	pounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
	*	==		27.0 CEE	***	******	######################################
	1 DALAPON	3.743	3.742	0.001	1072796	0 01100	0.01100
s	2 DCAA	10.605	10.600	0.005	7429174	0.02130	0.02130
·	3 MCPP	11 043	11 040	0.003	1554114	2.12000	2.120
	4 DICAMBA	11.218	11 218	0.000	2490159	0.01060	0 01060
	5 MCPA	12.011	12.006	0.005	1837675	2.14000	2.140
	6 DICHLOROPROP	12.841	12.837	0.004	1351403	0.02120	0.02120
	7 2.4-D	14.361	14.351	0.010	1075351	0 02110	0 02110
	8 PENTACHLOROPHENOL	15.589	15.586	0.003	1433994	0.00266	0.002660
	9 2.4,5-TP(SILVEX)	16.802	16 798	0.004	1466074	0.00525	0.005250
	10 2,4,5-T	17.916	17.910	0.006	1696184	0 00527	0.005270
	11 DINOSES	18.085	18.085	0.000	1022613	0 00317	0.003170
	12 2,4-DB	18 582	18.5BO	0.002	859969	0.02110	0 02110



Data File: /var/chem/gc1.i/2180.b/a-a20465.d

Report Date: 18-May-2000 12:35

#### STL-PITTSBURGH

Data file : /var/chem/gcl.i/2180.b/a-a20465.d

Lab Smp Id: MLherb

Inj Date : 18-MAY-2000 10:41
Operator : 01797 Inst ID: gcl.i

Smp Info : MLherb, 2180.b

Misc Info: 190-80-2

Comment

Method : /var/chem/gcl.i/2180.b/LONGH.m Meth Date : 18-May-2000 12:35 eppinged Qu Quant Type: ESTD Cal File: a-a20465.d

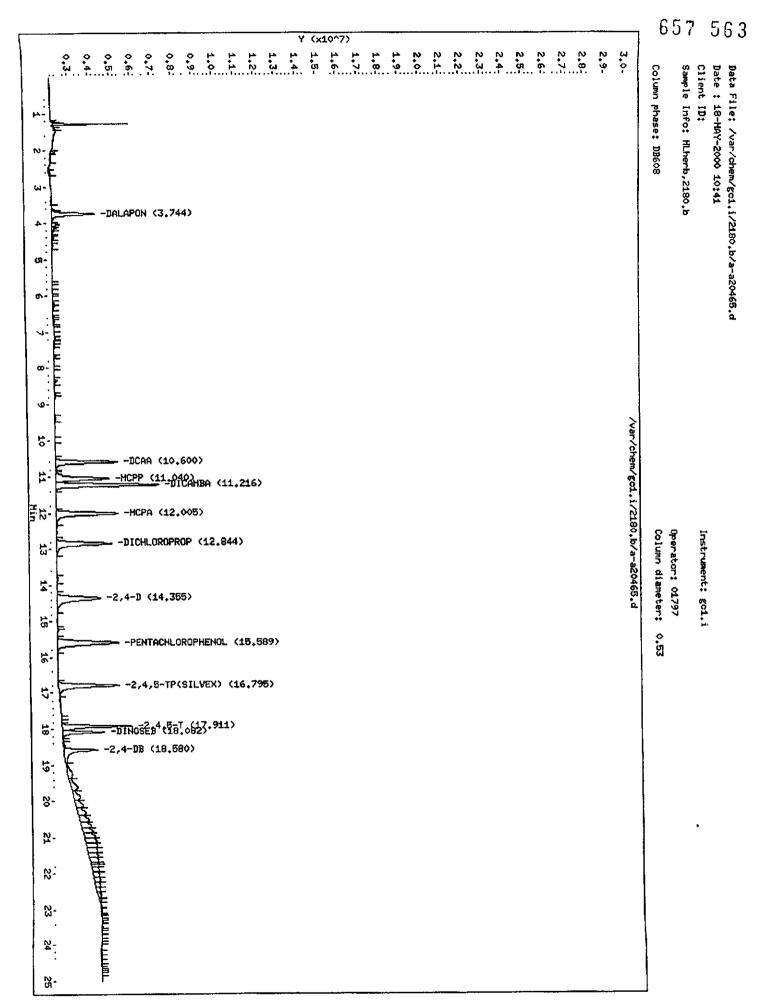
Cal Date : 18-MAY-2000 10:41 Calibration Sample, Level: 2 Als bottle: 3

Dil Factor: 1.00000 Integrator: Falcon

Compound Sublist: 1-all.sub

Target Version: 3.40

						MOUN	TTS
						CAL-AMT	ON-COL
Cor	mpounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
==		# # # # # # # # # # # # # # # # # # #	*****		디디 보 의 한 12 후 교		8 J J W W W W W
	1 DALAPON	3.744	3 742	0.002	2062640	0.02200	0.02157
\$	2 DCAA	10.600	10.600	0.000	14099110	0.04250	0 04144
•	3 MCPP	11.040	11 040	0.000	2566482	4.26000	3.843
	4 DICAMBA	11.216	11.218	-0.002	5012679	0.02130	0.02132
	5 MCPA	12 005	12.006	-0.001	3025338	4.28000	3 865
	6 DICHLOROPROP	12 844	12.837	0.007	2684133	0.04240	0.04225
	7 2,4-D	14.355	14 351	0.004	2107638	0 04250	0.04192
	8 PENTACHLOROPHENO	L 15.589	15 586	0 003	2979386	0 00532	0.005421
	9 2,4,5-TP(SILVEX)	16.795	16.798	-0.003	2928487	0.01050	0.01049
	10 2,4,5-T	17.911	17.910	0 001	3423539	0.01050	0.01057
	11 DINOSEB	18.082	18.085	-0 003	2041441	0.00635	0 006339
	12 2,4-DB	18 580	18.580	0 000	1633271	0.04220	0 04111



Data File: /var/chem/gcl.i/2180.b/a-a20466.d

Report Date: 18-May-2000 12:35

# STL-PITTSBURGH

Data file : /var/chem/gc1.i/2180.b/a-a20466.d Lab Smp Id: Mherb Inj Date : 18-MAY-2000 11:10

Inst ID: gcl.i Operator : 01797

Smp Info : Mherb, 2180.b

Misc Info: 190-80-3

Comment

Method : /var/chem/gc1.i/2180.b/LONGH.m Meth Date : 18-May-2000 12:35 eppinged Qu Quant Type: ESTD Cal File: a-a20466.d Cal Date : 18-MAY-2000 11:10

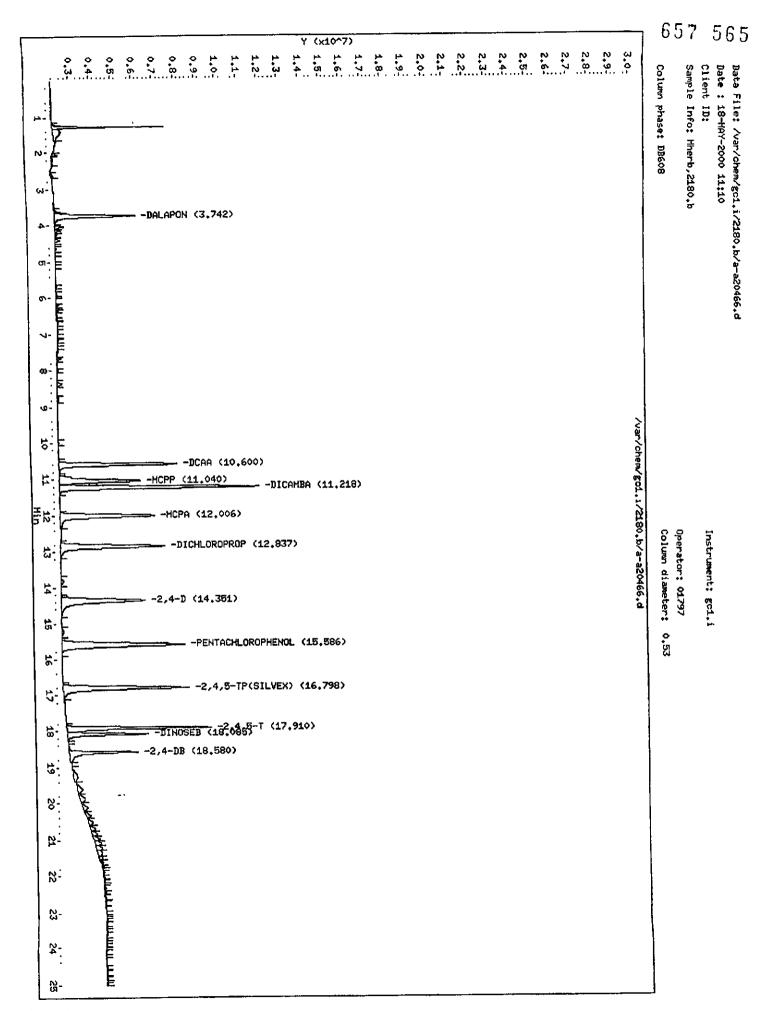
Calibration Sample, Level: 3 Als bottle: 4

Dil Factor: 1.00000

Compound Sublist: 1-all.sub Integrator: Falcon

Target Version: 3.40

					AMOUN	TS
					CAL-AMT	ON-COL
Соп	ebnuoqu	RT	EXP RT DLT RT	RESPONSE	( ng)	( ng)
	13.22.22.22.23.23.23.23.23.23.23.23.23.23	**	医圆性性畸胎 医动物性动脉	22225427	<b>医系数性数据</b>	
	1 DALAPON	3 742	3.742 0.00	0 3907001	0.04390	0.04182
Ś	2 DCAA	10 600	10.600 0.00	0 25977534	0.08510	0.07906
•	3 MCPP	11.040	11.040 0 00	0 3911550	8 52000	6.539
	4 DICAMBA	11.218	11.218 0.00	0 9630896	0.04250	0.04146
	5 MCPA	12.006	12.006 0.00	0 4590884	8 56000	6.552
	6 DICHLOROPROP	12.837	12.837 0.00	0 5048056	0 08480	0.08117
	7 2,4-D	14 351	14.351 0.00	0 4079345	0.08510	0.08242
	8 PENTACHLOROPHENOL	15 586	15.586 0.00	0 5970947	0.01064	0.01079
	9 2,4,5-TP(SILVEX)	16.798	16.798 0.00	0 6114466	0.02110	0.02163
	10 2,4,5-T	17 910	17 910 0.00	0 7000461	0 02110	0.02144
	11 DINOSEB	18 085	18 085 0.00	0 3966951	0.01270	0.01244
	12 2,4-DB	18 580	18.580 0.00	0 3355256	0.08450	0 08447



Data File: /var/chem/gc1.i/2180.b/a-a20467.d

Report Date: 18-May-2000 12:35

# STL-PITTSBURGH

Data file : /var/chem/gc1.i/2180.b/a-a20467.d Lab Smp Id: MHherb

Inj Date : 18-MAY-2000 11:39

Inst ID: gcl.i Operator : 01797

Smp Info : MHherb, 2180.b

Misc Info: 190-80-4

Comment

Method : /var/chem/gcl.i/2180.b/LONGH.m Meth Date : 18-May-2000 12:35 eppinged Qu Ouant Type: ESTD Cal File: a-a20467.d

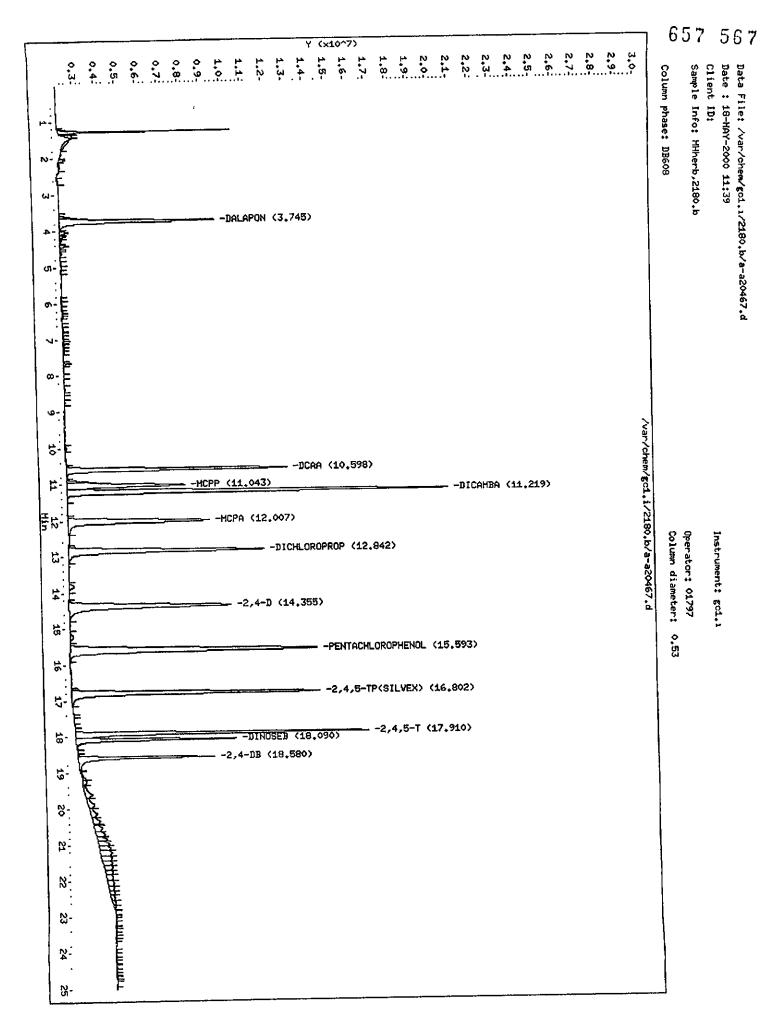
Cal Date : 18-MAY-2000 11:39 Calibration Sample, Level: 4

Als bottle: 5 Dil Factor: 1.00000

Compound Sublist: 1-all.sub Integrator: Falcon

Target Version: 3.40

						AMOUN	TS
						CAL-AMT	ON-COL
Co	mpounds .	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
		==	=====	#3##3#			
	1 DALAPON	3 745	3 742	0.003	7455549	0.08780	0.08166
\$	2 DCAA	10.598	10.600	-0.002	47903459	0.17000	0.1512
Þ	3 MCPP	11.043	11.040	0.003	5719515	17.0000	10 74
	4 DICAMBA	11 219	11 218	0.001	18486872	0.08510	0.08090
	S MCPA	12.007	12.006	0.001	6842481	17.1000	10.94
	6 DICHLOROPROP	12.842	12.837	0.005	9453587	0.17000	0 1561
		14.355	14.351	0.004	7843431	0.17000	0.1612
	7 2,4-D 8 PENTACHLOROPHENOL	15.593	15.586	0.007	11981629	0.02128	0.02156
	9 2,4,5-TP(SILVEX)	16.802	16.798	0.004	12016265	0.04210	0.04241
		17.910	17.910	0.000	14229900	0.04220	0.04322
	10 2,4,5-T	18.090	18.085	0.005	7757899	0.02540	0 02459
	11 DINOSEB 12 2,4-DB	18.580	18.580		6589066	0 16900	0 1666



657 568 Data File: /var/chem/gcl.i/2180.b/a-a20468.d Report Date: 18-May-2000 12:35

# STL-PITTSBURGH

Data file : /var/chem/gcl.i/2180.b/a-a20468.d Lab Smp Id: Hherb

Inj Date : 18-MAY-2000 12:08

Inst ID: gcl.i Operator : 01797

Smp Info : Hherb, 2180.b Misc Info : 190-80-5

Comment

Method : /var/chem/gcl.i/2180.b/LONGH.m Meth Date : 18-May-2000 12:35 eppinged Qu Quant Type: ESTD Cal File: a-a20468.d Cal Date : 18-MAY-2000 12:08

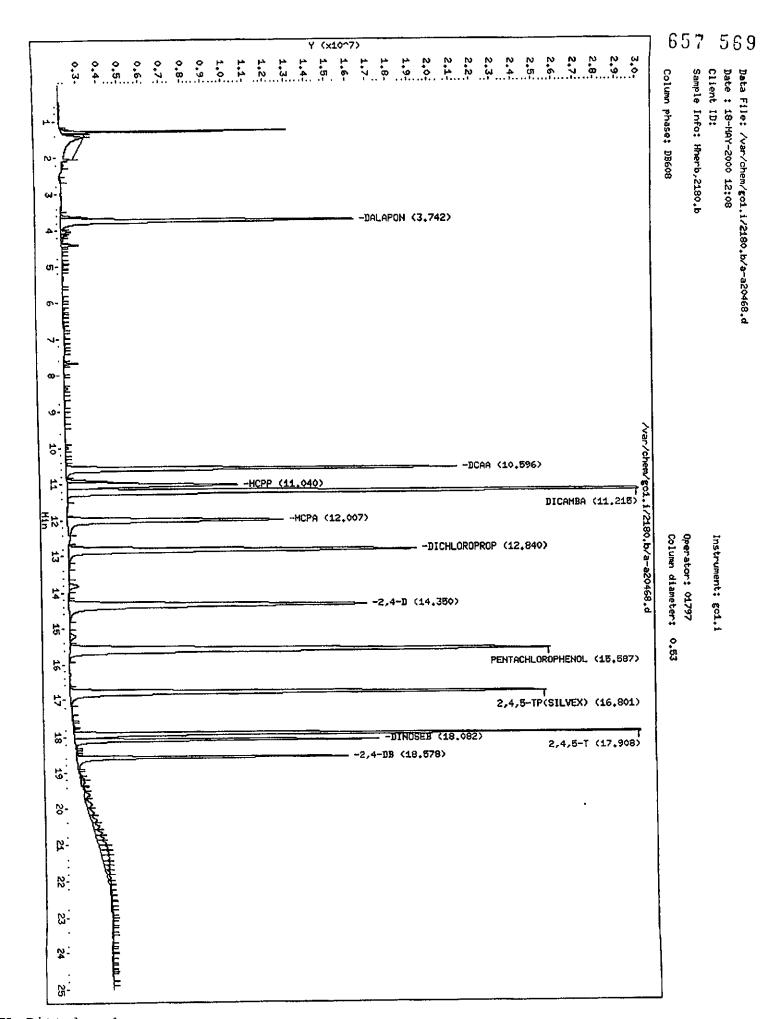
Calibration Sample, Level: 5 Als bottle: 6

Dil Factor: 1.00000

Compound Sublist: 1-all.sub Integrator: Falcon

Target Version: 3.40

						NUOMA	TS
						CAL-AMT	ON-COL
Com	oounds	RŤ	EXP RT D	LT RT	RESPONSE	(ng)	( ng)
	em 보고 프로디 박대용 등 본 참 화로 등 등 휴대용 등 등 경	***	医拉克莱耳亚 河		*======	6203426	2222E
	1 DALAPON	3.742	3.742	0.000	14136953	0.17600	0.1586
\$	2 DCAA	10.596	10.600	-0.004	84790711	0.34000	0 2795
•	3 MCPP	11.040	11 040	0 000	8263392	34.1000	17.41
	4 DICAMBA	11.215	11.218	-0.003	34577437	0.17000	0.1547
	5 MCPA	12.007	12.006	0 001	10445407	34.0000	18.59
	6 DICHLOROPROP	12.840	12.837	0.003	16846334	0.33900	0.2886
	7 2,4-D	14.350	14 351	-0.001	14484500	0.34000	0.3053
	8 PENTACHLOROPHENOL	15.587	15.586	0.001	23432763	0.04255	0.04224
	9 2.4,5-TP(SILVEX)	16.801	16 798	0 003	23172887	0.08400	0.08222
	10 2,4,5-T	17 908	17 910	-0 002	28087411	0 08440	0 08513
	11 DINOSEB	18.082	18.085	-0.003	14833311	0.05080	0.04773
	12 2,4-DB	18.578	18.580	-0.002	13169053	0.33800	0.3340



Data File: /var/chem/gc1.i/2180.b/a-a20529.d

Report Date: 20-May-2000 11:44

# STL-PITTSBURGH

Data file : /var/chem/gc1.i/2180.b/a-a20529.d

Lab Smp Id: Mherb

Inj Date : 20-MAY-2000 11:04

Operator: 01797 Inst ID: gcl.i

Smp Info : Mherb,2180.b Misc Info : 190-80-3

Comment

Method : /var/chem/gcl.i/2180.b/LONGH.m

Meth Date: 20-May-2000 11:44 g Quant Type: ESTD Cal Date: 18-MAY-2000 12:08 Cal File: a-a20468.d

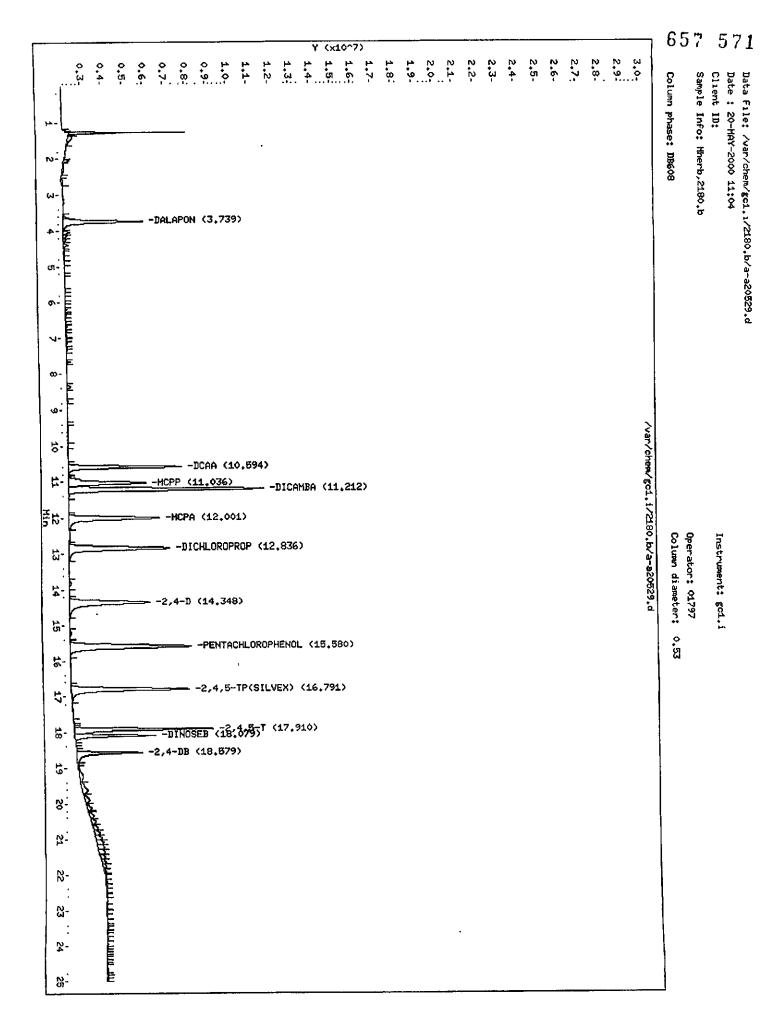
Als bottle: 2 Continuing Calibration Sample

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: all.sub

Target Version: 3.40

						AMOUN	TS
						CAL-AMT	ON-COL
Comp	ounds	RT	EXP RT	DLT RT	respon <b>s</b> e	( ng)	( ng)
====	- 有限的自主なではことが表示可能を表示されてこ	<b>=</b> =		****			### <b>##</b>
	1 DALAPON	3 739	3.739	0.000	3830936	0 04390	0.04299
Ş	2 DCAA	10 594	10.594	0 000	25434014	0 08510	0 08383
	3 MCPP	11.036	11.036	0.000	3736465	B 52000	7.872
	4 DICAMBA	11.212	11.212	0.000	9414834	0.04250	0.04212
	5 MCPA	12.001	12.001	0 000	4406821	8.56000	7 843
	6 DICHLOROPROP	12 836	12.836	0 000	4885129	0 08480	0.08368
	7 2,4-D	14.348	14.348	0 000	3910530	0.08510	0.08242
	8 PENTACHLOROPHENOL	15 580	15.580	0.000	5892468	0 01064	0.01062
	9 2,4,5-TP(SILVEX)	16.791	16.791	0 000	5714110	0.02110	0 02027
1	10 2,4,5-T	17.910	17.910	0.000	6737635	0 02110	0 02042
1	1 DINOSEB	18.079	18 079	0 000	3929837	0.01270	0 01264
1	12 2,4-DB	18 579	18.579	0 000	3201210	0.08450	0 08120



Data File: /var/chem/gc1.i/2180.b/a-a20550.d

Report Date: 22-May-2000 08:09

### STL-PITTSBURGH

Data file: /var/chem/gc1.i/2180.b/a-a20550.d

Lab Smp Id: Mherb

Inj Date : 20-MAY-2000 21:15

Operator: 01797 Inst ID: gcl.i

Smp Info : Mherb,2180.b Misc Info : 190-80-3

Comment :

Method : /var/chem/gc1.i/2180.b/LONGH.m

Meth Date : 22-May-2000 08:08 g Quant Type: ESTD Cal Date : 18-MAY-2000 12:08 Cal File: a-a20468.d

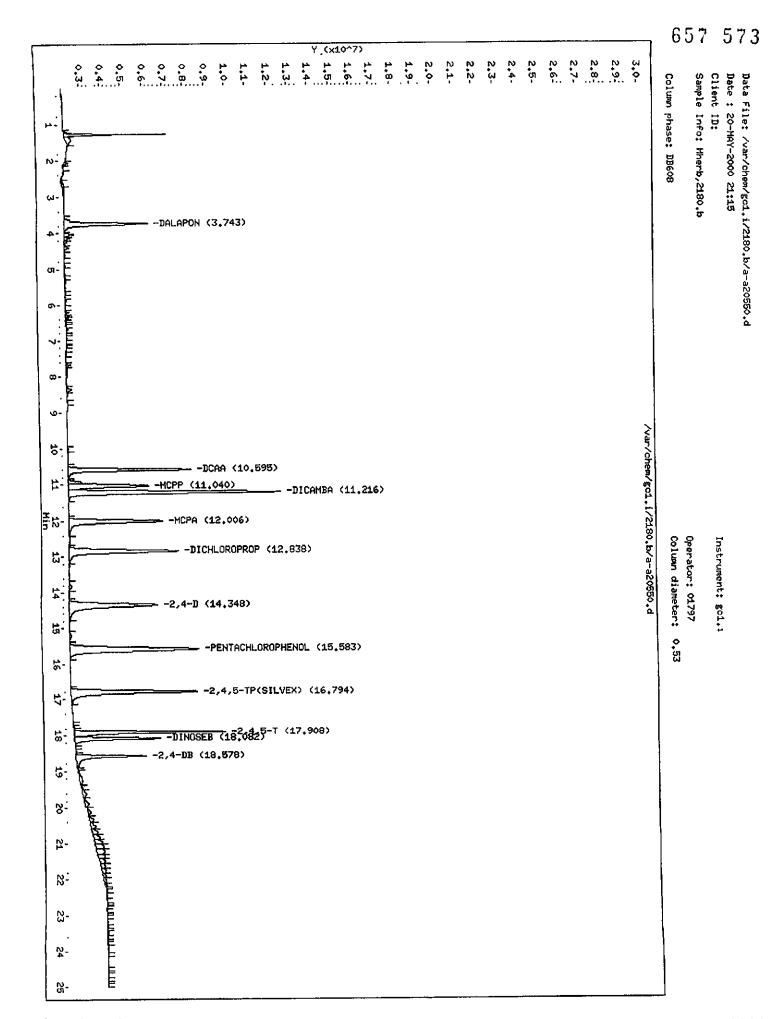
Als bottle: 23 Continuing Calibration Sample

Dil Factor: 1.00000 Integrator: Falcon

Integrator: Falcon Compound Sublist: all.sub

Target Version: 3.40

						AMOUN	TS
						CAL-AMT	ON-COL
Compo	ounds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( ng)
		<b>==</b>		DF====	*****		
1	DALAPON	3.743	3 739	0.004	4041087	0.04390	0.04535
\$ 2	2 DCAA	10.595	10 594	0.001	27612144	0.08510	0.09101
3	MCPP	11.040	11.036	0.004	3879025	8 52000	8 172
4	DICAMBA	11 216	11.212	0.004	10271232	0.04250	0 04596
5	MCPA	12.006	12.001	0.005	4562486	8 56000	8.120
•	DICHLOROPROP	12.838	12.836	0 002	5265248	0.08480	0.09019
7	7 2,4-D	14.348	14.348	0.000	4252615	0 08510	0.08963
8	PENTACHLOROPHENOL	15 583	15 580	0.003	6238749	0 01064	0.01124
9	2,4,5-TP(SILVEX)	16 794	16 791	0.003	6134288	0 02110	0.02176
10	2,4,5-T	17 908	17.910	-0 002	7294783	0.02110	0 02211
1.1	DINOSEB	18 082	18.079	0.003	4157301	0.01270	0.01338
13	2 2,4-DB	18.578	18.579	-0.001	3371666	0.08450	0 08552



HERBICIDE QC DATA



# UXB INTERNATIONAL METHOD BLANK COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATER Lab Sample ID:C0E180000 489

Method: SW846 8151A Luns 200

Herbicides (8151A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DDE4W101 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/20/00

Moisture %:NA

QC Batch: 0139489

Client Sample Id: INTRA-LAB BLANK

# CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L C	}
94-75-7	2,4-D	0.040	ן <u>ט</u>
93-72-1	2,4,5-TP (Silvex)	[0.010	<u> </u>

657 576 Data File: /var/chem/gc1.i/2180.b/a-a20536.d

Report Date: 22-May-2000 08:55

#### STL-PITTSBURGH

Data file : /var/chem/gc1.i/2180.b/a-a20536.d

Lab Smp Id: DDE4W101 Client Smp ID: PBLK

Inj Date : 20-MAY-2000 14:27

Operator: 01797 Smp Info: DDE4W101,2180.b Misc Info: 130194BLK Inst ID: gcl.i

Comment :

Method : /var/chem/gc1.i/2180.b/LONGH.m Meth Date : 22-May-2000 08:08 g Qu Quant Type: ESTD Cal File: a-a20468.d Cal Date : 18-MAY-2000 12:08

QC Sample: BLANK Als bottle: 9

Dil Factor: 1.00000 Integrator: Falcon Compound Sublist: all.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* 20\*Vt/Vo/Vi

Name	Value	Description
DF Vt Vo Vi	10.000 100.000	Dilution Factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected

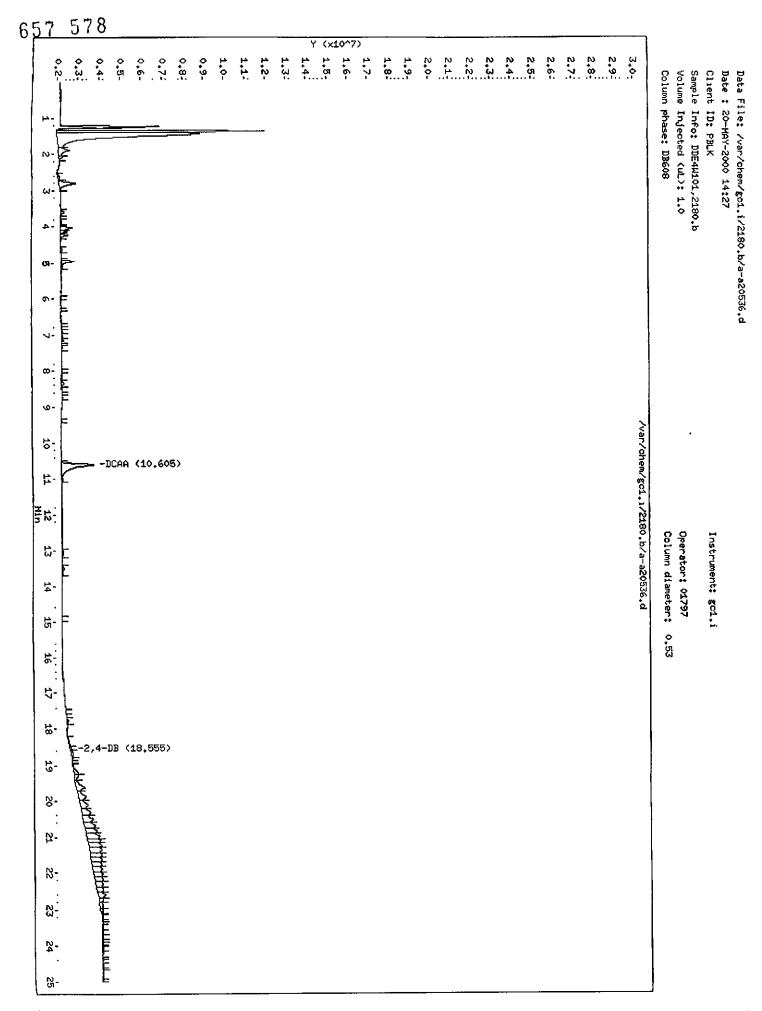
#### CONCENTRATIONS

						ON-COLUMN	FINAL
Com	pounds	RT	EXP RT D	LT RT	RESPONSE	( ng)	( mg/L)
2 C =		**		====	******	0000000	******
	1 DALAPON	Con	pound Not	Detecte	d		
\$	2 DCAA	10.605	10 594	0.011	11821468	0.03896 (	0.07793(a)
	3 MCPP	Con	pound Not	Detecte	ed.		
	4 DICAMBA	Con	pound Not	Detecte	d.		
	5 MCPA	Con	pound Not	Detecte	d.		
	6 DICHLOROPROP	Con	pound Not	Detecte	d.		
	7 2,4-D	Cot	npound Not	Detecte	d.		
	8 PENTACHLOROPHENOL	Con	pound Not	Detecte	ed		
	9 2,4,5-TP(SILVEX)	Con	pound Not	Detecte	d		
	10 2,4,5-T	Con	npound Not	Detecte	ed.		
	11 DINOSEB	Cor	npound Not	Detecte	eđ.		
	12 2,4-DB	18.555	18.579	-0 024	83111	0 00211	0.004216(a)

Data File: /var/chem/gcl.i/2180.b/a-a20536.d Report Date: 22-May-2000 08:55

#### QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).



#### UXB INTERNATIONAL CHECK SAMPLE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATL Lab Sample ID: C0E180000 489

Method: SW846 8151A w 5/22/00 Herbicides (8151A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DDE4W102 Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/20/00

Moisture %:NA

QC Batch: 0139489

Client Sample Id: CHECK SAMPLE

#### CONCENTRATION UNITS:

CA	S NO.	COMPOUND ((	ug/L or ug/kg	) mg/L	Q
94-75	-7 2,4	-D	١٥	133	l
93-72	-1 2,4	,5-TP (Silvex)	0.	0323	

Data File: /var/chem/gc1.i/2180.b/a-a20537.d

Report Date: 22-May-2000 08:55

#### STL-PITTSBURGH

Data file: /var/chem/gc1.i/2180.b/a-a20537.d

Client Smp ID: LCS Lab Smp Id: DDE4W102

Inj Date : 20-MAY-2000 14:56 Operator : 01797 Smp Info : DDE4W102,2180.b Inst ID: gcl.i

Misc Info : 130194LCS

Comment

Method : /var/chem/gcl.i/2180.b/LONGH.m

Meth Date: 22-May-2000 08:08 g Quant Type: ESTD Cal Date : 18-MAŶ-2000 12:08 Cal File: a-a20468.d

Als bottle: 10 QC Sample: LCS

Dil Factor: 1.00000 Compound Sublist: all.sub Integrator: Falcon

Target Version: 3.40

Concentration Formula: Amt \* DF \* 20\*Vt/Vo/Vi

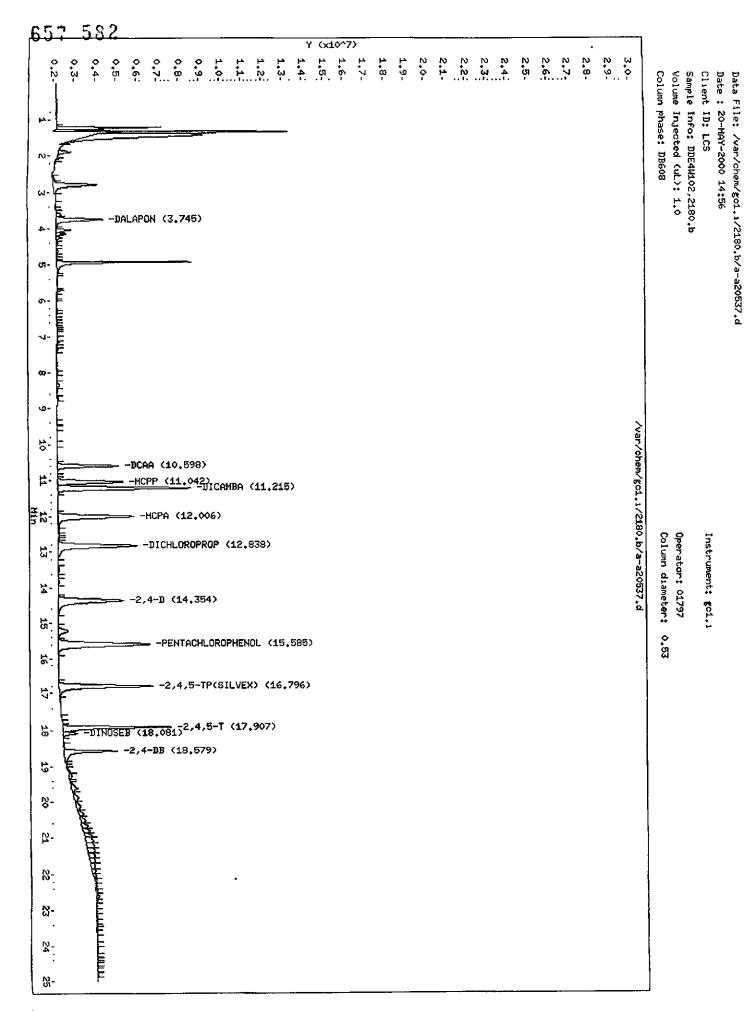
Name	Value	Description
DF Vt Vo Vi	10.000 100.000	Dilution Factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected

						CONCENTRA	ations
						ON-COLUMN	FINAL
Compo	unds	RT	EXP RT	DLT RT	RESPONSE	( ng)	( mg/L)
	=======================================	77.75			******	三三	9554444
1	DALAPON	3.745	3.739	0 006	2276647	0.02555	0 05110(aR)
\$ 2	DCAA	10.598	10.594	0 004	13509516	0.04453	0.08906(a)
3	МСРР	11.042	11.036	0.006	3165066	6 66777	13 34
4.	DICAMBA	11.215	11.212	0 003	6402544	0.02865	0.05729(a)
5	MCPA	12 006	12.001	0.005	3658101	6.51079	13 02
6	DICHLOROPROP	12 838	12.836	0.002	3790550	0.06493	0.1299(a)
7	2,4-D	14.354	14.348	0.006	3159564	0.06659	0 1332 (aR)
8	PENTACHLOROPHENOL	15.585	15.580	0 005	4402096	0 00793	0.01587(a)
9	2,4,5-TP(SILVEX)	16 796	16 791	0.005	4553679	0 01616	0.03231 (aR)
10	2,4,5-T	17.907	17.910	-0 003	5233112	0 01586	0.03172 (aR)
11	DINOSEB	18 081	18 079	0.002	693827	0 00223	0.004465(a)
12	2,4-DB	18.579	18.579	0.000	2599680	0.06594	0.1319(a)

Data File: /var/chem/gc1.i/2180.b/a-a20537.d Report Date: 22-May-2000 08:55

### QC Flag Legend

a - Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).R - Spike/Surrogate failed recovery limits.



# UXB INTERNATIONAL MATRIX SPIKE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WHEN Lab Sample ID:COE130194 001

Method: SW846 8151A W5/22/00

Herbicides (8151A)

Sample WT/Vol: 100 / mL Date Received: 05/13/00 Work Order: DD6A411D Date Extracted:05/18/00 Dilution factor: 1 Date Analyzed: 05/20/00

Moisture %:10

QC Batch: 0139489

Client Sample Id: DF/S1/0133/SDC/001 A

#### CONCENTRATION UNITS:

_	CAS NO.	COMPOUND	(ug/L or ug	/kg) mg/L	<u>Q</u>
Ī	94-75-7	2,4-D		0.133	
į	93-72-1	2,4,5-TP (Silvex)		0.0322	

Data File: /var/chem/gcl.i/2180.b/a-a20531.d

Report Date: 22-May-2000 08:55

#### STL-PITTSBURGH

Data file: /var/chem/gc1.i/2180.b/a-a20531.d Lab Smp Id: DD6A411D Clien Client Smp ID: DF/S1/0133/SDC/MS

Inj Date : 20-MAY-2000 12:02

Operator : 01797 Inst ID: qc1.i

Smp Info : DD6A411D,2180.b

Misc Info: 1301940018

Comment

Method : /var/chem/gc1.i/2180.b/LONGH.m Meth Date : 22-May-2000 08:08 g Qu Quant Type: ESTD Cal Date : 18-MAY-2000 12:08 Cal File: a-a20468.d

Als bottle: 4 QC Sample: MS

Dil Factor: 1.00000

Integrator: Falcon Compound Sublist: all.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* 20\*Vt/Vo/Vi

Name	Value	Description
DF Vt Vo Vi	10.000 100.000	Dilution Factor Volume of final extract (uL) Volume of sample extracted (mL) Volume injected

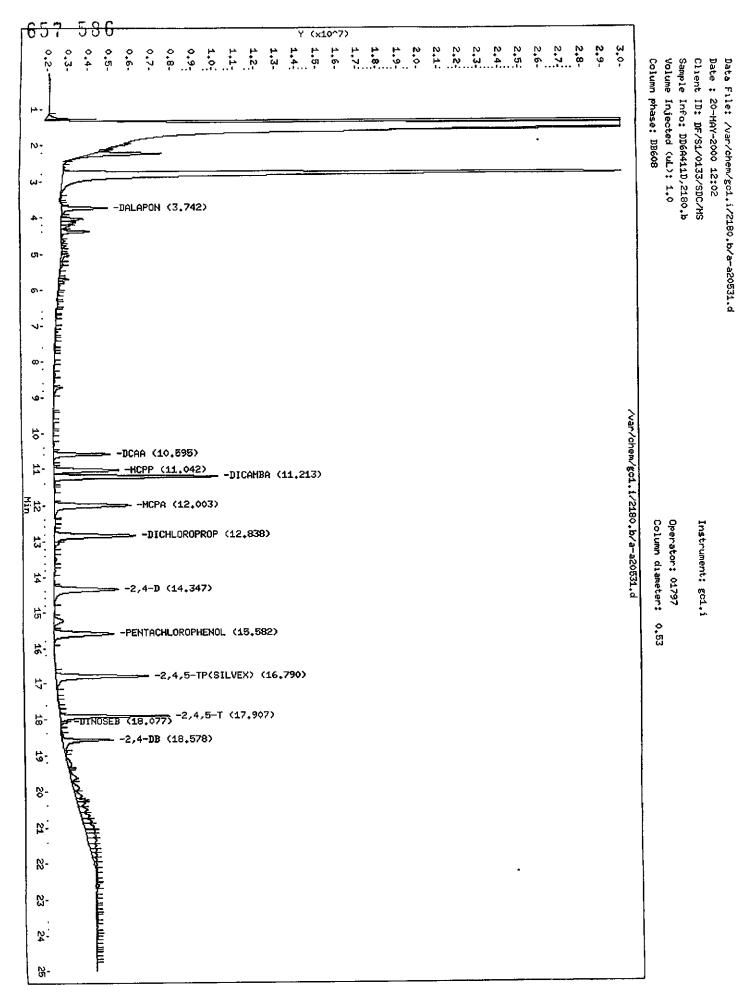
						CONCENTRA	ations	
						ON-COLUMN	FINAL	
Compou	nds	RT	EXP RT	DLT RT	RESPONSE	(ng)	( mg/L)	
======	~	==	22232					
1	DALAPON	3.742	3.739	0.003	2236189	0.02510	0.05019(aR)	
\$ 2	DCAA	10.595	10.594	0.001	12391207	0 04084	0.08168(a)	
3 1	MCPP	11 042	11.036	0.006	3180908	6.70115	13.40	
4	DICAMBA	11.213	11.212	0.001	7895838	0 03533	0 07066(a)	
5	MCPA	12.003	12.001	0.002	3732991	6 64408	13.29	
6	DICHLOROPROP	12.838	12.836	0.002	3944797	0 06757	0.1351(a)	
7	2.4-D	14.347	14.348	-0 001	3150651	0.06640	0 1328 (aR)	>
8	PENTACHLOROPHENOL	15 582	15.580	0.002	2896901	0 00522	0 01044(a)	
9	2,4,5-TP(SILVEX)	16.790	16.791	-0.001	4532904	0.01608	0.03216(aR)	)
10	2,4,5-T	17.907	17.910	-0.003	5292004	0 01604	0 03208 (aR)	
11	DINOSEB	18.077	18.079	-0.002	324719	0.00104	0.002090(a)	
12	2,4-DB	18 578	18.579	-0.001	2494486	0 06327	0.1265(a)	

Data File: /var/chem/gc1.i/2180.b/a-a20531.d Report Date: 22-May-2000 08:55

657 585

#### QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).R Spike/Surrogate failed recovery limits.



#### UXB INTERNATIONAL MATRIX SPIKE DUPLICATE COMPOUNDS

Lab Name: Severn Trent Laboratories, Inc. SDG Number:

Matrix: (soil/water) SOLID WATEA Lab Sample ID: COE130194 001

Method: SW846 8151A Lun staa to

Herbicides (8151A)

Date Received: 05/13/00 Sample WT/Vol: 100 / mL Date Extracted: 05/18/00 Work Order: DD6A411E Date Analyzed: 05/20/00 Dilution factor: 1

Moisture %:10

QC Batch: 0139489

Client Sample Id: DF/S1/0133/SDC/001 A

#### CONCENTRATION UNITS:

CAS NO.	COMPOUND	(ug/L or ug/kg) mg/L	<u>Q</u>
94-75-7	2,4-D	0.144	
93-72-1	2,4,5-TP (Silvex)	0.0339	ll

657 538

Data File: /var/chem/gc1.i/2180.b/a-a20532.d

Report Date: 22-May-2000 08:55

#### STL-PITTSBURGH

Data file: /var/chem/gc1.i/2180.b/a-a20532.d Lab Smp Id: DD6A411E Clien Client Smp ID: DF/S1/0133/SDC/MSD

Ing Date : 20-MAY-2000 12:31

Operator : 01797 Inst ID: gcl.i

Smp Info : DD6A411E,2180.b

Misc Info : 130194001D

Comment :

Method : /var/chem/gc1.i/2180.b/LONGH.m Meth Date : 22-May-2000 08:08 g Quant Type: ESTD Cal Date : 18-MAY-2000 12:08 Cal File: a-a20468.d

Als bottle: 5 QC Sample: MSD

Dil Factor: 1.00000 Integrator: Falcon Compound Sublist: all.sub

Target Version: 3.40

Concentration Formula: Amt \* DF \* 20\*Vt/Vo/Vi

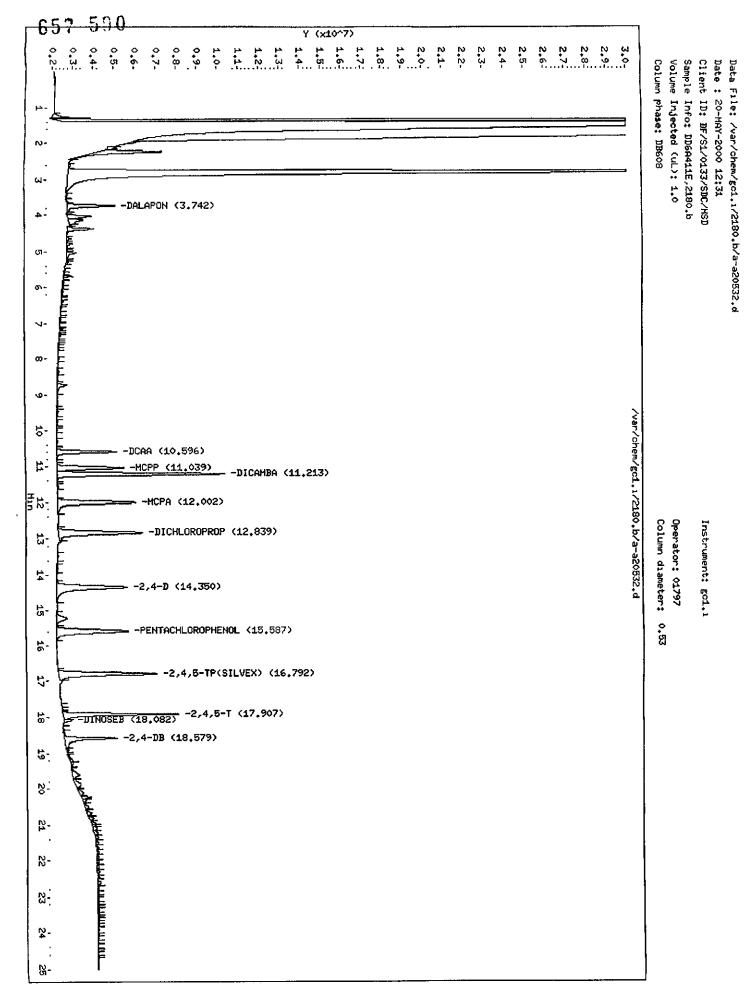
Name	Value	Description
DF Vt	10.000	Dilution Factor Volume of final extract (uL)
Vo Vi		Volume of sample extracted (mL) Volume injected

						CONCENTR	ations
						ON-COLUMN	FINAL
Compounds		RT	EXP RT	DLT RT	RESPONSE	( ng)	( mg/L)
**********		==	=====				F33555
1 DALAPON		3.742	3.739	0.003	2308136	0 02590	0.05181(aR)
\$ 2 DCAA		10.596	10 594	0.002	13014217	0 04290	0.08579(a)
3 MCPP		11.039	11 036	0.003	3251457	6 84977	13 70
4 DICAMBA		11.213	11.212	0.001	8123697	0.03635	0 07270(a)
5 MCPA		12.002	12 001	0.001	3820491	6 79981	13 60
6 DICHLOR	OPROP	12.839	12 836	0 003	4104129	0.07030	0.1406(a)
7 2,4-D		14.350	14.348	0 002	3427686	0.07224	0 1445 (aR)
8 PENTACH	LOROPHENOL	15 587	15.580	0 007	3454777	0.00623	0 01245(a)
9 2,4,5-T	P(SILVEX)	16 792	16.791	0.001	4774343	0.01694	0 03388 (aR)
10 2,4,5-T		17.907	17.910	-0.003	5612984	0.01701	U. U3402 (aR)
11 DINOSEB		18.082	18 079	0 003	413389	0.00133	0.002660(a)
12 2,4-DB		18 579	18.579	0 000	2589940	0.06570	0.1314(a)

Data File: /var/chem/gc1.i/2180.b/a-a20532.d Report Date: 22-May-2000 08:55

## QC Flag Legend

- a Target compound detected but, quantitated amount Below Limit Of Quantitation(BLOQ).R Spike/Surrogate failed recovery limits.



# HERBICIDE MISCELLANEOUS

C. Godium Sulfate Mfg. L. R. Bes.  Quanterra N:\QA\LOGBOOKS\sepfun extr doc May-00	3050	Extract(s) (Record line number from above)	24 Analyst	21 22 22	19	15. 16.	13	10.	4	50%	Separatory Funnel Extraction Worksheet  Date Extraction Began Date Completed S-(8-00 S-19-0 Lot Number Sa
Lot Number  May-00	00	Time 7	My hy						004	501 001ms	Sample ID  Sample ID  Client ID  A  A  A  A  A  A  A  A  A  A  A  A  A
op-00-0035	M. Mr. REHIZ-E	Extract(s) Received ( Analyst Location  Usasilians Cris Paper	DY DY IM								thod Solvent Solvent Solvent Polume Final Surre (ml.) (ml.) (ml.) (ml.)
Reviewed By Muse-		Date Time 7.28	ħγ ħγ						8	79	39 489 HeX-N442:  Mo-89-3 1, 0  Value of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the control of the contro
Page 16 of 80		Extract(s) Relinquished Location Analyst  Middig Grand Color Color	10 10 VV			Justanie			*	0-79-14 1,C	STL Piusb 450 Willia Piusburgh 412-820-8 No Matrix Sp Volume (t

|

5 By Propa Numb 593 DOGAY <5 = Extraction Fluid No. 1 5.7 mL Glacial Acetic Acid dil 500 mL + 64 3 mL of 1N Na OH dil to 1L (pH 4.93  $\pm$  0.05) >5 = Extraction Fluid No. 2 5 7 mL Glacial Acetic Acid dll to 1L (pH 2.88  $\pm$  0.05) DOGAS 2007 DOCAL (Record line number from above) = Sample determined to have free-liquid, % solid determination was performed. ROOK JEMP -Sample ID Extract(s) 5,00,/165. Cociso 194 WgtVol Quanterra incorporated 450 William Pitt Way Pittsburgh, Pennsylvania 15238 412/826-5477 FAX: 412/826-5571 5-160 Ú Date 126-00 13 Fluid Determination 25.2 भ अभ 1.3C init pH 8 3/21/5 Time Final pH 0 Lisam famos 111.655 Extract(s) Respired # S Extract Fluid #1, or #2 Malyst Conc : HNO3
to Acidity
(for Metals) ځ 3 rotten Location Filter lot # 8520701W S-16-50 0 - C - Sad 5-16-00 Date/Time On edde A4 5DDXA7 2. DOGA4 575 4DOM6 70 ø 3 DDGAS 12 7 20.9% d-127-80 Date Sample ID 12:45 Accument ARDS J-16-00 12:00 Solution Number 1 - Log Book Number pH Calibration 995404-17 Time u 5-1700 00.00 20.00 Wgt (g) Wet/Diy TCLP Extraction 0 Extract(s) Relinquis/jed SN9087030 Serial Number 4.00 4.00 9:30 2000 Analyst Fluid Vol. (4.0)企 038 Solution Number 2 - Log Book Number 30±2 Agitation Apparatus APMs. Gass Ċ δ 7.0) 96-MT-526 C44817 Environmedial Services Final pH After Tumbling 20 10.030 Location 2.0 (10.0) 26 Ρi STL

TCLP (Method 1311)

*"manterra* 

657 594 Sequence: C:\HPCHEM\2\SEQUENCE\2200.S

# Cracted: Dr. 5-70-00 Sequence Table (Front Injector):

Vial Information Part:

,			
	Vial	Vial Information	=======================================
1	1	RINSE	
2	2	190-80-3 529	$\sqrt{\omega}$
3	3	130194001	\ uA
4	4	130194001S	422/00
5	5	130194001D	2/2/
6	6	130194002	$c \rightarrow 0$
7	7	130194003	conf. regil
8	8	130194004 <b>535</b>	
9	9	130194BLK	
10	10	130194LCS	
11	11	100126003	
12	12	100126004 ZIOX	
13	13	100126005540)	
14	14	100126006	
15	15	100126BLK	
16	16	100126LCS	
17	17	100126LCD	
18	18	090125001 <b>54</b>	
19	19	090125002	
20	20	090125BLK	
21	21	090125LCS	
22	22	090125LCD	
23	23	190-80-3 55V	
24	24	100183001	
25	25	100183001S	

	Vial		
26	26	100183001D	
27	27	100183002	
28	28	100183BLK 555	
29	29	100183LCS	
30	30	110158001	
31	31	1101580018	
32	32	110158001D	
33	33	110158BLK 560	•
34	34	110158LCS	
35	35	110226001.	
36	36	110226002	
37	37	110226003	
38	38	110226004 565	
39	39	110226005	
40	40	110226006'	
41	41	1102260065	
42	42	110226006D	
43	43	110226BLK 570	
44	44	190-80-3	
45	45	110226LCS	
46	46	110226007	
47	47	110226008	
48	48	110226009 575	
49	49	110226010	
50	50	110226011	
51	51	110226012.	
52	52	120236010	
53	53	120166002 <b>580</b>	

Line	Vial	Vial Information
54	54	120166003
55	55	120166005
56	56	120166006
57	57	120166007
58	58	120137001 585
59	59	130190001
60	60	130190002
61	61	160131001
62	62	120137BLK
63	63	120137LCS 590
64	64	120137LCD
65	65	190-80-3

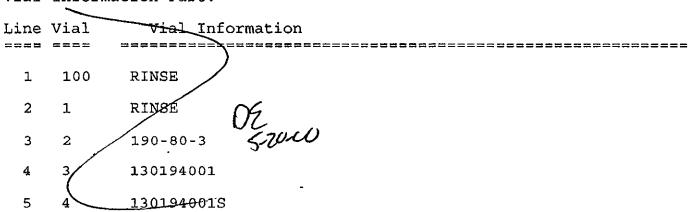
Method and Injection Info Part:

Line	Vial	SampleName	Method	Inj	SampleType	InjVolume	DataFile
====	====	===============	=======	===	========	==== <b>=</b> ====	========
					_		
1	1	HEXANE	HERB	1	Sample		
2	2	Mherb, 2180.b	HERB	1	Sample		
3	3	DD6A4104,2180.b	HERB	1	Sample		
4	4	DD6A411D,2180.b	HERB	1	Sample		
4 5 6	5	DD6A411E,2180.b	HERB	1	Sample		
6	6	DD6A5104,2180.b	HERB	1	Sample		
7	7	DD6A6104,2180.b	HERB	1	Sample		
8 9	8	DD6A7104,2180.b	HERB	1	Sample		
9	9	DDE4W101,2180.b	HERB	1	Sample		
10	10	DDE4W102,2180.b	HERB	1	Sample		
11	11	DD0LQ101,2180.b	HERB	1	Sample		
12	12	DD0LR101,2180.b	HERB	1	Sample		
13	13	DD0LT101,2180.b	HERB	1	Sample		
14	14	DD0LV101,2180.b	HERB	1	Sample		
15	15	DD1RV101,2180.b	HERB	1	Sample		
16	16	DD1RV102,2180.b	HERB	1	Sample		
17	17	DD1RV103,2180.b	HERB	1	Sample		
18	18	DCX46102,2180.b	HERB	1	Sample		
19	19	DCX4H102,2180.b	HERB	1	Sample		
20	20	DD05F101,2180.b	HERB	1	Sample		
21	21	DD05F102,2180.b	HERB	1	Sample		
22	22	DD05F103,2180.b	HERB	1	${ t Sample}$		
23	23	Mherb,2180.b	HERB	1	Sample		
24	24	DD16N102,2180.b	HERB	1	Sample		
25	25	DD16N106,2180.b	HERB	1	Sample		
26	26	DD16N107,2180.b	HERB	1	Sample		
27	27	DD17C102,2180.b	HERB	1	Sample		

		SampleName	Method	Inj	SampleType	-	
====	====	=======================================	=======	===	=======	======	=======
28	28	DD1T5101,2180.b	HERB	1	Sample		
29	29	DD1T5102,2180.b	HERB	1	Sample		
30	30	DD2LG101,2180.b	HERB	1	Sample		
31	31	DD2LG102,2180.b	HERB	1	Sample		
32	32	DD2LG103,2180.b	HERB	1	Sample		
33	33	DDE53101,2180.b	HERB	1	Sample		
34	34	DDE53102,2180.b	HERB	ī	Sample		
35	35	DD3CP102,2180.b	HERB	1	Sample		
36	36	DD3CQ102,2180.b	HERB	ī	Sample		
37	37	DD3CR101,2180.b	HERB	ī	Sample		
38	38	DD3CT101,2180.b	HERB	1	Sample		
39	39	DD3CW101,2180.b	HERB	1	Sample		
40	40	DD3D1101,2180.b	HERB	1	Sample		
41	41	DD3D1102,2180.b	HERB	1	Sample		
42	42	DD3D1103,2180.b	HERB	1	Sample		
43	43	DD931101,2180.b	HERB	1	Sample		
44	44	Mherb, 2180.b	HERB	1	Sample		
45	45	DD931102,2180.b	HERB	1	Sample		
46	46	DD3DD101,2180.b	HERB	1	Sample		
47	47	DD3DK101,2180.b	HERB	1	Sample		
48	48	DD3DM101,2180.b	HERB	1	Sample		
49	49	DD3DP101,2180.b	HERB	1	Sample		
50	50	DD3DQ101,2180.b	HERB	1.	Sample		
51	51	DD3DR101,2180.b	HERB	1	Sample		
52	52	DD59M101,2180.b	HERB	1	Sample		
53	53	DD4JG102,2180.b	HERB	1	Sample		
54	54	DD4JT102,2180.b	HERB	1	Sample		
55	55	DD4K0102,2180.b	HERB	1	Sample		
56	56	DD4K3102,2180.b	HERB	1	Sample		
57	57	DD4K5102,2180.b	HERB	1	Sample		
58	58	DD4A6102,2180.b	HERB	1	Sample		
59	59	DD693102,2180.b	HERB	1	Sample		
60	60	DD694102,2180.b	HERB	1	Sample		
61	61	DD7VM102,2180.b	HERB	1	Sample		
62	62	DDE54101,2180.b	HERB	1	Sample		
63	63	DDE54102,2180.b	HERB	1	Sample		
64	64	DDE54103,2180.b	HERB	1	Sample		
65	65	Mherb,2180.b	HERB	1	Sample		

### Sequence Table (Back Injector):

Vial Information Part:



657 598 Crontod: DE 5-19-00

Sequence Table (Front Injector):

### Vial Information Part:

VIGI	111101		ı
Line	Vial	Vial Information	:=====================================
1	1	RINSE	
2	2	190-80-1 464	Dola Comment
3	3	190-80-2	
4	4	190-80-3	
5	5	190-80-4	
6	6	190-80-5	•
7	7	060132015	
8	8	0601320155 <b>47</b> 0	
9	9	060132015D	
10	10	060132BLK	
11	11	060132LCS	
12	12	050196001	
13	13	0501960018475	
14	14	050196001D	
15	15	050196002	
16	16	050196003.	
17	17	050196004	•
18	18	050196005 480	
19	19	050196006 ·	
20	20	050196007-	
21	21	050196008.	
22	22	050196009	
23	23	050196010 485	
24	24	050196011	
25	25	050196BLK	

	Vial	Vial Information	50 / 58
26	26	050196LCS	
27	27	190-80-3	
28	28	050196012 490	
29	29	050196013	
30	30	050196014	
31	31	050196015	
32	32	050196016	
33	33	050196017455	
34	34	050196018	
35	35	050196019	
36	36	060132001	
37	37	060132001S	
38	38	060132001D SVO	
39	39	060132002	
40	40	060132003	
41	41	060132004	
42	42	060132005	
43	43	060132006 505	
44	44	060132007	
45	45	060132008	
46	46	060132BLK	
47	47	060132LCS	
48	48	190-80-3 <b>510</b>	
49	49	060132009	
50	50	060132010	
51	51	060132011	
52	52	060132012	
53	53	060132013 515	

b5/ Line	000 Vial ====	Vial Information
54	54	060132014
55	55	160130001
56	56	160130001S
57	57	160130001D
58	58	160130002 570
59	59	160130BLK
60	60	160130LCS
61	61	180156001
62	62	180156BLK
63	63	180156LCS 575
64	64	180156LCD
65	65	190-80-3

Method and Injection Info Part:

Line	Vial	SampleName	Method	Inj ===	SampleType	InjVolume	DataFile
Line ====  1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16	Vial ==== 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16	SampleName ====================================	Method =======  HERB HERB HERB HERB HERB HERB HERB HER	Inj === 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	SampleType ====================================	injvolume	Datarile
16 17 18 19 20 21 22 23 24 25 26 27	16 17 18 19 20 21 22 23 24 25 26 27	DCT80101,2180.b DCT8V101,2180.b DCT9V101,2180.b DCT90101,2180.b DCT93101,2180.b DCT94101,2180.b DCT94101,2180.b DCT9A101,2180.b DCT9C101,2180.b DDAVX101,2180.b DDAVX101,2180.b Mherb,2180.b	HERB HERB HERB HERB HERB HERB HERB HERB	1 1 1 1 1 1 1 1 1 1 1 1 1	Sample Sample Sample Sample Sample Sample Sample Sample Sample Sample Sample	•	

		SampleName	Method	Inj	SampleType	_	
	~						
28	28	DCT9E101,2180.b	HERB	1.	Sample		
29	29	DCT9J101,2180.b	HERB	1	Sample		
30	30	DCT9K101,2180.b	HERB	1	Sample		
3.1	31	DCT9L101,2180.b	HERB	1	Sample		
32	32	DCT9M101,2180.b	HERB	1	Sample		
33	33	DCT9Q101,2180.b	HERB	1	Sample		
34	34	DCT9R101,2180.b	HERB	1	Sample		
35	35	DCT9T101,2180.b	HERB	1	Sample		
36	36	DCVHK101,2180.b	HERB	1	Sample		
37	37	DCVHK102,2180.b	HERB	1	Sample		
38	38	DCVHK103,2180.b	HERB	1	Sample		
39	39	DCVHL101,2180.b	HERB	1	Sample		
40	40	DCVHM101,2180.b	HERB	1	Sample		
41	41	DCVHP101,2180.b	HERB	1	Sample		
42	42	DCVHQ101,2180.b	HERB	1	Sample		
43	43	DCVHR101,2180.b	HERB	1	Sample		
44	44	DCVHT101,2180.b	HERB	1	Sample		
45	45	DCVHV101,2180.b	HERB	1	Sample		
46	46	DDAW7101,2180.b	HERB	1	Sample		
47	47	DDAW7102,2180.b	HERB	1	Sample		
48	48	Mherb, 2180.b	HERB	1	Sample		
49	49	DCVHV101,2180.b	HERB	1	Sample		
50	50	DCVHW101,2180.b	HERB	1	Sample		
51	51	DCVHX101,2180.b	HERB	1	Sample		
52	52	DCVJ0101,2180.b	HERB	1	Sample		
53	53	DCVJ1101,2180.b	HERB	1	Sample		
54	54	DCVJ2101,2180.b	HERB	1	Sample		
55	55	DD7V9102,2180.b	HERB	1	Sample		
56	56	DD7V9104,2180.b	HERB	1	Sample		
57	57	DD7V9105,2180.b	HERB	1	Sample		
58	58	DD7VJ102,2180.b	HERB	1	Sample		
59	59	DD98R101,2180.b	HERB	1	Sample		
60	60	DD98R102,2180.b	HERB	1	Sample		
61	61	DDCHQ102,2180.b	HERB	1	Sample		
62	62	DDE86101,2180.b	HERB	1	Sample		
63	63	DDE86102,2180.b	HERB	1	Sample		
64	64	DDE86103,2180.b	HERB	1	Sample		
65	65	Mherb,2180.b	HERB	1	Sample		

## Sequence Table (Back Injector):

```
Vial Information Part.
                Vial Information
Line Vial
      100
              RINSE
  1
              RINSE
  2
      1
  3
      2
              190-80-1
      3
              190-80-2
  5
      4
              190-80-3
```

PSR024

5/18/00

13:49:54 MT

SAMPLE CUSTODIAN REMOVAL REQUEST

PAGE 001

REQUESTED BY: YUSHINSC

METHOD: QS Herbicides (8151A)

STORAGE LOCATION	WORK ORDER #	PICKED CNTR#	CONTROL #	CLIENT #	ANALYSIS	LOTID	SMP#	<u>sfx</u>	MATRIX DESCRIPTION		ry at CVD re	
2F CLP1	DD6A4-1-04		234894	399411	A-64-QS	C0E130194	001	SOLID		0	3	1
2F CLP1	DD6A5-1-04	<del></del>	234895	399411	A-64-QS	C0E130194	002	SOLID		0	3	1
2F CLP1	DD6A6-1-04	<del></del>	234896	399411	A-64-QS	COE130194	003	SOLID		0	3	1
2F CLP1	DD6A7-1-04		234897	399411	A-64-QS	C0E130194	004	SOLID		0	3	1

Gyskinski Gyskinski	Huskinski Huskinski	5-18-00 1540 5-18-00 221

\*\*\*\*\* END OF REPORT \*\*\*\*\*

METALS DATA

### Cover Page - Inorganic Analysis Data Package

	Client ID	Lab Sample ID:	
	DF/S1/0133/SDC/001 A	DD6A4	
	DF/S1/0133/SDC/001 AD	DD6A4D	
	DF/S1/0133/SDC/001 AS	DD6A4S	
	DF/S1/0133/SDC/001 B	DD6A5	
	DF/S1/0133/SDC/001 C	DD6A6	
	DF/S1/0133/SDC/001 D	DD6A7	
technically a	C0E130194 - TOTALS 6010B  this data package is in compliance with the nd for completeness, for other than condition	ns detailed above. Release of the data comb	
	y data package and in the computer-readable ry Manager or the Manager's designee, as ve		rized by
Signature:		Name:	<del></del>
Date:		Title:	<del></del>

REVIEWED BY: MTW DATE: 5-21-00

Version 3.63.3

Cover Page Equivalent

METALS RESULTS

. .

# Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A4 Client ID: DF/S1/0133/SDC/001 A

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 10.39

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	189.04	0.29	1.1	7.7		1	ICPST	5/17/00	9:15

Comments: Lot #. C0E130194 Sample #: 1

## Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A5 Client ID: DF/S1/0133/SDC/001 B

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 8.64

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	189.04	0.28	1.1	7.2		1	ICPST	5/17/00	9:32

657 608

## STL-Pittsburgh

# Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A6 Client ID: DF/S1/0133/SDC/001 C

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 10.27

Element	WL/ Mass	MDL	Report Limit	Сопс	Q	DF	Instr	Anal Date	Anal Time
Arsenic	189.04	0.29	1.1	8.1		1	ICPST	5/17/00	9:36

Comments: Lot #: C0E130194 Sample #: 3

# Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A7 Client ID: DF/S1/0133/SDC/001 D

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 10.81

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	189.04	0.29	1.1	7.7		1	ICPST	5/17/00	9:48

657 610

## Metals Data Reporting Form

Initial Calibration Verification Standard Units: ug/L Instrument: ICPST Acceptable Range: 90% - 110% Chart Number: T00517A.ARC Standard ID: 0014-079-6 Standard Source: Inorganic Ventures ICV3-1 5/17/00 8:48 AM WL/ % % % % % True Found Rec Found Rec Found Rec Conc Found Rec Found Rec Element Mass 261.47 104.6 189.042 250.0 Arsenic

657 611

## Metals Data Reporting Form

Continuing Calibration Verification Units: ug/L **ICPST** Instrument: Acceptable Range: 90% - 110% Chart Number: T00517A.ARC Inorganic Ventures Standard ID: 0014-084-5 Standard Source: CCV3-1 CCV3-2 5/17/00 5/17/00 9:40 AM 9:53 AM WL/ % % % % True Rec Found Rec Found Rec Found **Found** Found Rec Rec Element Mass Conc 507.31 101.5 517.33 **103.5** 189.042 500.0 Arsenic

657 612

**Version 3.63.3** 

# Metals Data Reporting Form

Initial Cali	ibration B	lank R	esults									
Instrument	:ICP	ST	_				Units:		ug/L			
Chart Num	ber: <u>T00</u>	517A.A	RC_									
Standard S	ource:						Standar	rd ID:				_
			ICB1 5/17/0 8:52 A	00								
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	Q	Found	Q	Found	Q
Amannia	100.042	1 10	2.6	: TT	1				l			

U Result is less than the MDL

Form 3 Equivalent B Result is between MDL and RL

5010

# Metals Data Reporting Form

Continuing	g Calibra	tion B	lank Resu	ılts								
Instrument	: <u>IC</u>	PST					Units:		ug/L			
Chart Num	ber: <u>T0</u>	0517A.A	ARC									
Standard S	ource:						Standa	ard ID	);		···	
			CCB1 5/17/00 9:44 AM		CCB2 5/17/00 9:57 AN	)						
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	Q	Found	Q	Found	Q
Arsenic	189.042	10	2.6	υ	2.6	U						

U Result is less than the MDL

Form 3 Equivalent B Result is between MDL and RL

#### Metals Data Reporting Form

Preparation Blank Results

Lab Sample ID: DD8AMB

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	189.042	0.26	1.0	0.26	Ŭ	I	ICPST	5/17/00	9:07

657 615

#### Metals Data Reporting Form

Interference Check Standard A **Instrument:** ICPST Units: ug/L Chart Number: T00517A.ARC Acceptable Range: 0% - 0% Standard Source: Inorganic Ventures Standard ID: 0014-088-12 ICSA 5/17/00 8:56 AM WL/ Reporting True Element Limit Mass Conc Found Found Found Found Found Arsenic 189.042 10 3(u)

MW 5-22-00

657 616

#### Metals Data Reporting Form

Interference Check Standard AB

Instrument: ICPST Units: ug/L

Chart Number: T00517A.ARC Acceptable Range: 80% - 120%

Standard Source: Inorganic Ventures Standard ID: 0014-075-12

			ICSA 5/17/0 9:01 A	00								-
Element	WL/ Mass	True Conc	Found	% Rec	Found	% Rec	Found	% Rec	Found	% Rec	Found	% Rec
Arsenic	189.042	1000	1051.0	105.1								

#### Metals Data Reporting Form

Matrix Spike Sample Results

Spike Sample ID: DD6A4S

Original Sample ID: DD6A4 Client ID: DF/S1/0133/SDC/001 AS

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 10.39

Element	WL/ Mass	OS Conc	Q	MS Conc	Q	Spike Level	% Rec	OS DF	MS DF	Instr	OS Anal Date	OS Anal Time	MS Anal Date	MS Anal Time
Arsenic	189.0	7.7		216		223.19	93.1	1	1	ICPST	5/17/00	9:15	5/17/00	9:24

Comments: Lot #: C0E130194 Sample #: 1

U Result is less than the MDL

Form 5A Equivalent

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

<sup>\*</sup> Duplicate analysis RPD was not within limits

657 618

#### Metals Data Reporting Form

Matrix Spike Duplicate Sample Results

Spike Sample ID: DD6A4D

Original Sample ID: DD6A4 Client ID: DF/S1/0133/SDC/001 AD

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 10.39

Element	WL/ Mass	OS Conc	Q	MSD Conc	Q	Spike Level	% Rec	OS DF	MSD DF	Instr	OS Anal Date	OS Anal Time	MSD Anal Date	MSD Anal Time
Arsenic	189.0	7.7		217		223.19	93.7	1	1	ICPST	5/17/00	9:15	5/17/00	9:28

Comments: Lot #: C0E130194 Sample #: 1

Version 3.63.3

Form 5A Equivalent

U Result is less than the MDL

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

<sup>\*</sup> Duplicate analysis RPD was not within limits

# Metals Data Reporting Form

Matrix Spike Duplicate RPD Report

Matrix Spike Duplicate Sample ID: DD6A4D

Matrix Spike Sample ID: DD6A4S Client ID: DF/S1/0133/SDC/001 AD

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: 10.39

Element	WL/ Mass	MS Cone	Q	MSD Cone	Q	RPD	MS DF	MSD DF	Instr	MS Anal Date	MS Anal Time	MSD Anal Date	MSD Anal Time
Arsenic	189.042	216		217		0.6%	1	1	ICPST	5/17/00	9:24	5/17/00	9:28

Comments: Lot #: C0E130194 Sample #: 1

- B Result is between MDL and RL
- N Spike recovery failed
- NC Percent recovery was not calculated

#### Metals Data Reporting Form

Laboratory Control Sample Results

Lab Sample ID: DD8AMC

Matrix: Soil Units: mg/kg Prep Date: 5/16/00 Prep Batch: 0137277

Weight: 1.0 Volume: 100 Percent Moisture: NA

Element	WL/ Mass	Spike Level	Conc	Percent Recovery	Q	Range	DF	Instr	Anal Date	Anal Time
Arsenic	189.042	200	198	98.7		80-120	1	ICPST	5/17/00	9:11

# Metals Data Reporting Form

Serial Dilution RPD Report

Serial Dilu	ition Sampl	e LD:	· · · · · · · · · · · · · · · · · · ·	Ē	D6A4P						<del></del>			
Original S	ample ID:		D:	D6A	4		Client II	):	DF/S	1/0133/5	SDC/001 A	4		
Matrix:	Soil	ı	Units:	m	ıg/kg		Prep Da	te:	5/1	6/00	Pr	ep Bato	<b>:h:</b> 01	37277
Weight: _	1.0	•	Volume:		100		Percent I	Moistı	ıre: _	10.3	9			***
Element	W) Ma	1	OS Conc	Q	Serial Dilution Conc	Q	Percent Diff	OS DF	Ser Dil DF	Instr	OS Anal Date	OS Anal Time	Ser Dil Anal Date	Ser Dil Anal Time
Arsenic	189	.042	7.7		7.8			1	5	ICPST	5/17/00	9:15	5/17/00	9:19

1

ICPST

5/17/00

9:15

5/17/00

9:19

### Metals Data Reporting Form

**Instrument Detection Limits** 

Instrument: ICPST Units: ppb

Element	Wavelength /Mass	Reporting Limit	MDL	Date of MDL
Arsenic	189.04	10	2.6	4/1/00

# Metals Data Reporting Form

Inter-Element Correction Factors

nstrument: _	ICPST_	Date of IEC's: 3/23/00
Interfering Element	Wavelength /Mass	Correction Factor(s)
Aluminum	308.215	Pb(0.000521), Tl(-0.000024)
Aluminum	308.215	Pb(-0.00021)
Chromium	267.716	Sb(0.006657)
Chromium	267.716	As(-0.002866), Sb(0.011574)
Cobalt	228.616	Pb(0.000021), Se(-0.000486)
Cobalt	228.616	Cd(-0.000119), Fe(0.078354), Ni(-0.000459), Pb(-0.000632), Se(0.000542), T1(0.002839)
Iron	271.441	Pb(0.000053), Sb(0.000025), Se(-0.000357)
Iron	271.441	Cd(0.000097), Cr(-0.00001), Pb(0.00011), Sb(0.000021), Se(0.000012), V(-0.000261), Zn(0.000116)
Magnesium	279.078	Fe(-0.00063)
Manganese	257.61	Se(0.000566)
Manganese	257.61	Se(0.000509), Tl(-0.00542)
Molybdenum	202.03	As(-0.000973), Pb(-0.000443), Sb(-0.002114), Se(0.000412)
Molybdenum	202.03	Pb(-0.00095), Sb(-0.009783)
Nickel	231.604	Pb(0.000115)
Nickel	231.604	Pb(0.00033), Sb(-0.001031), Zn(0.004462)
Vanadium	292.402	Pb(-0.000387)
Vanadium	292.402	Al(0.024171), Be(-0.008664), Cr(-0.000104), Fe(0.013874), Pb(0.00004), Sb(-0.007914), Se(0.000373), Tl(0.001639)

# Metals Data Reporting Form

Linear Dynamic Ranges

Instrument: \_ ICPST

Units: ppb

Element	Wavelength	Linear	Date of Linear
	/Mass	Range	Range
Arsenic	189.04	10000	3/15/00

### Metals Data Reporting Form

Preparation Log

Preparation Batch: 0137277 Instrument: ICP Matrix: Soil

Sample ID	Prep Date	Weight (g)	Volume (ml)	% Moisture
DD8AMB	5/16/00	1.0	100	NA
DD8AMC	5/16/00	1.0	100	NA
DD6A4	5/16/00	1.0	100	10.39
DD6A4D	5/16/00	1.0	100	10.39
DD6A4S	5/16/00	1.0	100	10.39
DD6A5	5/16/00	1.0	100	8.64
DD6A6	5/16/00	1.0	100	10.27
DD6A7	5/16/00	1.0	100	10.81

657 626

# Metals Data Reporting Form

Instrument Runlog

Instrument: ICPST

Chart Number: T00517A.ARC

	Date of	Time of
Sample Name	Analysis	Analysis
ZZZZZZ	5/17/00	8:09 AM
ZZZZZZ	5/17/00	8:13 AM
ZZZZZZ	5/17/00	8:17 AM
ZZZZZZ	5/17/00	8:21 AM
STD1	5/17/00	8:36 AM
STD6	5/17/00	8:40 AM
STD7	5/17/00	8:44 AM
ICV3-1	5/17/00	8:48 AM
ICB1	5/17/00	8:52 AM
ICSA	5/17/00	8:56 AM
ICSAB	5/17/00	9:01 AM
DD8AMB	5/17/00	9:07 AM
DD8AMC	5/17/00	9:11 AM
DD6A4	5/17/00	9:15 AM
DD6A4P	5/17/00	9:19 AM
DD6A4S	5/17/00	9:24 AM
DD6A4D	5/17/00	9:28 AM
DD6A5	5/17/00	9:32 AM
DD6A6	5/17/00	9:36 AM
CCV3-1	5/17/00	9:40 AM
CCB1	5/17/00	9:44 AM
DD6A7	5/17/00	9:48 AM
CCV3-2	5/17/00	9:53 AM
CCB2	5/17/00	9:57 AM
ZZZZZZ	5/17/00	10:01 AM
ZZZZZZ	5/17/00	10:05 AM
ZZZZZZ	5/17/00	10:09 AM
ZZZZZZ	5/17/00	10:13 AM
ZZZZZZ	5/17/00	10:18 AM
ZZZZZZ	5/17/00	10:22 AM
ZZZZZZ	5/17/00	10·26 AM
ZZZZZZ	5/17/00	10:30 AM
ZZZZZZ	5/17/00	10:34 AM
ZZZZZZ	5/17/00	10:38 AM
ZZZZZZ	5/17/00	10:42 AM
ZZZZZZ	5/17/00	10:47 AM
ZZZZZZ	5/17/00	10:51 AM
ZZZZZZ	5/17/00	10:55 AM
ZZZZZZ	5/17/00	10:59 AM
ZZZZZZ	5/17/00	11:03 AM
ZZZZZZ	5/17/00	11:07 AM

# Metals Data Reporting Form

Instrument Runlog

Instrument: ICPST

Chart Number: T00517A.ARC

	Date of	Time of		
Sample Name	Analysis	Analysis		
ZZZZZZ	5/17/00	11:12 AM		
ZZZZZZ	5/17/00	11:16 AM		
ZZZZZZ	5/17/00	11:20 AM		
ZZZZZZ	5/17/00	11:24 AM		
ZZZZZZ	5/17/00	11:28 AM		
ZZZZZZ	5/17/00	11:32 AM		
ZZZZZZ	5/17/00	11:36 AM		
ZZZZZZ	5/17/00	11:41 AM		
ZZZZZZ	5/17/00	11:45 AM		
ZZZZZZ	5/17/00	11:49 AM		
ZZZZZZ	5/17/00	11:53 AM		
ZZZZZZ	5/17/00	11:57 AM		
ZZZZZZ	5/17/00	12:01 PM		
ZZZZZZ	5/17/00	12:06 PM		
ZZZZZZ	5/17/00	12:10 PM		
ZZZZZZ	5/17/00	12:14 PM		
ZZZZZZ	5/17/00	12:18 PM		
ZZZZZZ	5/17/00	12:22 PM		
ZZZZZZ	5/17/00	12:26 PM		
ZZZZZZ	5/17/00	12:31 PM		
ZZZZZZ	5/17/00	12:35 PM		
ZZZZZZ	5/17/00	12·39 PM		
ZZZZZZ	5/17/00	12:43 PM		
ZZZZZZ	5/17/00	12:47 PM		
ZZZZZZ	5/17/00	12:51 PM		
ZZZZZZ	5/17/00	12:56 PM		
ZZZZZZ	5/17/00	1:00 PM		
ZZZZZZ	5/17/00	1:04 PM		
ZZZZZZ	5/17/00	1:08 PM		
ZZZZZZ	5/17/00	1:12 PM		
ZZZZZZ	5/17/00	1:16 PM		
ZZZZZZ	5/17/00	1:20 PM		
ZZZZZZ	5/17/00	1:25 PM		
ZZZZZZ	5/17/00	1:29 PM		
ZZZZZZ	5/17/00	1:33 PM		
ZZZZZZ	5/17/00	1:37 PM		
ZZZZZZ	5/17/00	1:41 PM		
ZZZZZZ	5/17/00	1:45 PM		
ZZZZZZ	5/17/00	1:50 PM		
ZZZZZZ	5/17/00	1:54 PM		
ZZZZZZ	5/17/00	1:58 PM		

#### Metals Data Reporting Form

Instrument Runlog

Instrument: ICPST

Chart Number: T00517A.ARC

Sample Name	Date of Analysis	Time of Analysis	
ZZZZZZ	5/17/00	2:02 PM	
ZZZZZZ	5/17/00	2:06 PM	

.

#### METALS RAW DATA

Analysis Report (ONC) Averages

05/17/00 02:11:48 PM page 3

1	STD1	00331
2	STD6	5.34553
3 4	STD7 ICV3-1 0014-079-6	.26147
5	ICB1	00053
6 7	ICSA 0014-088-12 ICSAB 0014-075-12	00138 1.0510
8	DD8AMB	.00018
9	DD8AMC	1.9749
	DD6A4 DD6A4P5	.06865 .01392
11 12	DD6A45	1.9310
13	DD6A4D	1.9422
	DD6A5 DD6A6	.06566 .07272
	CCV3-1 0014-084-5	.50731
17	CCB1	00043
18	DD6A7	.06883 .51733
	CCV3-2 CCB2	.00027
21		00040
	DDOLEC	1.9427 .00149
	DCQF8 DCQFL	.00090
	DCQFR	.00157
	DCQFX	.00099 .00182
	DCQG5 DCQGK	00083
	DCQGM	.01887
	DCQGN	.00481 .51603
	CCV3-3 CCB3	00182
	DCQGQ	.00260
	DCQGV	.01129 .00052
35 36		.00121
37		.00005
38		.00003 .00021
39 40		00001
41		2.0957
42		2.1676 .50734
	CCV3-4 CCB4	00006
45	DCVMK	.00057
46		.00151 .00106
47 48		.00108
49	DCVMP	.00039
	DD8DGBF	.00019 2.0111
	DD8DGCF DCQF8F	.00063
	DCQFLF	.00130

Ave	r	a	g	e
-----	---	---	---	---

-.00069

#	Sample Name	AS
54	DCQFRF	00057
55	CCV3-5	.52236
56	CCB5	00031
57	DCQFXF	00026
58	DCQG5F	.00201
59	DCQGKF	.00072
60	DCQGMF	.01650
61	DCQGNF	.00446
62	DCQGQF	.00337
63	DCQGVF	.00743
64	DCVMDF	00067
65	DCVMEF	.00055
66	DCVMFF	00029
67	CCV3-6	.50773
68	CCB6	.00003
69	DCVMGF	00042
70	DCVMHF	.00068
71	DCVMHP5F	00057
72	DCVMHSF	2.1854
73	_	2.1046
74	DCVMKF	.00066
75		.00034
76	DCVMMF	.00055
77	DCVMNF	.00074
78	DCVMPF	.00123
79	CCV3 - 7	.51537

80 CCB7

Analysis Report Summary

05/17/00 02:11:48 PM page 1 B66.700

				4071	$\mathcal{U}$			
#	Sample Name	File	Method	Date	Time	OpID	Type	Mode
1	STD1	T00517A	METTRA	05/17/00	08:36		X	IR
2	STD6	T00517A	METTRA	05/17/00	08:40		X	IR
3	STD7	T00517A	METTRA	05/17/00	08:44		X	IR
4	ICV3-1 0014-079-6	T00517A	METTRA	05/17/00	08:48	RJG	S	CONC
5	ICB1	T00517A	METTRA	05/17/00	08:52		S	CONC
6	ICSA 0014-088-12	T00517A	METTRA	05/17/00	08:56		Q	CONC
7	ICSAB 0014-075-12	T00517A	METTRA	05/17/00	09:01		Q	CONC
8	DD8AMB	T00517A	METTRA	05/17/00	09:07		ŝ	CONC
9	DD8AMC	T00517A	METTRA	05/17/00	09:11		S	CONC
_	DD6A4	T00517A	METTRA	05/17/00	09:15		S	CONC
	DD6A4P5	T00517A	METTRA	05/17/00	09:19		S	CONC
		T00517A	METTRA	05/17/00	09:24		s	CONC
	DD6A4S	T00517A	METTRA	05/17/00	09:28		s	CONC
	DD6A4D	T00517A	METTRA	05/17/00	09:32		s	CONC
	DD6A5	T00517A	METTRA	05/17/00	09:36		Š	CONC
	DD6A6		METTRA	05/17/00	09:40		S	CONC
	CCV3-1 0014-084-5	T00517A	METTRA	05/17/00	09:44		s	CONC
	CCB1	T00517A	METTRA	05/17/00	09:48		S	CONC
	DD6A7	T00517A	METTRA	05/17/00	09:53		S	CONC
	CCV3-2	T00517A	METTRA	05/17/00	09:57		s	CONC
	CCB2	T00517A		05/17/00	10:01		s	CONC
	DD0L6B	T00517A	METTRA		10:01		S	CONC
	DD0L6C	T00517A	METTRA	05/17/00	10:03		S	CONC
	DCQF8	T00517A	METTRA	05/17/00			S	CONC
	DCQFL	T00517A	METTRA	05/17/00	10:13			
	DCQFR	T00517A	METTRA	05/17/00	10:18		S	CONC
	DCQFX	T00517A	METTRA	05/17/00	10:22		S	CONC
	DCQG5	T00517A	METTRA	05/17/00	10:26		S	CONC
	DCQGK	T00517A	METTRA	05/17/00	10:30		S	CONC
	DCQGM	T00517A	METTRA	05/17/00	10:34		S	CONC
	DCQGN	T00517A	METTRA	05/17/00	10:38		S	CONC
	CCV3-3	T00517A	METTRA	05/17/00	10:42		S	CONC
	CCB3	T00517A	METTRA	05/17/00	10:47		S	CONC
	DCQGQ	T00517A	METTRA	05/17/00	10:51		S	CONC
	DCQGV	T00517A	METTRA	05/17/00	10:55		S	CONC
	DCVMD	T00517A	METTRA	05/17/00	10:59		S	CONC
36	DCVME	T00517A	METTRA	05/17/00	11:03		S	CONC
37	DCVMF	T00517A	METTRA	05/17/00	11:07		S	CONC
	DCVMG	T00517A	METTRA	05/17/00	11:12		s	CONC
39	DCVMH	T00517A	METTRA	05/17/00	11:16		S	CONC
40	DCVMHP5	T00517A	METTRA	05/17/00	11:20		S	CONC
41	DCVMHS	T00517A	METTRA	05/17/00	11:24		S	CONC
42	DCVMHD	T00517A	METTRA	05/17/00	11:28		S	CONC
43	CCV3-4	T00517A	METTRA	05/17/00	11:32		S	CONC
44	CCB4	T00517A	METTRA	05/17/00	11:36		S	CONC
45	DCVMK	T00517A	METTRA	05/17/00	11:41		S	CONC
46	DCVML	T00517A	METTRA	05/17/00	11:45		S	CONC
	DCVMM	T00517A	METTRA	05/17/00	11:49		S	CONC
	DCVMN	T00517A	METTRA	05/17/00	11:53		S	CONC
49		T00517A	METTRA	05/17/00	11:57		S	CONC
	DD8DGBF	T00517A	METTRA	05/17/00	12:01		S	CONC
	DD8DGCF	T00517A	METTRA	05/17/00	12:06		S	CONC
	DCQF8F	T00517A	METTRA	05/17/00	12:10		S	CONC
	DCQFLF	T00517A	METTRA	05/17/00	12:14	RJG	S	CONC

Analysis Report Summary

05/17/00 02:11:48 PM

page 2

		<u>-</u>		, ,				F2.
#	Sample Name	File	Method	Date	Time	OpID	Туре	Mode
	DCQFRF	T00517A	METTRA	05/17/00	12:18		S	CONC
	CCV3-5	T00517A	METTRA	05/17/00	12:22		S	CONC
	CCB5	T00517A	METTRA	05/17/00	12:26	RJG	S	CONC
57	DCQFXF	T00517A	METTRA	05/17/00	12:31		S	CONC
58	DCQG5F	T00517A	METTRA	05/17/00	12:35		S	CONC
59	DCQGKF	T00517A	METTRA	05/17/00	12:39	RJG	S	CONC
60	DCQGMF	T00517A	METTRA	05/17/00	12:43	RJG	S	CONC
61	DCQGNF	T00517A	METTRA	05/17/00	12:47	RJG	S	CONC
62	DCQGQF	T00517A	METTRA	05/17/00	12:51	RJG	S	CONC
63	DCQGVF	T00517A	METTRA	05/17/00	12:56	RJG	S	CONC
64	DCVMDF	T00517A	METTRA	05/17/00	13:00	RJG	S	CONC
65	DCVMEF	T00517A	METTRA	05/17/00	13:04	RJG	S	CONC
66	DCVMFF	T00517A	METTRA	05/17/00	13:08	RJG	S	CONC
67	CCV3-6	T00517A	METTRA	05/17/00	13:12	RJG	S	CONC
68	CCB6	T00517A	METTRA	05/17/00	13:16	RJG	S	CONC
69	DCVMGF	T00517A	METTRA	05/17/00	13:20	RJG	S	CONC
70	DCVMHF	T00517A	METTRA	05/17/00	13:25	RJG	S	CONC
71	DCVMHP5F	T00517A	METTRA	05/17/00	13:29	RJG	S	CONC
72	DCVMHSF	T00517A	METTRA	05/17/00	13:33	RJG	S	CONC
73	DCVMHDF	T00517A	METTRA	05/17/00	13:37	RJG	S	CONC
74	DCVMKF	T00517A	METTRA	05/17/00	13:41	RJG	S	CONC
75	DCVMLF	T00517A	METTRA	05/17/00	13:45	RJG	S	CONC
76	DCVMMF	T00517A	METTRA	05/17/00	13:50	RJG	S	CONC
77	DCVMNF	T00517A	METTRA	05/17/00	13:54	RJG	S	CONC
78	DCVMPF	T00517A	METTRA	05/17/00	13:58	RJG	S	CONC
79	CCV3-7	T00517A	METTRA	05/17/00	14:02	RJG	S	CONC
80	CCB7	T00517A	METTRA	05/17/00	14:06	RJG	S	CONC

Standard: STD1 Method: METTRA Standar Run Time: 05/17/00 08:36:32

Elem	AG	AL	AS	BA	BE	CA	CD
Avge	.00072	.07210	00331	.00088	03461	.00345	.00163
SDev	.00053	.00156	.00132	.00023	.00006	.00001	.00074
%RSD	74.177	2.1635	39.822	26.526	.17853	.29131	45.448
#1	.00109	.07321	00425	.00072	03465	.00344	.00215
#2	.00034	.07100	00238	.00105	03456	.00346	.00110
Elem	CO	CR	CU	FE	MG	MN	MO
Avge	00011	.00168	.00526	00063	.00054	.00234	.00157
SDev	.00016	.00083	.00034	.00001	.00045	.00014	.00019
%RSD	141.42	49.568	6.4064	.87747	82.457	6.0229	11.986
#1	00023	.00109	.00502	00063	.00086	.00224	.00143
#2	.00000	.00227	.00550	00062	.00023	.00244	.00170
Elem	NI	PB/1	PB/2	SB/1	SB/2	SE/1	SE/2
Avge	.00023	.04316	00680	02575	.01080	07401	.06504
SDev	.00016	.00566	.00905	.00389	.00164	.00081	.00432
%RSD	70.051	13.115	133.07	15.090	15.178	1.0954	6.6418
#1	.00011	.04716	00040	02301	.00964	07458	.06810
#2		.03915	01320	02850	.01196	07344	.06199
Elem	TL	<b>*</b> *					
Avge SDev %RSD	04198 .01073 25.558	V .00000 .00000 .00000	ZN 00014 .00008 55.830				

657 635

05/17/00 08:40:38 AM

Standardization Rpt.

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Wavlen	371.030						
Avge	17539			<b></b> +-			<b>~</b> =
SDev	153.9021		<b></b>				
%RSD	.8774785		<b>-</b> -	<del>-</del> -		<b></b>	
#1	17430						
#2	17648						

page 1

Method: METTRA Standard: STD6 Run Time: 05/17/00 08:40:43 AS Elem AG CD PB/1 PB/2 SB/1 SB/2 10.572 5.3455 15.432 Avqe 6.5911 8.4785 9.9025 6.3469 .0010 .000 SDev .060 .0020 .0097 .0099 .0002 %RSD .56931 .01836 .00313 .03057 .11386 .09966 .00354 #1 10.529 5.3462 15.432 6.5897 8.4853 9.9095 6.3468 #2 5.3448 15.433 10.614 6.5926 8.4717 9.8955 6.3471 Elem SE/1 SE/2 TL4.6925 4.8374 4.9331 Avge .0067 .0058 .0131 SDev %RSD .13852 .11688 .27858 #1 4.8421 4.9372 4.7018 #2 4.8326 4.9290 4.6833 3 7 IntStd 4 5 6 Mode Counts NOTUSED NOTUSED NOTUSED NOTUSED NOTUSED NOTUSED Elem Y Wavlen 371.030 \_ \_ - -Avge 17569 13.08147 - -SDev - -\_ \_ - -∜RSD .0744569 \_ \_ #1 17560 #2 17578 \_\_\_ \_ \_

05/17/00 08:48:29 AM

							rage .
Method: Run Time	METTRA : 05/17/00		d: STD7	004-021	.9		
Elem Avge SDev %RSD	AL 6.7609 .0134 .19878	BA 12.752 .010 .07526	BE 12.996 .009 .06836	CA 5.8933 .0006 .01066	CO 3.4503 .0003 .00977	CR 14.061 .004 .03162	CU 3.0190 .0040 .13282
#1 #2	6.7704 6.7514	12.759 12.745	12.990 13.002	5.8928 5.8937	3.4501 3.4506	14.058 14.065	3.0218 3.0162
Elem Avge SDev %RSD	FE 4.6489 .0010 .02237	MG 15.174 .002 .01527	MN 11.767 .003 .02263	MO 2.2912 .0125 .54468	NI 2.8219 .0008 .02681	V .92300 .00027 .02967	ZN 3.1068 .0013 .04121
#1 #2	4.6482 4.6497	15.176 15.172	11.769 11.765	2.2824 2.3000	2.8224 2.8213	.92281 .92320	3.1077 3.1059
IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 17237 20.43566 .1185556	2 NOTUSED	3 NOTUSED   	4 NOTUSED   	5 NOTUSED   	6 NOTUSED    	7 NOTUSED    
#1 #2	17252 17223			- <del>-</del> -			

Standardization Report 05/17/00 08:48:30 AM page 1

Method: METTRA Slope = Conc(SIR)/IR

	_					
Element		High std	Low std	Slope	Y-intercept	Date Standardized
AG	328.068	STD6	STD1	.189197	000135	05/17/00 08:44:53
AL	308.215	STD7	STD1	7.48962	540022	05/17/00 08:44:53
AS	189.042	STD6	STD1	.186956	.000619	05/17/00 08:44:53
BA	493.409	STD7	STD1	.313691	000277	05/17/00 08:44:53
BE	313.042	STD7	STD1	.304307	.010532	05/17/00 08:44:53
CA	317.933	STD7	STD1	16.9784	058565	05/17/00 08:44:53
CD	226.502	STD6	STD1	.064805	000106	05/17/00 08:44:53
CO	228.616	STD7	STD1	1.15927	.000133	05/17/00 08:44:53
CR	267.716	STD7	STD1	.284435	000477	05/17/00 08:44:53
CU	324.753	STD7	STD1	1.32726	<b>-</b> .006979	05/17/00 08:44:53
FE	271.441	STD7	STD1	10.8195	.006786	05/17/00 08:44:53
MG	279.078	STD7	STD1	6.59043	003583	05/17/00 08:44:53
MN	257,610	STD7	STD1	.340002	000795	05/17/00 08:44:53
MO	202.030	STD7	STD1	1.74699	002738	05/17/00 08:44:53
NI	231.604	STD7	STD1	1.41697	000322	05/17/00 08:44:53
PB/1	220.351	STD6	STD1	.152719	006591	05/17/00 08:44:53
PB/2	220.352	STD6	STD1	.117851	.000802	05/17/00 08:44:53
PB	220.353	NONE	NONE	.000000	.000000	*NOT STANDARDIZED
SB/1	206.831	STD6	STD1	.100723	.002594	05/17/00 08:44:53
SB/2	206.832	STD6	STD1	.157825	001704	05/17/00 08:44:53
SB	220.353	NONE	NONE	.000000	.000000	*NOT STANDARDIZED
SE/1	196.021	STD6	STD1	.203609	.015069	05/17/00 08:44:53
SE/2	196.022	STD6	STD1 .	.205420	013362	05/17/00 08:44:53
SE, Z	220.353	NONE	NONE	.000000	.000000	*NOT STANDARDIZED
TL	190.864	STD6	STD1	.422430	.017734	05/17/00 08:44:53
v	292.402	STD7	STD1	4.31954	.000000	05/17/00 08:44:53
$z\overline{ ext{N}}$	213.856	STD7	STD1	1.29505	.000184	05/17/00 08:44:53
en F.A	223.030		~ ~ ~ ~			,,

5036

page 1

Method: M Run Time: Comment: Mode: COM	05/17/00 STL PITTS		me: ICV3-1 METALS ANA			erator: RJO	: Waxa
Elem	AG	AL	AS	BA	BE	CA	CD 5(700)
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.50563	12.321	.26148	1.0206	1.0055	25.700	.25295
SDev	.00007	.007	.00202	.0014	.0018	.050	.00099
%RSD	.01405	.06047	.77420	.13641	.17922	.19522	.39071
#1	.50568	12.316	.26005	1.0196	1.0068	25.735	.25365
#2	.50557	12.327	.26291	1.0216	1.0042	25.664	.25225
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.55000	13.750	.27500	1.1000	1.1000	27.500	.27500
Low	.45000	11.250	.22500	.90000	.90000	22.500	.22500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.0313	1.0380	1.0007	12.806	24.875	1.0083	1.0242
SDev	.0014	.0012	.0005	.011	.028	.0002	.0015
%RSD	.13537	.11452	.05429	.08878	.11316	.01640	.14606
#1	1.0322	1.0388	1.0011	12.814	24.895	1.0081	1.0253
#2	1.0303	1.0372		12.798	24.855	1.0084	1.0232
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	1.1000	1.1000	13.750	27.500	1.1000	1.1000
Low	.90000	.90000	.90000	11.250	22.500	.90000	.90000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.0322	.25330	.25395	.25374	.26198	.25680	.25852
SDev	.0054	.00203	.00373	.00181	.00357	.00036	.00143
%RSD	.51870	.80332	1.4700	.71428	1.3626	.14162	.55364
#1	1.0360	.25186	.25659	.25502	.26450	.25705	.25953
#2	1.0284	.25474	.25131	.25245	.25945	.25654	.25751
Errors High Low	LC Pass 1.1000 .90000	NOCHECK	NOCHECK	LC Pass .27500 .22500	NOCHECK	NOCHECK	LC Pass .27500 .22500
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.25932	.26259	.26150	.51692	1.0215	1.0359	
SDev	.00230	.00214	.00219	.00409	.0050	.0039	
%RSD	.88528	.81395	.83750	.79134	.49062	.37943	
#1	.26094	.26410	.26305	.51981	1.0251	1.0387	
#2	.25770	.26108	.25995	.51403	1.0180	1.0331	
Errors High Low	NOCHECK	NOCHECK	LC Pass .27500 .22500	LC Pass .55000 .45000	LC Pass 1.1000 .90000	LC Pass 1.1000 .90000	

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 17515 47.16485 .2692849	2 NOTUSED   	3 NOTUSED	4 NOTUSED    	5 NOTUSED	6 NOTUSED   	7 NOTUSED    
#1 #2	17482 17548			- <del>-</del> -		₩ =	
#2	17548			<del></del>		<b></b>	

Method: METTRA Sample Name: ICB1 Operator: RJG

Run Time: 05/17/00 08:52:44

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00004	.01184	00054	.00047	.00059	.00819	.00012
SDev	.00036	.01314	.00013	.00005	.00002	.00102	.00003
%RSD	840.14	110.93	23.235	10.251	2.5674	12.458	25.758
#1	.00021	.02113	00063	.00050	.00058	.00747	.00010
#2	00030	.00255	00045	.00044	.00060	.00891	.00015
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.01000	.20000	.00500	5.0000	.00500
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem Units Avge SDev %RSD	CO ppm .00013 .00000	CR ppm .00018 .00015 84.927	CU ppm 00019 .00015 79.295	FE ppm .00618 .00436 70.560	MG ppm .00788 .00139 17.619	MN ppm .00037 .00003 9.0935	MO ppm .00374 .00116 31.050
#1 #2	.00013	.00029 .00007	00030 00008	.00926 .00309	.00886 .00689	.00039	.00456 .00292
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.05000	.01000	.02500	.10000	5.0000	.01500	.04000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00016	00038	.00107	.00059	.00119	.00051	.00074
SDev	.00023	.00102	.00081	.00088	.00141	.00142	.00048
%RSD	143.25	270.73	75.137	148.19	118.07	279.75	65.015
#1	.00000	.00034	.00165	.00121	.00219	00050	.00040
#2	00032	00109	.00050	00003	.00020	.00151	.00108
Errors High Low	LC Pass .04000 04000	NOCHECK	NOCHECK	LC Pass .00300 00300	NOCHECK	NOCHECK	LC Pass .06000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00290	.00082	00042	.00283	.00000	.00070	
SDev	.00375	.00042	.00097	.00226	.00000	.00005	
%RSD	129.26	50.415	233.30	79.711	70.560	7.2821	
#1 #2	00555 00025	.00112	00110 .00027	.00123	.00000	.00066 .00074	
Errors High Low	NOCHECK	NOCHECK	LC Pass .00500 00500	LC Pass .01000 01000	LC Pass .05000 05000	LC Pass .02000 02000	

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	Y						
Wavlen	371.030					<b>-</b>	
Avge	17549		<b></b>	<del></del>			
SDev	91.49989						
%RSD	.5213981					<b>→</b> =	
#1	17484						
#2	17614						

05/17/00 09:00:59 AM

Method: METTRA Sample Name: ICSA 0014-088-12 Operator: RJG Run Time: 05/17/00 08:56:53

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00035	520:05	00138	.00172	00076	491.09	00204
SDev	.00018	2.05	.00142	.00003	.00008	.90	.00043
%RSD	52.634	.39335	102.93	1.8192	9.9415	.18227	21.013
#1	00022	518.60	00239	.00174	00071	491.72	00234
#2	00048	521.49	00038	.00170	00081	490.46	00173
Errors Value Range	NOCHECK	QC Pass 500.00 20.000	NOCHECK	NOCHECK	NOCHECK	QC Pass 500.00 20.000	NOCHECK
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00048	.00303	00290	201.49	526.39	.00733	00030
SDev	.00030	.00001	.00034	.27	.24	.00006	.00023
%RSD	63.564	.41888	11.816	.13390	.04546	.81924	76.431
#1	00026	.00304	00314	201.30	526.22	.00737	00045
#2	00069		00265	201.68	526.56	.00729	00014
Errors Value Range	NOCHECK	NOCHECK	NOCHECK	QC Pass 200.00 20.000	QC Pass 500.00 20.000	NOCHECK	NOCHECK
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00148	00238	00201	00214	.00269	.00051	.00123
SDev	.00106	.00411	.00233	.00018	.00078	.00765	.00485
%RSD	71.578	172.61	115.86	8.6090	29.063	1499.5	392.42
#1	.00223	.00053	00366	00227	.00213	.00592	.00466
#2	.00073	00529	00036	00201	.00324	00490	00219
Errors Value Range	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00112	.00160	.00069	.00438	00043	.00099	
SDev	.00031	.00352	.00224	.00131	.00192	.00010	
%RSD	27.397	219.92	324.03	29.924	440.95	9.9890	
#1	00091	00089	00089	.00530	.00092	.00092	
#2	00134	.00408	.00228	.00345	00179	.00106	
Errors Value Range	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK	

65 Analysis	7 644 Report	QC Sta	ndard	05/17	/00 09:00:	59 AM	page 2
IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16092 30.15838 .1874079	2 NOTUSED   	3 NOTUSED   	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1 #2	16071 16114		<del></del>		 		een aan Ann ma

05/17/00 09:05:09 AM

Method: METTRA Sample Name: ICSAB 0014-075-12 Operator: RJG

Run Time: 05/17/00 09:01:03

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

		•					
Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.0755	517.96	1.0510	.52058	.49396	489.21	.93824
SDev	.0010	.14	.0022	.00082	.00112	1.13	.00271
%RSD	.08839	.02665	.21325	.15804	.22683	.23186	.28936
#1	1.0761	518.05	1.0526	.52116	.49475	490.01	.94016
#2	1.0748	517.86	1.0495	.52000	.49317	488.41	.93632
Errors	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass
Value	1.0000	500.00	1.0000	.50000	.50000	500.00	1.0000
Range	20.000	20.000	20.000	20.000	20.000	20.000	20.000
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.49381	.50610	.52961	201.42	523.27	.50548	1.0099
SDev	.00192	.00106	.00155	.23	.60	.00065	.0029
%RSD	.38893	.21007	.29184	.11363	.11461	.12945	.29065
#1	.49517	.50686	.53070	201.58	523.70	.50595	1.0078
#2	.49245	.50535	.52851	201.25	522.85	.50502	1.0120
Errors	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass
Value	.50000	.50000	.50000	200.00	500.00	.50000	1.0000
Range	20.000	20.000	20.000	20.000	20.000	20.000	20.000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.98333	.99402	.97830	.98353	1.0660	1.0444	1.0515
SDev	.00576	.00131	.00457	.00348	.0137	.0021	.0059
%RSD	.58603	.13153	.46742	.35437	1.2825	.20002	.56543
#1	.98740	.99494	.98153	.98600	1.0756	1.0459	1.0557
#2	.97925	.99309	.97507	.98107	1.0563	1.0430	1.0473
Errors Value Range	QC Pass 1.0000 20.000	NOCHECK	NOCHECK	QC Pass 1.0000 20.000	NOCHECK	NOCHECK	QC Pass 1.0000 20.000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	1.0388	1.0655	1.0566	1.0453	.50972	1.0499	
SDev	.0033	.0111	.0063	.0049	.00029	.0042	
%RSD	.31640	1.0404	.59619	.47298	.05646	.39710	
#1	1.0364	1.0733	1.0611	1.0488	.50992	1.0529	
#2	1.0411	1.0577	1.0521	1.0418	.50952	1.0470	
Errors Value Range	NOCHECK	NOCHECK	QC Pass 1.0000 20.000	QC Pass 1.0000 20.000	QC Pass .50000 20.000	QC Pass 1.0000 20.000	

657 Analysis	646 Report	QC Star	ndard	05/17	page 2		
IntStd	1	2	3	4	5	6	7
Mode	Counts	NOTUSED	NOTUSED	NOTUSED	NOTUSED	NOTUSED	NOTUSED
Elem	Y						
Wavlen	371.030						
Avge	16157					<b></b>	
SDev	8.060603		<b>-</b> -				
%RSD	.0498891	<del>-</del> -					
#1	16151	~ <b>~</b>		~		<b>**</b> ••	
<u>#</u> 2	16163						

Method: METTRA Sample Name: DD8AMB Operator: RJG

Run Time: 05/17/00 09:07:29

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00015	.02040	.00018	.00036	.00027	.11647	.00004
SDev	.00032	.00451	.00092	.00002	.00006	.00082	.00021
%RSD	219.51	22.102	503.44	4.0687	22.920	.70217	586.10
#1 #2	.00037 00008	.02358 .01721	.00083 00047	.00037	.00032	.11705 .11590	00011 .00018
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.01000	.20000	.00500	5.0000	.00500
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00186	.00013	.00032	.03921	.02477	.00136	.00032
SDev	.00011	.00031	.00030	.00152	.00890	.00000	.00047
%RSD	5.8149	228.02	91.747	3.8807	35.914	.11819	147.94
#1	00179	.00035	.00011	.04029	.03107	.00136	.00066
#2	00194	00008		.03814	.01848	.00136	00001
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.05000	.01000	.02500	.10000	5.0000	.01500	.04000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00017	.00318	.00071	.00154	.00012	.00195	.00134
SDev	.00173	.00121	.00177	.00078	.00148	.00466	.00262
%RSD	1036.8	38.061	248.58	50.786	1262.9	239.41	195.62
#1	.00105	.00404	00054	.00098	00093	.00524	.00319
#2	00139		.00197	.00209	.00116	00135	00051
Errors High Low	LC Pass .04000 04000	NOCHECK	NOCHECK	LC Pass .00300 00300	NOCHECK	NOCHECK	LC Pass .06000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00264	.00205	.00049	00106	.00311	.00474	
SDev	.00052	.00265	.00194	.00072	.00015	.00004	
%RSD	19.683	129.35	398.25	67.444	4.9398	.81214	
#1 #2	00300 00227	.00017 .00392	00088 .00186	00056 00157	.00322	.00476 .00471	
Errors High Low	NOCHECK	NOCHECK	LC Pass .00500 00500	LC Pass .01000 01000	LC Pass .05000 05000	LC Pass .02000 02000	

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 17411 124.4149 .7145815	2 NOTUSED    	NOTUSED	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    	
#1	17499		<b></b>					
#2	17323						~ ~	

Method: METTRA Sample Name: DD8AMC Run Time: 05/17/00 09:11:38 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04755	1.8435	1.9749	1.9793	.04989	L.15359	.04887
SDev	.00048	.0228	.0189	.0175	.00032	.00003	.00021
%RSD	1.0158	1.2389	.95519	.88179	.65038	.02095	.42930
#1	.04789	1.8596	1.9883	1.9917	.05012	L.15357	.04902
#2	.04721	1.8273	1.9616	1.9670	.04966	L.15361	.04872
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Low	LC Pass
High	.06000	2.4000	2.4000	2.4000	.06000	60.000	.06000
Low	.04000	1.6000	1.6000	1.6000	.04000	40.000	.04000
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.50055	.20240	.24155	1.0032	L.01035	.48996	L00087
SDev	.00389	.00163	.00292	.0076	.00467	.00411	.00020
%RSD	.77676	.80793	1.2087	.75884	45.149	.83930	23.303
#1	.50330	.20356	.24362	1.0086	L.01365	.49287	L00101
#2	.49781	.20124	.23949	.99781	L.00704	.48705	L00073
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Low	LC Pass	LC Low
High	.60000	.24000	.30000	1.2000	60.000	.60000	1.2000
Low	.40000	.16000	.20000	.80000	40.000	.40000	.80000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.50351	.48250	.48294	.48279	.00085	.00037	L.00053
SDev	.00490	.00475	.00732	.00646	.00009	.00023	.00018
%RSD	.97387	.98494	1.5155	1.3390	10.562	62.184	34.551
#1	.50698	.48586	.48812	.48736	.00091	.00053	L.00066
#2	.50005	.47914	.47776	.47822	.00078	.00021	L.00040
Errors High Low	LC Pass .60000 .40000	NOCHECK	NOCHECK	LC Pass .60000 .40000	NOCHECK	NOCHECK	LC Low .60000 .40000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	1.9493	1.9559	1.9538	1.9770	.49343	.50515	
SDev	.0103	.0175	.0151	.0203	.00581	.00481	
%RSD	.52985	.89354	.77271	1.0245	1.1768	.95243	
#1	1.9566	1.9683	1.9644	1.9913	.49753	.50855	
#2	1.9420	1.9436	1.9431	1.9627	.48932	.50174	
Errors High Low	NOCHECK	NOCHECK	LC Pass 2.4000 1.6000	LC Pass 2.4000 1.6000	LC Pass .60000 .40000	LC Pass .60000 .40000	

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Wavlen Avge SDev	371.030 17286 110.5559						 
%RSD	.6395624		<b></b>				
#1	17208						
#2	17364						

Method: METTRA Sample Name: DD6A4 Run Time: 05/17/00 09:15:47 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00120	140.16	.06866	2.9092	.00487	H1287.8	00065
SDev	.00050	.25	.00308	.0089	.00006	6.9	.00024
%RSD	41.657	.17969	4.4842	.30657	1.2570	.53274	37.092
#1	.00155	140.34	.07083	2.9155	.00482	H1292.7	00081
#2	.00085	139.98	.06648	2.9029	.00491	H1283.0	00048
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem Units Avge SDev %RSD	CO ppm .03398 .00009 .27102	CR ppm .12943 .00096 .74200	CU ppm .14369 .00080 .55317	FE ppm 86.045 .343	MG ppm 15.632 .063 .40145	MN ppm 2.0709 .0078 .37855	MO ppm .00766 .00078 10.143
#1	.03405	.13011	.14425	86.288	15.677	2.0764	.00711
#2		.12875	.14313	85.803	15.588	2.0653	.00821
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.07758	.25495	.25136	.25256	.01560	.01043	.01215
SDev	.00111	.00629	.00285	.00400	.00185	.00136	.00152
%RSD	1.4302	2.4673	1.1342	1.5823	11.885	13.002	12.524
#1	.07836	.25940	.25338	.25538	.01429	.00947	.01108
#2	.07679	.25050	.24935	.24973	.01691	.01139	.01323
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00286	.00475	.00412	.00776	.19260	.39674	
SDev	.00117	.00047	.00070	.00224	.00161	.00197	
%RSD	40.706	9.8567	16.993	28.869	.83583	.49612	
#1 #2	.00204 .00369	.00442 .00508	.00363	.00934 .00617	.19374 .19147	.39813 .39534	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Report

05/17/00 09:19:52 AM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	Y	<b></b>					
Wavlen	371.030					<del></del>	
Avge	17466						
SDev	120.1736				<del></del>	TT 04	
%RSD	.6880245						
#1	17382			<b>→</b> ¬			
#2	17551	- <del>-</del>					

Method: METTRA Sample Name: DD6A4P5 Operator: RJG

Run Time: 05/17/00 09:19:56

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00032	28.700	.01392	.62452	.00120	304.53	00021
SDev	.00007	.159	.00228	.00300	.00009	.84	.00007
%RSD	20.907	.55375	16.396	.48094	7.4890	.27519	34.454
#1	.00037	28.588	.01554	.62239	.00127	303.94	00027
#2	.00027	28.813	.01231	.62664	.00114	305.12	00016
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00608	.02783	.02889	18.648	3.4105	.44267	.00365
SDev	.00004	.00004	.00051	.092	.0154	.00188	.00018
%RSD	.63536	.15640	1.7484	.49231	.45214	.42395	5.0674
#1	.00606	.02779	.02853	18.583	3.3996	.44134	.00352
#2	.00611	.02786	.02925	18.713	3.4214	.44400	.00378
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01608	.05540	.05667	.05624	.00357	.00323	.00334
SDev	.00133	.00182	.00164	.00049	.00005	.00169	.00111
%RSD	8.2517	3.2777	2.8926	.86876	1.3217	52.484	33.333
#1	.01514	.05668	.05551	.05590	.00354	.00442	.00413
#2	.01702	.05412	.05782	.05659	.00360		.00255
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00001	.00122	.00082	.00380	.04382	.08486	
SDev	.00283	.00287	.00097	.00410	.00025	.00033	
%RSD	24794.	234.75	118.54	107.71	.58063	.38684	
#1	.00202	00081	.00013	.00670	.04400	.08510	
#2	00199	.00326	.00151	.00091	.04364	.08463	

05/17/00 09:24:01 AM page 2

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Wavlen	371.030					<b></b>	<del></del>
Avge	17243						
SDev	112.0046	<b></b>	<del>-</del> -				<b>→ →</b>
%RSD	.6495600	· -	<b></b>				
#1	17322						
#2	17164				<b></b>		

Method: METTRA Sample Name: DD6A4S Operator: RJG

Run Time: 05/17/00 09:24:05

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

AG	AL	AS	BA	BE	CA	CD
ppm	ppm	ppm	ppm	ppm	ppm	ppm
.05021	152.11	1.9310	5.1051	.04841	H1447.9	.04070
.00002	.39	.0095	.0119	.00023	4.9	.00006
.03987	.25810	.49099	.23397	.47135	.33643	.13952
.05019	152.38	1.9377	5.1136	.04857	H1451.3	.04066
.05022	151.83	1.9243	5.0967	.04825	H1444.4	.04074
LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
01000	20000	01000	20000	00500	-5.0000	00500
CO	CR	CU	FE	MG	MN	MO
ppm	ppm	ppm	ppm	ppm	ppm	ppm
.45670	.31688	.38190	90.334	14.941	2.4209	.00605
.00100	.00070	.00125	.287	.047	.0060	.00081
.21861	.22161	.32799	.31817	.31216	.24787	13.470
.45740	.31737	.38278	90.537	14.974	2.4252	.00548
.45599	.31638	.38101	90.131	14.908	2.4167	.00663
LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
100.00	20.000	10.000	500.00	600.00	10.000	20.000
05000	01000	02500	10000	-5.0000	01500	04000
NI	PB/1	PB/2	PB	SB/1	SB/2	SB
ppm	ppm	ppm	ppm	ppm	ppm	ppm
.49771	.69388	.68255	.68632	.01570	.00952	.01157
.00200	.00321	.00510	.00447	.00007	.00104	.00072
.40119	.46245	.74747	.65151	.44041	10.974	6.2171
.49912	.69615	.68616	.68948	.01565	.00878	.01107
.49630	.69161	.67894	.68316	.01574	.01026	.01208
LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
SE/1	SE/2	SE	TL	V_	ZN	
ppm	ppm	ppm	ppm	ppm	ppm	
1.8597	1.8806	1.8736	1.8837	.64130	.86654	
.0105	.0096	.0099	.0139	.00256	.00381	
.56671	.51143	.52970	.74052	.39916	.44031	
1.8671	1.8874	1.8807	1.8936	.64311	.86924	
1.8522	1.8738	1.8666	1.8738	.63949	.86384	
NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	
	ppm .05021 .00002 .03987 .05019 .05022 LC Pass 2.000001000 CO ppm .45670 .00100 .21861 .45740 .45599 LC Pass 100.0005000 NI ppm .49771 .00200 .40119 .49912 .49630 LC Pass 100.0004000 SE/1 ppm 1.8597 .0105 .56671 1.8671 1.8522	ppm         ppm           .05021         .39           .03987         .25810           .05019         .52.38           .05022         .51.83           LC Pass         LC Pass           2.0000         600.00          01000        20000           CO         CR           ppm         .95670           .45670         .31688           .00100         .00070           .21861         .22161           .45740         .31737           .45599         .31638           LC Pass         LC Pass           100.00        01000           NI         PB/1           ppm         .49771         .69388           .00200         .00321           .40119         .46245           .49912         .69615           .49630         .69161           LC Pass         NOCHECK           100.00        04000           SE/1         SE/2           ppm         1.8806           .0105         .0096           .56671         .51143           1.8522         1.8738	ppm         ppm         ppm           .05021         152.11         1.9310           .00002         .39         .0095           .03987         .25810         .49099           .05019         152.38         1.9377           .05022         151.83         1.9243           LC Pass         LC Pass           2.0000         600.00         10.000          01000        20000        01000           CO         CR         CU           ppm         ppm         ppm           .45670         .31688         .38190           .00100         .00070         .00125           .21861         .22161         .32799           .45740         .31737         .38278           .45599         .31638         .38101           LC Pass         LC Pass         LC Pass           100.00        01000        02500           NI         PB/1         PB/2           ppm         ppm         .69388         .68255           .00200         .00321         .00510           .49912         .69615         .68616           .49630         .69161         .67894	ppm         ppm         ppm         ppm           .05021         152.11         1.9310         5.1051           .00002         .39         .0095         .0119           .03987         .25810         .49099         .23397           .05019         152.38         1.9377         5.1136           .05022         151.83         1.9243         5.0967           LC Pass         LC Pass         LC Pass           2.0000         600.00         10.000         10.000          01000        20000         10.000         10.000          01000        20000        01000        20000           CO         CR         CU         FE           ppm         ppm         ppm         ppm           .45670         .31688         .38190         90.334           .00100         .00070         .00125         .287           .21861         .22161         .32799         .31817           .45740         .31737         .38278         90.537           .45599         .31638         .38101         90.131           LC Pass         LC Pass         LC Pass           100.00        05000	ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm         ppm <td>  DPM</td>	DPM

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Wavlen	371.030	<b></b>					
Avge	17884						
SDev	27.43630				~ -	100 400	
%RSD	.1534155						
#1	17864		<del></del>				
#2	17903		<b></b> -				

Method: METTRA Sample Name: DD6A4D Operator: RJG

Run Time: 05/17/00 09:28:13

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04889	137.63	1.9422	4.7988	.04909	H1436.9	.04080
SDev	.00018	.09	.0046	.0091	.00006	3.4	.00002
%RSD	.36322	.06573	.23601	.19007	.13224	.23870	.04242
#1	.04877	137.69	1.9455	4.8053	.04914	H1439.4	.04079
#2	.04902	137.56	1.9390	4.7924	.04904	H1434.5	.04081
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.46245	.30433	.37588	79.731	15.011	2.1200	.00536
SDev	.00216	.00049	.00069	.103	.025	.0034	.00110
%RSD	.46768	.15921	.18282	.12874	.16322	.15987	20.588
#1	.46398	.30467	.37637	79.803	15.028	2.1224	.00458
#2	.46092	.30398	.37540	79.658	14.993	2.1176	.00614
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.50168	.68423	.67223	.67623	.01522	.01085	.01231
SDev	.00280	.00276	.00007	.00097	.00120	.00134	.00129
%RSD	.55847	.40371	.01036	.14290	7.8861	12.359	10.516
#1	.50366	.68618	.67228	.67691	.01438	.00990	.01139
#2	.49970	.68228	.67218	.67554	.01607	.01180	.01322
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	1.8877	1.8987	1.8951	1.9084	.62814	.85790	
SDev	.0021	.0064	.0049	.0099	.00157	.00084	
%RSD	.11134	.33457	.26052	.51718	.25035	.09808	
#1	1.8892	1.9032	1.8985	1.9154	.62925	.85850	
#2	1.8862	1.8942	1.8916	1.9014	.62703	.85731	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Report

05/17/00 09:32:18 AM page 2

IntSto Mode Elem Wavler Avge SDev %RSD	Counts Y	2 NOTUSED   	3 NOTUSED   	4 NOTUSED	5 NOTUSED    	6 NOTUSED	7 NOTUSED   
#1	17422						
#2	17506	<b></b>					

Method: METTRA Sample Name: DD6A5 Run Time: 05/17/00 09:32:22 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00391	127.85	.06566	3.1142	.00491	H1435.3	00004
SDev	.00011	.36	.00140	.0074	.00007	3.3	.00003
%RSD	2.8931	.27846	2.1357	.23838	1.4359	.23297	71.273
#1	.00383	127.60	.06665	3.1090	.00496	H1433.0	00006
#2	.00399	128.10	.06467	3.1195	.00486	H1437.7	00002
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.03353	.12691	.14768	79.249	15.439	1.9048	.00612
SDev	.00069	.00025	.00126	.178	.052	.0063	.00017
%RSD	2.0538	.19393	.85207	.22437	.33440	.33259	2.7892
#1	.03305	.12673	.14679	79.123	15.402	1.9003	.00624
#2	.03402	.12708	.14857	79.374	15.475	1.9092	.00600
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.07419	.29287	.29205	.29232	.01797	.00858	.01171
SDev	.00079	.00255	.00024	.00069	.00130	.00230	.00110
%RSD	1.0662	.87192	.08306	.23555	7.2396	26.781	9.3893
#1	.07475	.29107	.29222	.29183	.01705	.01021	.01249
#2	.07363	.29468	.29187	.29281	.01889	.00696	.01093
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.01259	.00682	.00874	.01430	.17856	.40338	
SDev	.00092	.00130	.00056	.00372	.00030	.00012	
%RSD	7.2938	19.013	6.3958	25.980	.16796	.03052	
#1	.01194	.00774	.00914	.01168	.17835	.40330	
#2	.01324	.00590	.00835	.01693	.17877	.40347	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 17250 85.73669 .4970193	2 NOTUSED   	3 NOTUSED    	4 NOTUSED	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1	17311						
#2	17190						~ =

Method: METTRA Sample Name: DD6A6 Run Time: 05/17/00 09:36:31 Method: METTRA Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00172	191.83	.07272	3.0718	.00646	H1105.6	00042
SDev	.00022	1.62	.00101	.0325	.00013	15.6	.00024
%RSD	12.710	.84544	1.3884	1.0575	1.9832	1.4112	56.921
#1	.00156	192.98	.07201	3.0948	.00655	H1116.7	00059
#2	.00187	190.68	.07344	3.0489	.00637	H1094.6	00025
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04377	.25659	.19915	120.63	21.405	2.7404	.00629
SDev	.00072	.00331	.00219	1.43	.269	.0328	.00030
%RSD	1.6468	1.2887	1.0977	1.1884	1.2585	1.1951	4.8117
#1	.04428	.25893	.20070	121.64	21.595	2.7635	.00650
#2	.04326	.25426	.19761	119.61	21.214	2.7172	.00607
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.10824	.36393	.36109	.36204	.01991	.00591	.01058
SDev	.00178	.01097	.00126	.00281	.00673	.00897	.00374
%RSD	1.6457	3.0143	.34791	.77757	33.802	151.74	35.405
#1	.10950	.37169	.36020	.36403	.01515	.01226	.01322
#2	.10698	.35617	.36198	.36004	.02467	00043	.00793
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00538	.00329	.00040	.00877	.27507	.77073	
SDev	.00595	.00447	.00100	.00047	.00388	.01041	
%RSD	110.55	136.19	251.44	5.3899	1.4094	1.3505	
#1	00118	.00012	00031	.00911	.27781	.77809	
#2	00959	.00645	.00111	.00844	.27233	.76337	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 17521 176.7063 1.008543	2 NOTUSED   	3 NOTUSED   	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    	
#1	17396		va. ***		<b>-</b>			
#2	17646							

Sample Name: CCV3-1 0014-084-5 Method: METTRA Operator: RJG

Run Time: 05/17/00 09:40:40

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units Avge	ppm 1.0140	ppm 24.095	ppm .50732	ppm 1.9871	ppm 1.9848	ppm	ppm
SDev %RSD	.0264	.677	.01358	.0550	.0512	49.665 1.206	.48360 .01228
	2.6073	2.8081	2.6763	2.7666	2.5790	2.4290	2.5396
#1 #2	.99535 1.0327	23.617 24.574	.49772 .51692	1.9482 2.0260	1.9486 2.0210	48.812 50.518	.47492 .49228
Errors High Low	LC Pass 1.1000 .90000	LC Pass 27.500 22.500	LC Pass .55000 .45000	LC Pass 2.2000 1.8000	LC Pass 2.2000 1.8000	LC Pass 55.000 45.000	LC Pass .55000 .45000
Elem	CO	CR	CU	FE	MG	MN	MO
Units Avge	ppm 1.9965	ppm 2.0006	ppm 1.9687	ppm 24.707	ppm 48.928	ppm 1.9647	ppm 1.9945
SDev %RSD	.0518 2.5934	.0526 2.6273	.0523 2.6590	.651 2.6331	1.246 2.5476	.0517	.0669
						2.6316	3.3554
#1 #2	1.9599 2.0331	1.9634 2.0378	1.9317 2.0057	24.247 25.167	48.046 49.809	1.9281 2.0012	1.9471 2.0418
Errors High	LC Pass 2.2000	LC Pass 2.2000	LC Pass 2,2000	LC Pass 27.500	LC Pass	LC Pass	LC Pass
Low	1.8000	1.8000	1.8000	22.500	55.000 45.000	2.2000 1.8000	2.2000 1.8000
Elem	NI	PB/1	PB/2	<b>D</b> D	an /-	GD /0	an.
* *		•	-	PB	SB/1	SB/2	SB
Units Avge	ppm 1.9820	ppm .49640	ppm .49721	ppm .49694	SB/1 ppm .51554	58/2 ppm .51085	ppm
	ppm	ppm .49640 .01153	ppm .49721 .01298	ppm .49694 .01250	ppm .51554 .01268	ppm .51085 .01340	ppm .51241 .01316
Avge SDev %RSD	ppm 1.9820 .0490 2.4710	ppm .49640 .01153 2.3227	ppm .49721 .01298 2.6108	ppm .49694 .01250 2.5150	ppm .51554 .01268 2.4597	ppm .51085 .01340 2.6238	ppm .51241 .01316 2.5688
Avge SDev	ppm 1.9820 .0490	ppm .49640 .01153	ppm .49721 .01298	ppm .49694 .01250	ppm .51554 .01268	ppm .51085 .01340	ppm .51241 .01316
Avge SDev %RSD #1 #2 Errors	ppm 1.9820 .0490 2.4710 1.9473 2.0166	ppm .49640 .01153 2.3227	ppm .49721 .01298 2.6108	ppm .49694 .01250 2.5150 .48810 .50578	ppm .51554 .01268 2.4597	ppm .51085 .01340 2.6238	ppm .51241 .01316 2.5688 .50310 .52172
Avge SDev %RSD #1 #2	ppm 1.9820 .0490 2.4710 1.9473 2.0166	ppm .49640 .01153 2.3227 .48824 .50455	ppm .49721 .01298 2.6108 .48803 .50639	ppm .49694 .01250 2.5150 .48810 .50578	ppm .51554 .01268 2.4597 .50657 .52451	ppm .51085 .01340 2.6238 .50137 .52033	ppm .51241 .01316 2.5688 .50310 .52172
Avge SDev %RSD #1 #2 Errors High Low Elem	ppm 1.9820 .0490 2.4710 1.9473 2.0166 LC Pass 2.2000 1.8000 SE/1	ppm .49640 .01153 2.3227 .48824 .50455 NOCHECK	ppm .49721 .01298 2.6108 .48803 .50639	ppm .49694 .01250 2.5150 .48810 .50578 LC Pass .55000	ppm .51554 .01268 2.4597 .50657 .52451 NOCHECK	ppm .51085 .01340 2.6238 .50137 .52033	ppm .51241 .01316 2.5688 .50310 .52172 LC Pass .55000
Avge SDev %RSD #1 #2 Errors High Low Elem Units	ppm 1.9820 .0490 2.4710 1.9473 2.0166 LC Pass 2.2000 1.8000	ppm .49640 .01153 2.3227 .48824 .50455 NOCHECK	ppm .49721 .01298 2.6108 .48803 .50639 NOCHECK	ppm .49694 .01250 2.5150 .48810 .50578 LC Pass .55000 .45000	ppm .51554 .01268 2.4597 .50657 .52451 NOCHECK V_ppm	ppm .51085 .01340 2.6238 .50137 .52033 NOCHECK	ppm .51241 .01316 2.5688 .50310 .52172 LC Pass .55000
Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge SDev	ppm 1.9820 .0490 2.4710 1.9473 2.0166 LC Pass 2.2000 1.8000 SE/1 ppm .51381 .00917	ppm .49640 .01153 2.3227 .48824 .50455 NOCHECK SE/2 ppm .51742 .01395	ppm .49721 .01298 2.6108 .48803 .50639 NOCHECK SE ppm .51622 .01236	ppm .49694 .01250 2.5150 .48810 .50578 LC Pass .55000 .45000 TL ppm 1.0107 .0275	ppm .51554 .01268 2.4597 .50657 .52451 NOCHECK V_ ppm 1.9836 .0501	ppm .51085 .01340 2.6238 .50137 .52033 NOCHECK ZN ppm 2.0066 .0524	ppm .51241 .01316 2.5688 .50310 .52172 LC Pass .55000
Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge SDev %RSD	ppm 1.9820 .0490 2.4710 1.9473 2.0166 LC Pass 2.2000 1.8000 SE/1 ppm .51381 .00917 1.7855	ppm .49640 .01153 2.3227 .48824 .50455 NOCHECK SE/2 ppm .51742 .01395 2.6966	ppm .49721 .01298 2.6108 .48803 .50639 NOCHECK SE ppm .51622 .01236 2.3946	ppm .49694 .01250 2.5150 .48810 .50578 LC Pass .55000 .45000 TL ppm 1.0107 .0275 2.7176	ppm .51554 .01268 2.4597 .50657 .52451 NOCHECK V_ ppm 1.9836	ppm .51085 .01340 2.6238 .50137 .52033 NOCHECK ZN ppm 2.0066	ppm .51241 .01316 2.5688 .50310 .52172 LC Pass .55000
Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge SDev	ppm 1.9820 .0490 2.4710 1.9473 2.0166 LC Pass 2.2000 1.8000 SE/1 ppm .51381 .00917	ppm .49640 .01153 2.3227 .48824 .50455 NOCHECK SE/2 ppm .51742 .01395	ppm .49721 .01298 2.6108 .48803 .50639 NOCHECK SE ppm .51622 .01236	ppm .49694 .01250 2.5150 .48810 .50578 LC Pass .55000 .45000 TL ppm 1.0107 .0275	ppm .51554 .01268 2.4597 .50657 .52451 NOCHECK V_ ppm 1.9836 .0501	ppm .51085 .01340 2.6238 .50137 .52033 NOCHECK ZN ppm 2.0066 .0524 2.6138 1.9695	ppm .51241 .01316 2.5688 .50310 .52172 LC Pass .55000
Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge SDev %RSD #1 #2	ppm 1.9820 .0490 2.4710 1.9473 2.0166 LC Pass 2.2000 1.8000 SE/1 ppm .51381 .00917 1.7855 .50732 .52030	ppm .49640 .01153 2.3227 .48824 .50455 NOCHECK SE/2 ppm .51742 .01395 2.6966 .50756 .52729	ppm .49721 .01298 2.6108 .48803 .50639 NOCHECK SE ppm .51622 .01236 2.3946 .50748 .50748	ppm .49694 .01250 2.5150 .48810 .50578 LC Pass .55000 .45000 TL ppm 1.0107 .0275 2.7176 .99129 1.0301	Ppm .51554 .01268 2.4597 .50657 .52451 NOCHECK V_ ppm 1.9836 .0501 2.5242 1.9482 2.0190	ppm .51085 .01340 2.6238 .50137 .52033 NOCHECK ZN ppm 2.0066 .0524 2.6138 1.9695 2.0437	ppm .51241 .01316 2.5688 .50310 .52172 LC Pass .55000
Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge SDev %RSD	ppm 1.9820 .0490 2.4710 1.9473 2.0166 LC Pass 2.2000 1.8000 SE/1 ppm .51381 .00917 1.7855	ppm .49640 .01153 2.3227 .48824 .50455 NOCHECK SE/2 ppm .51742 .01395 2.6966 .50756	ppm .49721 .01298 2.6108 .48803 .50639 NOCHECK SE ppm .51622 .01236 2.3946 .50748	ppm .49694 .01250 2.5150 .48810 .50578 LC Pass .55000 .45000 TL ppm 1.0107 .0275 2.7176	Ppm .51554 .01268 2.4597 .50657 .52451 NOCHECK V_ ppm 1.9836 .0501 2.5242 1.9482	ppm .51085 .01340 2.6238 .50137 .52033 NOCHECK ZN ppm 2.0066 .0524 2.6138 1.9695	ppm .51241 .01316 2.5688 .50310 .52172 LC Pass .55000

Analysis Report

05/17/00 09:44:45 AM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 17200 458.8073 2.667558	2 NOTUSED    	3 NOTUSED    	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1	17524						
#2	16875						

page 1

Method: METTRA Sample Name: CCB1 Operator: RJG

Run Time: 05/17/00 09:44:49

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

		140001. 1					
Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00056	.03680	00043	.00078	.00087	.08783	.00012
SDev	.00002	.00145	.00211	.00003	.00012	.01532	.00004
%RSD	3.0728	3.9466	485.70	4.4381	13.398	17.441	32.211
#1	.00057	.03577	00192	.00081	.00095	.07700	.00014
#2	.00055	.03782	.00106	.00076	.00078	.09867	.00009
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.01000	.20000	.00500	5.0000	.00500
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem Units Avge SDev %RSD	CO ppm .00013 .00000	CR ppm .00063 .00000	CU ppm .00025 .00013 49.320	FE ppm .01283 .00499 38.915	MG ppm .01374 .00024 1.7614	MN ppm .00072 .00005 6.3063	MO ppm .00712 .00214 29.990
#1	.00013	.00063	.00016	.00930	.01357	.00075	.00863
#2	.00013	.00063	.00034	.01636	.01391	.00069	.005 <b>61</b>
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.05000	.01000	.02500	.10000	5.0000	.01500	.04000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00018	00009	.00107	.00069	.00359	00078	.00068
SDev	.00024	.00239	.00097	.00144	.00165	.00141	.00149
%RSD	131.84	2695.0	90.093	209.97	45.947	180.94	220.08
#1	.00035	.00160	.00176	.00171	.00242	00177	00038
#2	.00001	00178	.00039	00033	.00475	.00022	.00173
Errors High Low	LC Pass .04000 04000	NOCHECK	NOCHECK	LC Pass .00300 00300	NOCHECK	NOCHECK	LC Pass .06000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00025	.00106	.00079	.00347	.00166	.00108	
SDev	.00369	.00223	.00026	.00252	.00235	.00003	
%RSD	1465.8	210.25	32.530	72.497	141.06	2.6176	
#1 #2	.00286 00236	00052 .00264	.00061 .00097	.00169 .00526	.00332	.00110 .00106	
Errors High Low	NOCHECK	NOCHECK	LC Pass .00500 00500	LC Pass .01000 01000	LC Pass .05000 05000	LC Pass .02000 02000	

657 666

Analysis Report

05/17/00 09:48:54 AM page 2

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED 	6 NOTUSED	7 NOTUSED
Wavlen	371.030						
Avge	16930						- <del>-</del>
SDev	32.49101	<b></b>	<b></b>				
%RSD	.1919095				- <i>-</i>	<b></b>	<del>-</del> -
#1	16907	<b>-</b> -					-
#2	16953						

Operator: RJG

page 1

Method: METTRA Sample Name: DD6A7

Run Time: 05/17/00 09:48:58

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

		- 44404	•				
Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00105	183.35	.06884	3.5965	.00609	H1326.3	.00102
SDev	.00010	.20	.00131	.0038	.00003	2.4	.00038
%RSD	9.6315	.10861	1.8970	.10528	.42101	.18408	37.736
#1	.00113	183.21	.06976	3.5939	.00611	H1328.0	.00075
#2	.00098	183.50	.06791	3.5992	.00608	H1324.6	.00129
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.03861	.16309	.16963	109.53	19.607	2.3656	.00632
SDev	.00031	.00007	.00002	.29	.024	.0002	.00064
%RSD	.80135	.04190	.01381	.26070	.12428	.00956	10.100
#1	.03883	.16304	.16962	109.73	19.624	2.3658	.00587
#2	.03839	.16313	.16965	109.33	19.590	2.3655	.00677
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.09609	.30336	.29820	.29992	.01889	.00853	.01198
SDev	.00070	.00594	.00680	.00255	.00617	.00499	.00128
%RSD	.72840	1.9588	2.2791	.85170	32.635	58.526	10.666
#1	.09659	.29916	.30301	.30173	.02325	.00500	.01108
#2	.09560	.30756	.29340	.29811	.01453	.01206	.01289
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00296	00010	.00092	.01072	.25677	.55060	
SDev	.00885	.00342	.00067	.00991	.00057	.00350	
%RSD	298.51	3328.3	72.605	92.453	.22343	.63663	
#1	00329	.00231	.00045	.01773	.25717	.55307	
#2	.00922	00252	.00139	.00371	.25636	.54812	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

05/17/00 09:53:03 AM

page 2

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED 	7 NOTUSED
Wavlen	371.030						
Avge	17264		- <del>-</del>				
SDev	10.99606						
%RSD	.0636931						
#1	17256						<del></del>
#2	17272			<del></del>			

05/17/00 09:57:13 AM

Method: METTRA Sample Name: CCV3-2 Run Time: 05/17/00 09:53:07 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.0389	24.678	.51733	2.0345	2.0210	50.391	.49066
SDev	.0015	.029	.00238	.0005	.0032	.093	.00161
%RSD	.14379	.11699	.45943	.02512	.15947	.18507	.32795
#1	1.0378	24.658	.51901	2.0341	2.0233	50.457	.49180
#2	1.0400	24.699	.51565	2.0348	2.0187	50.325	.48952
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	27.500	.55000	2.2000	2.2000	55.000	.55000
Low	.90000	22.500	.45000	1.8000	1.8000	45.000	.45000
Elem Units Avge SDev %RSD	CO ppm 2.0283 .0008 .04105	CR ppm 2.0321 .0003 .01319	CU ppm 2.0229 .0005 .02358	FE ppm 25.080 .018 .07375	MG ppm 49.793 .001	MN ppm 1.9960 .0004 .01834	MO ppm 2.0408 .0126 .61639
#1	2.0289	2.0323	2.0232	25.093	49.793	1.9962	2.0319
#2	2.0277	2.0320	2.0225	25.067	49.792	1.9957	2.0497
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.2000	2.2000	2.2000	27.500	55.000	2.2000	2.2000
Low	1.8000	1.8000	1.8000	22.500	45.000	1.8000	1.8000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	2.0130	.50153	.50052	.50086	.52162	.51771	.51901
SDev	.0158	.00198	.00188	.00060	.00407	.00334	.00088
%RSD	.78667	.39562	.37635	.11894	.78055	.64592	.16852
#1	2.0242	.50013	.50185	.50128	.52450	.51535	.51840
#2	2.0018	.50293	.49919	.50043	.51874	.52008	.51963
Errors High Low	LC Pass 2.2000 1.8000	NOCHECK	NOCHECK	LC Pass .55000 .45000	NOCHECK	NOCHECK	LC Pass .55000 .45000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.52221	.52508	.52412	1.0224	2.0215	2.0500	
SDev	.00392	.00283	.00058	.0043	.0065	.0065	
%RSD	.75119	.53814	.11035	.42159	.31992	.31807	
#1	.51944	.52708	.52453	1.0255	2.0260	2.0546	
#2	.52499	.52308	.52372	1.0194	2.0169	2.0454	
Errors High Low	NOCHECK	NOCHECK	LC Pass .55000 .45000	LC Pass 1.1000 .90000	LC Pass 2.2000 1.8000	LC Pass 2.2000 1.8000	

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16698 70.74659 .4236912	2 NOTUSED   	3 NOTUSED   	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    	
#1	16748						<del>-</del> -	
#2	16648						<del>-</del> -	

Method: METTRA Sample Name: CCB2 Run Time: 05/17/00 09:57:17 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00025	.03783	.00028	.00063	.00080	.06431	.00013
SDev	.00062	.00434	.00030	.00009	.00025	.01381	.00016
%RSD	251.58	11.472	108.97	13.617	31.493	21.481	119.98
#1	00019	.04090	.00049	.00057	.00097	.05454	.00024
#2	.00068	.03476	.00006	.00069	.00062	.07408	
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.01000	.20000	.00500	5.0000	.00500
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00025	.00017	.00023	.00942	.01022	.00054	.00655
SDev	.00010	.00053	.00081	.00550	.00289	.00008	.00230
%RSD	38.430	312.41	359.51	58.414	28.249	15.185	35.050
#1	00032	00021	00035	.00553	.00818	.00048	.00817
#2	00018	.00055	.00080	.01331	.01226	.00059	.00493
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.05000	.01000	.02500	.10000	5.0000	.01500	.04000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00104	.00028	.00096	.00074	.00330	.00035	.00133
SDev	.00085	.00077	.00005	.00022	.00134	.00186	.00080
%RSD	81.888	271.45	5.4235	29.964	40.669	529.66	59.671
#1	.00044	00026	.00100	.00058	.00425	00097	.00077
#2	.00164	.00083	.00093		.00235	.00167	.00190
Errors High Low	LC Pass .04000 04000	NOCHECK	NOCHECK	LC Pass .00300 00300	NOCHECK	NOCHECK	LC Pass .06000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00289	.00153	.00198	.00722	.00000	.00101	
SDev	.00102	.00124	.00049	.00131	.00000	.00008	
%RSD	35.221	80.946	24.597	18.141	58.414	7.9652	
#1 #2	.00361 .00217	.00065 .00241	.00164 .00233	.00814 .00629	.00000	.00107 .00095	
Errors High Low	NOCHECK	NOCHECK	LC Pass .00500 00500	LC Pass .01000 01000	LC Pass .05000 05000	LC Pass .02000 02000	

Analysis Report

05/17/00 10:01:22 AM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16726 121.8350 .7284215	2 NOTUSED    	3 NOTUSED   	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    	
#1	16812	<b>-</b> -		- <del>-</del>				
#2	16640	<b>-</b> -						

Operator: RJG

Method: METTRA Sample Name: DD0L6B Run Time: 05/17/00 10:01:26

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00001	.09875	00041	.00157	.00044	.57004	.00007
SDev	.00070	.00263	.00037	.00000	.00009	.00195	.00012
%RSD	5167.5	2.6647	91.116	.20973	20.116	.34150	166.22
#1	00051	.10062	00067	.00157	.00050	.56867	00001
#2	.00048	.09689	00014	.00157		.57142	.00016
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.01000	.20000	.00500	5.0000	.00500
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00086	.00128	.00177	H.11812	.02006	.00173	.00153
SDev	.00037	.00022	.00076	.01394	.00015	.00004	.00025
%RSD	42.378	16.831	42.716	11.804	.76422	2.4671	16.119
#1	00112	.00113	.00123	H.10826	.02017	.00170	.00136
#2	00060	.00144	.00230	H.12798	.01995	.00176	.00170
Errors	LC Pass	LC Pass	LC Pass	LC High	LC Pass	LC Pass	LC Pass
High	.05000	.01000	.02500	.10000	5.0000	.01500	.04000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00036	.00374	.00209	.00264	.00157	.00241	.00213
SDev	.00000	.00047	.00078	.00036	.00271	.00126	.00006
%RSD	1.0689	12.639	37.255	13.692	171.99	52.340	2.7793
#1	.00036	.00341	.00264	.00289	00034	.00330	.00209
#2	.00036	.00408	.00154		.00349	.00152	.00217
Errors High Low	LC Pass .04000 04000	NOCHECK	NOCHECK	LC Pass .00300 00300	NOCHECK	NOCHECK	LC Pass .06000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00236	.00200	.00055	.00219	.00120	.01858	
SDev	.00140	.00093	.00015	.00176	.00056	.00013	
%RSD	59.149	46.431	28.094	80.364	46.703	.67491	
#1 #2	00335 00138	.00266 .00134	.00065 .00044	.00344	.00081 .00160	.01849 .01867	
Errors High Low	NOCHECK	NOCHECK	LC Pass .00500 00500	LC Pass .01000 01000	LC Pass .05000 05000	LC Pass .02000 02000	

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16584 89.55535 .5400114	2 NOTUSED	3 NOTUSED    	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1	16647						
#2	16521		<b></b>			<del>-</del> -	

Operator: RJG

Method: METTRA Sample Name: DD0L6C

Run Time: 05/17/00 10:05:36

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem Units	AG ppm	AL ppm	AS ppm	BA	BE	CA	CD
Avge SDev %RSD	.04749 .00005 .09884	1.7827 .0030 .16987	1.9427 .0070 .35981	ppm 1.9160 .0011 .05988	ppm .04891 .00010 .20169	ppm L.06471 .00259 4.0017	ppm .04693 .00038 .82099
#1	.04746	1.7849	1.9476	1.9168	.04898	L.06288	.04721
#2	.04753	1.7806	1.9377	1.9152	.04884	L.06654	.04666
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Low	LC Pass
High	.06000	2.4000	2.4000	2.4000	.06000	60.000	.06000
Low	.04000	1.6000	1.6000	1.6000	.04000	40.000	.04000
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.48230	.19437	.23408	.98738	L00225	.46940	L00067
SDev	.00138	.00004	.00027	.00394	.00027	.00069	.00035
%RSD	.28622	.02236	.11389	.39898	11.831	.14744	52.347
#1	.48327	.19434	.23426	.98459	L00244	.46989	L00092
#2	.48132	.19440	.23389	.99016	L00207	.46891	L00042
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Low	LC Pass	LC Low
High	.60000	.24000	.30000	1.2000	60.000	.60000	1.2000
Low	.40000	.16000	.20000	.80000	40.000	.40000	.80000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.48402	.46340	.46976	.46764	.00127	00030	L.00022
SDev	.00393	.00022	.00409	.00266	.00040	.00007	.00009
%RSD	.81241	.04717	.87183	.56857	31.915	22.917	40.299
#1	.48680	.46325	.47265	.46952	.00155	00035	L.00028
#2	.48124	.46356	.46686	.46576	.00098	00026	L.00016
Errors High Low	LC Pass .60000 .40000	NOCHECK	NOCHECK	LC Pass .60000 .40000	NOCHECK	NOCHECK	LC Low .60000 .40000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	1.9553	1.9501	1.9518	1.8958	.47568	.49323	
SDev	.0071	.0086	.0034	.0026	.00139	.00083	
%RSD	.36352	.44003	.17197	.13457	.29195	.16814	
#1	1.9502	1.9562	1.9542	1.8976	.47666	.49381	
#2	1.9603	1.9440	1.9494	1.8940	.47470	.49264	
Errors High Low	NOCHECK	NOCHECK	LC Pass 2.4000 1.6000	LC Pass 2.4000 1.6000	LC Pass .60000 .40000	LC Pass .60000 .40000	

	IntStd Mode Elem Wavlen Avge SDev %RSD	7 NOTUSED    
#1 17336		

Operator: RJG

Method: METTRA Sample Name: DCQF8

Run Time: 05/17/00 10:09:45

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

			_				
Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00064	.13078	.00149	.03175	.00046	105.47	00023
SDev	.00033	.00916	.00025	.00022	.00008	.93	.00002
%RSD	51.111	7.0035	17.000	.70992	16.421	.88297	7.3091
#1	00087	.13726	.00131	.03191	.00051	106.13	~.00022
#2	00041	.12431	.00167	.03159	.00041	104.81	00024
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00060	.00049	00005	6.0703	4.2168	.19105	.00126
SDev	.00010	.00048	.00020	.0484	.0256	.00134	.00018
%RSD	17.355	97.328	429.44	.79797	.60638	.70313	14.171
#1	00067	.00015	00019	6.1046	4.2349	.19200	.00139
#2	00053	.00083	.00010	6.0361	4.1987	.19010	.00113
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00100	00127	.00274	.00141	.00162	00015	.00044
SDev	.00126	.00387	.00029	.00110	.00099	.00133	.00121
%RSD	126.31	305.43	10.479	77.966	61.204	871.03	277.33
#1	.00011	00400	.00294	.00063	.00232	.00079	.00130
#2	.00188	.00147	.00254	.00218		00109	00042
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00321	.00253	.00062	.00314	.00067	.00848	
SDev	.00078	.00157	.00131	.00239	.00055	.00001	
%RSD	24.403	62.094	211.18	76.293	81.275	.14141	
#1 #2	00265 00376	.00364	.00155 00031	.00144	.00029 .00106	.00849	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Report

05/17/00 10:13:51 AM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16609 123.8141 .7454842	2 NOTUSED    	3 NOTUSED    	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED   
#1	16521	<b></b>					
#2	16696						

Method: METTRA Sample Name: DCQFL Operator: RJG

Run Time: 05/17/00 10:13:54

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00025	.32454	.00091	.02168	.00053	85.463	00012
SDev	.00027	.00275	.00030	.00015	.00006	.310	.00004
%RSD	109.28	.84751	33.476	.68089	12.325	.36295	36.649
#1	00044	.32649	.00069	.02158	.00057	85.682	00015
#2	00006	.32260	.00112	.02179	.00048	85.244	00009
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00030	.00083	00075	3.3002	4.5531	.17389	.00221
SDev	.00017	.00012	.00014	.0029	.0142	.00041	.00007
%RSD	57.220	14.471	18.687	.08707	.31287	.23676	3.3589
#1	00042	.00091	00085	3.3022	4.5631	.17418	.00215
#2	00018	.00074	00065	3.2982	4.5430	.17360	.00226
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00052	.00055	.00139	.00111	.00211	.00038	.00096
SDev	.00012	.00335	.00044	.00082	.00039	.00019	.00001
%RSD	22.844	612.13	31.377	74.078	18.637	48.905	.57856
#1	.00061	00182	.00170	.00053	.00239	.00025	.00096
#2	.00044	.00291	.00108	.00169	.00183	.00052	.00096
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZŅ	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00051	.00110	.00090	.00242	.00163	.02040	
SDev	.00169	.00392	.00318	.00138	.00000	.00016	
%RSD	332.76	355.95	351.61	57.083	.04103	.78766	
#1	.00170	.00387	.00315	.00145	.00163	.02051	
#2	00069	00167	00134	.00340	.00163	.02029	
Errors	NOCHECK	NOCHECK	LC Pass	LC Pass	LC Pass	LC Pass	

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16784 1.733240 .0103268	2 NOTUSED   	3 NOTUSED   	4 NOTUSED    	5 NOTUSED	6 NOTUSED	7 NOTUSED   	
#1.	16785		<b>-</b> -					
#2	16783			<b>~</b> -				

Method: METTRA Sample Name: DCQFR Run Time: 05/17/00 10:18:03 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

110ac. co	no coll.	I GCCOI. I					
Elem Units Avge SDev %RSD	AG ppm 00027 .00034 125.56	AL ppm .28597 .00537 1.8785	AS ppm .00157 .00078 49.305	BA ppm .04296 .00004 .10077	BE ppm .00030 .00014 47.020	CA ppm 121.92 .27	CD ppm 00025 .00012 47.635
#1	00052	.28977	.00102	.04299	.00040	122.12	00034
#2	00003	.28217	.00212	.04293		121.73	00017
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00108	.00116	00106	7.0520	5.0155	.10568	.00176
SDev	.00045	.00010	.00016	.0018	.0058	.00011	.00016
%RSD	42.288	8.1885	14.774	.02491	.11586	.10479	9.1506
#1	00075	.00123	00117	7.0532	5.0196	.10576	.00165
#2	00140	.00110	00095	7.0508	5.0114	.10561	.00187
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00089	.00007	.00079	.00055	.00216	.00063	.00114
SDev	.00012	.00179	.00009	.00065	.00082	.00089	.00087
%RSD	13.483	2501.8	10.944	118.30	38.116	142.68	76.549
#1	.00098	00119	.00073	.00009	.00275	.00126	.00176
#2	.00081	.00133	.00085	.00101	.00158	00001	.00052
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem Units Avge SDev %RSD	SE/1 ppm 00047 .00186 398.99	SE/2 ppm 00033 .00397 1208.4	SE ppm 00037 .00327 872.08	TL ppm 00142 .00014 9.9644	V_ppm .00184 .00000 .02491	ZN ppm .01101 .00012 1.0409	
#1	00179	00313	00268	00152	.00184	.01109	
#2	.00085	.00248	.00194	00132	.00184	.01093	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis	Report
----------	--------

05/17/00 10:22:08 AM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 <b>NOTUSE</b> D	6 NOTUSED	7 NOTUSED
Elem	Y			m			
Wavlen	371.030	- <del>-</del>					<b>→</b> →
Avge	16312		<b></b>				
SDev	34.96657						~ -
%RSD	.2143581					**	** ***
#1	16337		<b></b>			==	
#2	16288	<b></b>	~	- ~			

Operator: RJG

Method: METTRA Sample Name: DCQFX

Run Time: 05/17/00 10:22:12

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00039	.43543	.00100	.06223	.00049	161.18	00043
SDev	.00014	.01030	.00086	.00030	.00002	1.61	.00001
%RSD	37.268	2.3666	86.250	.48075	4.7122	.99902	1.7692
#1	00028	.44271	.00160	.06244	.00051	162.32	00044
#2	00049	.42814	.00039	.06202		160.05	00043
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00115	.00176	00023	12.517	9.1671	.05692	.00150
SDev	.00023	.00002	.00022	.067	.0838	.00048	.00048
%RSD	20.180	1.2453	95.632	.53906	.91407	.84963	31.868
#1	00131	.00175	00007	12.565	9.2264	.05727	.00183
#2	00099	.00178	00038	12.469	9.1079	.05658	.00116
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00041	00092	.00191	.00097	.00277	00232	00062
SDev	.00043	.00072	.00101	.00091	.00276	.00109	.00019
%RSD	105.21	78.024	52.904	94.241	99.561	47.053	30.616
#1	.00071	00041	.00262	.00161	.00082	00155	00076
#2	.00010	00142	.00119	.00032	.00472	00309	00049
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00308	.00362	.00139	00043	.00183	.00795	
SDev	.00002	.00029	.00019	.00171	.00018	.00000	
%RSD	.49502	8.0432	13.623	399.48	9.6868	.05355	
#1 #2	00310 00307	.00383	.00152 .00126	.00078 00164	.00170 .00195	.00794 .00795	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Report

05/17/00 10:26:17 AM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	Y			<b>-</b> -			
Wavlen	371.030 16509					- <del>-</del>	
Avge SDev	109.1776						
%RSD	.6613175	_ <b>_</b>			<del></del>	<del>-</del> -	
0102	•••						
#1	16432						
#2	16586						

page 1

Method: METTRA Sample Name: DCQG5 Run Time: 05/17/00 10:26:21 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

		ractor. r					
Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00019	.57472	.00183	.14390	00008	H697.95	00015
SDev	.00057	:01282	.00021	.00018	.00024	.55	.00019
%RSD	297.94	2.2303	11.535	.12748	279.26	.07870	129.62
#1	00059	.58378	.00197	.14377	.00008	H698.34	00001
#2	.00021	.56565	.00168	.14403	00025	H697.56	00028
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00039	.00206	.00009	7.5649	43.934	.33511	.00102
SDev	.00053	.00029	.00084	.0063	.018	.00052	.00097
%RSD	134.89	14.254	951.59	.08288	.04145	.15426	95.722
#1	00077	.00186	00051	7.5605	43.921	.33474	.00033
#2	00002	.00227	.00068	7.5694	43.947	.33547	.00171
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00115	00002	.00042	.00027	.00301	.00090	.00161
SDev	.00053	.00328	.00183	.00013	.00437	.00004	.00143
%RSD	45.536	17747.	439.15	46.162	145.01	4.7954	88.839
#1 #2	.00078 .00153	00234 .00230	.00171 00088	.00036	.00610 00008	.00087	.00261 .00060
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00200	00096	00131	.00401	.00422	.00892	
SDev	.00272	.00189	.00217	.00337	.00001	.00010	
%RSD	136.10	197.34	166.11	84.083	.16790	1.1107	
#1 #2	00008 00393	.00038 00230	.00023 00284	.00162 .00639	.00422	.00899 .00885	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

657 686

Analysis	Report
----------	--------

05/17/00 10:30:26 AM

page 2

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED 	6 NOTUSED 	7 NOTUSED
Wavlen	371.030						
Avge	15355						
SDev	37.22876				an ad		
%RSD	.2424588			<b></b>			<del></del>
#1	15381		<del>-</del> -				<del></del>
#2	15328				<del>-</del> -	<del>-</del>	

5084

Method: METTRA Sample Name: DCQGK Operator: RJG

Run Time: 05/17/00 10:30:30

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	<b>a</b> n
Units Avge	ppm .00003	ppm .20708	ppm 00083	ppm .06514	ppm	mqq	ppm D
SDev %RSD	.00021	.00734	.00112	.00033	.00031 .00001	121.59 .68	00032 .00005
	752.13	3.5426	134.13	.51194	3.5520	.56332	16.704
#1	.00018	.21227	00162	.06538	.00030	122.07	00028
#2	00012	.20189	00004	.06491		121.10	00036
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00697	.00073	00028	10.734	9.0930	.32877	00047
SDev	.00003	.00056	.00019	.052	.0374	.00163	.00045
%RSD	.49339	76.957	67.878	.48401	.41169	.49554	95.955
#1	.00695	.00113	00041	10.771	9.1195	.32992	00079
#2	.00700		00014	10.697	9.0665	.32761	00015
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01076	00073	.00197	.00107	.00348	00144	.00020
SDev	.00021	.00151	.00122	.00131	.00262	.00099	.00153
%RSD	1.9710	205.52	61.894	123.11	75.236	68.654	779.38
#1 #2	.01091 .01061	.00033 00180	.00283	.00200	.00534	00074 00215	.00128
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00240	.00249	.00086	00106	00067	.02806	
SDev	.00202	.00213	.00210	.00089	.00001	.00040	
%RSD	84.192	85.776	243.08	83.490	.79584	1.4319	
#1	00382	.00098	00062	00169	00066	.02834	
#2	00097	.00400	.00234	00044	00067	.02777	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Report

05/17/00 10:34:36 AM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16186 38.43167 .2374425	2 NOTUSED    	3 NOTUSED    	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1 #2	16158 16213		 	<del>-</del> -	 		<del></del>

5086

Method: METTRA Sample Name: DCQGM Operator: RJG

Run Time: 05/17/00 10:34:39

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00035	.08882	.01887	.06891	.00024	149.35	00121
SDev	.00022	.00646	.00024	.00044	.00018	.66	.00045
%RSD	63.357	7.2707	1.2624	.64171	75.740	.44009	37.102
#1	00051	.09338	.01871	.06860	.00036	148.88	00153
#2	00019	.08425	.01904	.06923	.00011	149.81	00089
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04515	.00055	00276	86.645	19.977	.50791	00154
SDev	.00034	.00005	.00027	.611	.123	.00335	.00014
%RSD	.74999	8.1201	9.7110	.70486	.61430	.65978	9.3622
#1	.04491	.00052	00295	86.213	19.890	.50554	00144
#2	.04539	.00059	00257	87.077	20.064	.51028	00164
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.07581	00384	00059	00167	.00195	.00124	.00148
SDev	.00099	.00076	.00017	.00037	.00158	.00151	.00153
%RSD	1.3065	19.741	28.628	21.835	80.681	121.83	103.74
#1	.07511	00330	00047	00141	.00084	.00017	.00039 ·
#2	.07651	00438	00071	00193	.00306	.00231	.00256
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00212	.00010	00064	00184	00094	.07293	
SDev	.00203	.00342	.00295	.00569	.00022	.00028	
%RSD	95.763	3394.1	462.53	309.92	23.491	.38862	
#1.	00068	.00252	.00145	.00219	00079	.07273	
#2	00356	00232	00273	00586	00110	.07313	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Ana	lvs	is	Report
LT TOT	J, 7 🔾		

05/17/00 10:38:45 AM

page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16049 129.8245 .8089285	2 NOTUSED    	NOTUSED	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1	16141						
#2	15957						<del></del>

Method: METTRA Sample Name: DCQGN Operator: RJG

Run Time: 05/17/00 10:38:48

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem Units Avge SDev %RSD	AG ppm 00043 .00010 24.403	AL ppm .06261 .00030 .47106	AS ppm .00481 .00142 29.501	BA ppm .03103 .00004 .12063	BE ppm .00063 .00004 5.9483	CA ppm 101.36 .06	CD ppm 00040 .00006 14.779
#1	00050	.06282	.00581	.03101	.00066	101.40	00045
#2	00035	.06240	.00381	.03106		101.32	00036
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01836	.00047	00153	18.093	7.0432	.22740	00083
SDev	.00084	.00019	.00016	.030	.0054	.00015	.00031
%RSD	4.5712	41.268	10.671	.16395	.07660	.06450	36.950
#1	.01777	.00060	00164	18.072	7.0394	.22730	00105
#2	.01895		00141	18.114	7.0470	.22751	00061
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.02452	00055	.00119	.00061	00041	.00088	.00045
SDev	.00003	.00105	.00040	.00008	.00085	.00264	.00148
%RSD	.11190	188.84	33.552	13.680	208.65	300.06	327.46
#1	.02454	.00019	.00090	.00067	00101	.00274	.00150
#2	.02450	00129	.00147	.00055	.00019	00099	00059
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00004	.00190	.00128	.00150	.00079	.04052	
SDev	.00424	.00104	.00210	.00384	.00001	.00004	
%RSD	11152.	54.653	164.34	255.61	.89395	.09167	
#1	.00303	.00264	.00277	00121	.00080	.04054	
#2	00296	.00117	00021	.00422	.00079	.04049	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

05/17/00 10:42:54 AM

page 2

IntStd Mode Elem Wavlen Avge SDev	1 Counts Y 371.030 16492 62.22539	2 NOTUSED   	NOTUSED	4 NOTUSED   	5 NOTUSED   	6 NOTUSED   	7 NOTUSED   
%RSD #1	.3773157						
#2	16448					- <del>-</del>	

Operator: RJG

Method: METTRA Sample Name: CCV3-3

Run Time: 05/17/00 10:42:57

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.0327	24.082	.51604	1.9899	2.0019	49.735	.48436
SDev	.0171	.431	.01154	.0335	.0354	.895	.00932
%RSD	1.6516	1.7887	2.2363	1.6838	1.7700	1.8004	1.9234
#1	1.0448	24.386	.52420	2.0136	2.0269	50.369	.49095
#2	1.0206	23.777	.50788	1.9662	1.9768	49.102	.47777
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	27.500	.55000	2.2000	2.2000	55.000	.55000
Low	.90000	22.500	.45000	1.8000	1.8000	45.000	.45000
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.9922	2.0011	1.9979	24.376	49.199	1.9674	2.0243
SDev	.0331	.0345	.0357	.410	.873	.0330	.0168
%RSD	1.6633	1.7222	1.7852	1.6830	1.7749	1.6760	.83103
#1	2.0156	2.0255	2.0231	24.666	49.817	1.9907	2.0362
#2	1.9688	1.9768	1.9727	24.086	48.582	1.9441	2.0124
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.2000	2.2000	2.2000	27.500	55.000	2.2000	2.2000
Low	1.8000	1.8000	1.8000	22.500	45.000	1.8000	1.8000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	2.0268	.49269	.49686	.49547	.51689	.51578	.51615
SDev	.0451	.00744	.00664	.00691	.00535	.00867	.00756
%RSD	2.2239	1.5096	1.3367	1.3940	1.0354	1.6799	1.4650
#1	2.0586	.49795	.50 <b>15</b> 6	.50036	.52067	.52190	.52149
#2	1.9949	.48743	.49217	.49059	.51310	.50965	.51080
Errors High Low	LC Pass 2.2000 1.8000	NOCHECK	NOCHECK	LC Pass .55000 .45000	NOCHECK	NOCHECK	LC Pass .55000 .45000
Elem Units Avge SDev %RSD	SE/1 ppm .51966 .00866 1.6656	SE/2 ppm .51506 .00587 1.1399	SE ppm .51659 .00680 1.3160	TL ppm 1.0147 .0191 1.8837	V_ppm 2.0018 .0411 2.0509	ZN ppm 2.0473 .0422 2.0630	
#1	.52578	.51921	.52140	1.0282	2.0309	2.0772	
#2	.51354	.51091	.51178	1.0012	1.9728	2.0175	
Errors High Low	NOCHECK	NOCHECK	LC Pass .55000 .45000	LC Pass 1.1000 .90000	LC Pass 2.2000 1.8000	LC Pass 2.2000 1.8000	

657 694 Analysis Report

05/17/00 10:47:03 AM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	Y				<del></del>		
Wavlen	371.030		<b>→ →</b>			~ -	
Avge	16447		<b></b>	<b>-</b> →			
SDev	227.3003		<del>-</del> -				
%RSD	1.381998	~~					
#1	16286			<b>-</b>			
#2	16608	** **					

5092

Method: METTRA Sample Name: CCB3 Operator: RJG

Run Time: 05/17/00 10:47:06

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00023	.06514	00183	.00059	.00124	.02803	.00029
SDev	.00005	.00492	.00184	.00023	.00006	.01072	.00011
%RSD	20.386	7.5481	100.69	38.881	4.4752	38.221	38.081
#1	.00020	.06862	00313	.00042	.00120	.02046	.00021
#2	.00027	.06166	00053	.00075	.00128	.03561	.00037
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.01000	.20000	.00500	5.0000	.00500
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00049	.00055	00126	.00978	.01393	.00060	.00618
SDev	.00053	.00013	.00050	.00883	.00250	.00020	.00237
%RSD	108.27	23.019	39.363	90.294	17.980	33.679	38.346
#1 .	00086	.00064	00161	.00354	.01216	.00045	.00785
#2	00011	.00046	00091	.01603	.01570		.00450
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.05000	.01000	.02500	.10000	5.0000	.01500	.04000
Low	05000	01000	02500	10000	~5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00076	.00074	.00145	.00121	.00203	00073	.00019
SDev	.00018	.00247	.00081	.00028	.00524	.00361	.00067
%RSD	23.740	332.06	55.929	23.224	258.50	497.87	349.27
#1	.00063	.00249	.00087	.00141	00168	.00183	.00066
#2	.00089	00100	.00202	.00101	.00573	~.00328	00028
Errors High Low	LC Pass .04000 04000	NOCHECK	NOCHECK	LC Pass .00300 00300	NOCHECK	NOCHECK	LC Pass .06000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00216	.00131	.00159	.00130	.00066	.00107	
SDev	.00134	.00120	.00124	.00192	.00019	.00002	
%RSD	61.949	91.303	78.064	147.19	28.145	2.1679	
#1	.00121	.00046	.00071	.00266	.00079	.00105	
#2	.00310	.00216	.00247	00005	.00053	.00109	
Errors High Low	NOCHECK	NOCHECK	LC Pass .00500 00500	LC Pass .01000 01000	LC Pass .05000 05000	LC Pass .02000 02000	

05/17/00 10:51:12 AM

page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	Y						
Wavlen	371.030 16367						
Avge SDev	54.34088						
%RSD	.3320083				= **		<b>-</b>
#1	16329						
#2	16406						

05/17/00 10:55:21 AM

Method: METTRA Sample Name: DCQGQ Operator: RJG

Run Time: 05/17/00 10:51:16

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00157	2.5237	.00261	.01748	.00852	97.223	.00332
SDev	.00041	.0022	.00281	.00008	.00001	.325	.00009
%RSD	25.850	.08566	107.56	.45778	.08431	.33413	2.8443
#1	00128	2.5222	.00459	.01742	.00852	96.993	.00325
#2	00185	2.5252	.00062	.01754	.00851	97.452	
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem Units Avge SDev %RSD	CO ppm .20464 .00085 .41469	CR ppm .00589 .00043 7.3742	CU ppm 00168 .00035 21.038	FE ppm 38.370 .175 .45503	MG ppm 33.662 .142	MN ppm .35757 .00182 .50805	MO ppm .00017 .00055 326.99
#1	.20404	.00620	00193	38.247	33.562	.35628	.00056
#2	.20524	.00558	00143	38.494	33.763	.35885	00022
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.27532	.00010	.00252	.00172	.00166	00045	.00025
SDev	.00126	.00300	.00164	.00010	.00514	.00356	.00067
%RSD	.45854	2975.8	65.170	5.6789	309.84	796.53	261.70
#1	.27443	.00222	.00136	.00165	00198	.00207	,00072
#2	.27622	00202	.00368	.00178	.00530	00297	00022
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem Units Avge SDev %RSD	SE/1 ppm 00263 .00270 102.43	SE/2 ppm 00079 .00304 382.46	SE ppm 00141 .00292 207.86	TL ppm .00138 .00039 28.457	V_ ppm .00235 .00006 2.6194	ZN ppm .57395 .00027	
#1 #2	00073 00454	.00135 00294	.00066 00348	.00110 .00166	.00230	.57376 .57414	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED 	7 NOTUSED 
371.030	<b></b>					
18031	<del></del>	<b></b>	<del></del>			
37.40622						
.2074510		<b>-</b>				
18005	<b></b>	<del>-</del> -				
18058					<del>-</del> -	
	Y 371.030 18031 37.40622 .2074510	Counts NOTUSED Y 371.030 18031 37.406222074510 18005	Counts NOTUSED NOTUSED Y 371.030 18031 37.406222074510	Counts NOTUSED NOTUSED NOTUSED Y 371.030 18031 37.406222074510	Counts NOTUSED NOTUSED NOTUSED Y 371.030 18031 37.406222074510	Counts NOTUSED NOTUSED NOTUSED NOTUSED  Y

page 1

Method: METTRA Sample Name: DCQGV Operator: RJG

Run Time: 05/17/00 10:55:25

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

		<u> </u>					
Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00010	.07433	.01129	.01995	.00059	84.026	00021
SDev	.00049	.00770	.00029	.00006	.00002	.320	.00000
%RSD	494.52	10.358	2.5467	.30461	3.8874	.38137	1.4924
#1	.00025	.07978	.01109	.01999	.00061	84.253	00021
#2	00045	.06889	.01150	.01990	.00057	83.800	00020
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00024	.00046	00150	9.8584	17.122	.52565	.00059
SDev	.00002	.00010	.00028	.0146	.040	.00070	.00047
%RSD	9.9913	21.984	18.411	.14766	.23366	.13240	79.057
#1	00022	.00053	00130	9.8687	17.151	.52615	.00092
#2	00026		00169	9.8481	17.094	.52516	.00026
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00107	.00327	.00054	.00145	.00253	00111	.00010
SDev	.00000	.00325	.00020	.00121	.00168	.00074	.00006
%RSD	.39198	99.472	36.582	83.896	66.235	66.978	60.979
#1	.00107	.00557	.00068	.00230	.00371	00163	.00015
#2	.00107	.00097	.00040	.00059	.00134	00058	.00006
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00207	.00185	.00054	.00136	.00204	.00906	
SDev	.00193	.00161	.00172	.00049	.00000	.00003	
%RSD	93.406	87.149	316.80	36.274	.10797	.30736	
#1	00344	.00071	00067	.00101	.00204	.00904	
#2	00070	.00298	.00176	.00171	.00204	.00907	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

	657	700
Analysis	Report	

05/17/00 10:59:31 AM page 2

IntStd Mode Elem Wavlen	1 Counts Y 371.030	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED 	7 NOTUSED
Avge	16286						
SDev	48.93124				** **		<del>-</del> -
%RSD	.3004460					·**	
#1	16252						
#2	16321					<del></del>	

Operator: RJG

Method: METTRA Sample Name: DCVMD

Run Time: 05/17/00 10:59:35

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00011	.73794	.00052	.21204	.00001	543.19	.00192
SDev	.00017	.00372	.00086	.00124	.00014	.82	.00002
%RSD	155.81	.50354	165.50	.58418	982.54	.15062	.93592
#1	.00022	.74057	.00113	.21116	.00012	542.61	.00194
#2	00001	.73532	00009	.21291	00009	543.77	.00191
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01582	.00208	.00027	1.3616	82.139	1.4969	00050
SDev	.00024	.00002	.00036	.0014	.278	.0074	.00072
%RSD	1.5156	1.0584	134.48	.10083	.33825	.49124	143.16
#1	.01599	.00207	.00001 .	1.3607	81.942	1.4917	.00001
#2	.01565	.00210		1.3626	82.335	1.5021	00101
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01177	.00127	00002	.00041	.00325	00002	.00107
SDev	.00153	.00041	.00027	.00004	.00088	.00096	.00093
%RSD	13.041	32.183	1376.0	10.524	27.220	4074.3	87.689
#1	.01285	.00155	00021	.00038	.00387	.00066	.00173
#2	.01068	.00098	.00017		.00262	00070	.00041
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00043	00115	~.00063	.00136	.00517	.02823	
SDev	.00375	.00374	.00125	.00185	.00239	.00011	
%RSD	862.83	324.14	199.66	135.32	46.184	.38125	
#1	00221	.00149	.00026	.00267	.00686	.02815	
#2	.00308	00380	00151	.00006	.00348	.02830	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

05/17/00 11:03:40 AM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 15242 54.44722 .3572231	2 NOTUSED   	3 NOTUSED   	4 NOTUSED    	5 NOTUSED   	6 NOTUSED   	7 NOTUSED   
#1 #2	15280 15203						

Method: METTRA Sample Name: DCVME Run Time: 05/17/00 11:03:44 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG ppm00006 .00008 127.76	AL	AS	BA	BE	CA	CD
Units		ppm	ppm	ppm	ppm	ppm	ppm
Avge		1.6279	.00121	.10997	.00060	172.30	00044
SDev		.0036	.00079	.00030	.00019	.50	.00010
%RSD		.22361	65.282	.27290	31.855	.28811	23.791
#1	00012	1.6253	.00065	.10976	.00074	172.65	00036
#2	00001	1.6304	.00177	.11018	.00047	171.95	00051
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00059	.00495	00004	8.1960	5.3524	.07090	.00171
SDev	.00025	.00010	.00055	.0121	.0039	.00004	.00000
%RSD	43.278	1.9211	1266.7	.14808	.07226	.05653	.23372
#1	00041	.00488	00043	8.1875	5.3497	.07087	.00171
#2	00077	.00502	.00034	8.2046	5.3551	.07093	.00170
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00231	.00355	.00114	.00194	.00070	.00038	.00049
SDev	.00037	.00025	.00137	.00100	.00288	.00074	.00047
%RSD	16.006	6.9612	120.91	51.500	411.09	193.52	94.836
#1	.00205	.00337	.00016	.00123	00134	.00091	.00016
#2	.00258		.00211	.00265	.00274	00014	.00082
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00012	.00016	.00015	00133	.00294	.01778	
SDev	.00233	.00026	.00095	.00222	.00000	.00013	
%RSD	1898.0	157.55	632.85	167.69	.08313	.74979	
#1	.00177	.00035	.00082	.00025	.00294	.01787	
#2	00153	00002	00052	00290	.00294	.01769	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

657 704

Analysis Report

05/17/00 11:07:50 AM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem Wavlen	Y 371.030						
Avge	16122						
SDev	14.46047					<b></b>	
%RSD	.0896953	<del></del> -				TTD 400-	,
#1	16112						
#2	16132		* *	• •			

Method: METTRA Sample Name: DCVMF Operator: RJG

Run Time: 05/17/00 11:07:54

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00006	.40368	.00005	.08033	00025	586.52	.00005
SDev	.00011	.01698	.00150	.00020	.00030	4.88	.00015
%RSD	185.47	4.2052	2827.5	.24751	120.36	.83181	320.12
#1	00002	.41569	00101	.08047	00004	589.97	00006
#2	.00014	.39168	.00112	.08019	00047	583.07	.00016
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00904	.00266	00089	.51918	298.21	3.4514	00070
SDev	.00029	.00004	.00087	.00499	1.41	.0101	.00086
%RSD	3.2401	1.5661	97.762	.96114	.47393	.29376	121.77
#1	.00925	.00269	00151	.51566	299.21	3.4586	00010
#2	.00884	.00263	00028	.52271	297.21	3.4442	00131
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00817	00054	.00062	.00023	.00100	00042	.00005
SDev	.00099	.00146	.00083	.00104	.00124	.00093	.00021
%RSD	12.124	269.96	134.87	451.82	123.57	221.22	401.49
#1	.00747	.00049	.00120	.00097	.00013	.00024	.00020
#2	.00887	00158		00051	.00187	00108	00010
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00426	00112	.00067	.00009	.00812	.01186	
SDev	.00093	.00014	.00041	.00153	.00003	.00015	
%RSD	21.908	12.769	60.261	1756.7	.43031	1.2858	
#1	.00360	00122	.00039	.00117	.00815	.01197	
#2	.00492	00102	.00096	00099	.00810	.01175	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

05/17/00 11:11:59 AM

page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 14605 66.29126 .4539090	2 NOTUSED    	3 NOTUSED    	4 NOTUSED   	5 NOTUSED   	6 NOTUSED    	7 NOTUSED    
#1	14558					<del></del>	
#2	14651						

page 1

Method: METTRA Sample Name: DCVMG Run Time: 05/17/00 11:12:03 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00039	.10012	.00004	.07133	.00057	162.38	00023
SDev	.00015	.01145	.00146	.00017	.00003	.33	.00003
%RSD	37.443	11.433	3997.2	.24020	5.8401	.20377	11.856
#1	00029	.10821	.00107	.07121	.00059	162.62	00025
#2	00049	.09203	00100	.07146	.00055	162.15	00021
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00154	.00097	.00105	1.6086	6.1583	.08305	.00594
SDev	.00004	.00003	.00055	.0086	.0162	.00007	.00021
%RSD	2.5911	3.4798	52.628	.53299	.26311	.08046	3.4636
#1	00151	.00099	.00066	1.6147	6.1698	.08310	.00608
#2	00156	.00094	.00144	1.6025	6.1469	.08300	.00579
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00063	.00042	.00073	.00062	.00185	.00077	.00113
SDev	.00013	.00074	.00017	.00036	.00186	.00396	.00326
%RSD	20.354	175.48	22.945	57.361	100.67	515.39	288.91
#1	.00072	00010	.00061	.00037	.00317	.00357	.00343
#2	.00054	.00095	.00084		.00053	00203	00118
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00226	00026	00093	.00003	.00095	.01365	
SDev	.00068	.00379	.00230	.00355	.00038	.00020	
%RSD	29.991	1440.9	248.36	13809.	39.978	1.4357	
#1	00178	00294	00256	.00253	.00122	.01379	
#2	00274	.00242	.00070	00248	.00068	.01351	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

05/17/00 11:16:08 AM

page 2

			•				
IntStd	1	2	3	4	5	6	7
Mode	Counts	NOTUSED	NOTUSED	NOTUSED	NOTUSED	NOTUSED	NOTUSED
Elem	Y		<b></b>				
Wavlen	371.030						
Avge	16316			~ -			~ _
SDev	101.8945						
%RSD	.6245066	<b></b> _	<b></b> -			<b></b>	
#1	16244						
#2	16388				<b>~ ~</b>		

Method: METTRA Sample Name: DCVMH Run Time: 05/17/00 11:16:12 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00020	.10930	.00021	.06884	00002	H639.20	00003
SDev	.00030	.01006	.00159	.00174	.00027	11.32	.00007
%RSD	148.76	9.2066	758.22	2.5249	1634.1	1.7705	289.12
#1	.00041	.11641	00092	.06761	.00018	Н631.20	.00003
#2	00001	.10218	.00134	.07007	00021	Н647.20	00008
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00310	.00105	00025	.10851	41.767	1.0215	00028
SDev	.00003	.00043	.00059	.01581	.785	.0198	.00022
%RSD	.97650	41.154	235.59	14.568	1.8803	1.9380	79.155
#1	.00308	.00136	00067	.09733	41.211	1.0075	00012
#2		.00075	.00017	.11969	42.322	1.0355	00043
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00442	.00007	00104	00067	.00178	.00071	.00107
SDev	.00097	.00113	.00203	.00097	.00072	.00096	.00088
%RSD	21.891	1684.2	194.43	144.81	40.216	136.40	82.745
#1	.00374	00073	.00039	.00002	.00229	.00139	.00169
#2	.00511	.00087	00248	00136	.00128	.00003	.00044
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00276	.00125	00009	00031	.00131	.01005	
SDev	.00182	.00253	.00229	.00065	.00062	.00015	
%RSD	65.876	202.06	2670.7	208.01	47.320	1.5414	
#1	00148	.00304	.00154	00077	.00087	.00994	
#2	00405	00054	00171	.00015	.00174	.01016	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Report

05/17/00 11:20:18 AM

page 2

IntStd Mode .Elem	1 Counts Y	2 NOTUSED	3 · NOTUSED	4 NOTUSED	5 NOTUSED 	6 NOTUSED	7 NOTUSED
Wavlen	371.030						
Avge	15247						
SDev	155.5635						
%RSD	1.020302						
#1	15357						
#2	15137					<del>-</del> -	

Operator: RJG

Method: METTRA Sample Name: DCVMHP5 Run Time: 05/17/00 11:20:22

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem Units Avge SDev %RSD	AG ppm 00026 .00035 131.08	AL ppm .06249 .01120 17.928	AS ppm 00001 .00048 4541.8	BA ppm .01349 .00005 .39305	BE ppm .00049 .00016 32.016	CA ppm 130.16 .19	CD ppm 00013 .00012 94.748
#1	00002	.07041	.00033	.01345	.00060	130.02	00004
#2	00051	.05457	00035	.01352	.00038	130.29	00022
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00004	00023	00178	.02389	8.1770	.19950	00065
SDev	.00010	.00005	.00027	.00843	.0270	.00050	.00022
%RSD	223.98	21.346	14.928	35.286	.33030	.24933	34.736
#1	00012	00027	00197	.02985	8.1579	.19915	00049
#2	.00003	00020	00160	.01793	8.1961	.19985	00081
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00003	00026	.00101	.00058	.00227	.00083	.00131
SDev	.00025	.00037	.00200	.00121	.00065	.00125	.00105
%RSD	951.29	141.38	198.64	207.16	28.515	150.38	80.166
#1	00015	00000	00041	00027	.00273	.00172	.00205
#2	.00020	00052	.00242	.00144	.00181	00005	.00057
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00082	00129	00113	.00076	.00014	.00239	
SDev	.00228	.00063	.00118	.00055	.00018	.00003	
%RSD	278.52	48.724	103.92	72.132	133.49	1.2996	
#1 #2	00243 .00079	00174 00085	00197 00030	.00037 .00114	.00001	.00237 .00241	

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Wavlen	371.030						
Avge	16290						
SDev	22.45064						
%RSD	.1378179						
#1	16306						
#2	16274		<b></b>				

5110

Method: METTRA Sample Name: DCVMHS Run Time: 05/17/00 11:24:31 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.05159	2.1979	2.0957	2.0126	.04880	H610.94	.04616
SDev	.00024	.0079	.0160	.0119	.00008	1.21	.00011
%RSD	.46024	.36157	.76493	.59063	.16475	.19785	.23409
#1	.05176	2.2035	2.1071	2.0042	.04886	H610.08	.04623
#2	.05142	2.1923	2.0844	2.0210	.04875	H611.79	.04608
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.48727	.19566	.25610	1.0618	39.977	1.4600	.00032
SDev	.00053	.00028	.00059	.0003	.079	.0042	.00118
%RSD	.10948	.14288	.22892	.02784	.19657	.28892	367.85
#1	.48689	.19546	.25568	1.0620	39.921	1.4570	00051
#2	.48765	.19586	.25651	1.0616	40.032	1.4630	.00116
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.49522	.48262	.47989	.48080	.00077	.00093	.00088
SDev	.00080	.00097	.00072	.00016	.00157	.00077	.00001
%RSD	.16182	.20074	.14964	.03252	204.20	82.986	.98500
#1	.49579	.48193	.48040	.48091	00034	.00147	.00087
#2	.49466	.48330	.47938	.48069	.00188	.00038	.00088
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	2.1003	2.1015	2.1011	2.1279	.49649	.53677	
SDev	.0040	.0040	.0040	.0315	.00106	.00237	
%RSD	.19187	.19139	.19156	1.4822	.21425	.44238	
#1	2.0975	2.0987	2.0983	2.1502	.49574	.53845	
#2	2.1032	2.1044	2.1040	2.1056	.49724	.53509	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

657 714

Analysis Report

05/17/00 11:28:36 AM

page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 15708 2.898862 .0184545	2 NOTUSED    	NOTUSED	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    	
#1 #2	15710 15706		 					

Method: METTRA Sample Name: DCVMHD Run Time: 05/17/00 11:28:40 Operator: RJG Method: METTRA

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.05300	2.2843	2.1676	2.0710	.05004	H627.73	.04749
SDev	.00055	.0280	.0189	.0074	.00042	3.37	.00043
%RSD	1.0279	1.2242	.87386	.35878	.83567	.53676	.89753
#1	.05262	2.3041	2.1810	2.0762	.05033	H630.11	.04779
#2	.05339	2.2645	2.1542	2.0657	.04974	H625.35	.04719
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.50416	.20223	.26395	1.1039	41.048	1.4999	00011
SDev	.00404	.00079	.00074	.0115	.233	.0065	.00001
%RSD	.80212	.39023	.28143	1.0438	.56739	.43211	9.3290
#1	.50702	.20279	.26447	1.1120	41.213	1.5045	00010
#2	.50130	.20167	.26342	1.0957	40.884	1.4953	00011
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem Units Avge SDev %RSD	NI ppm .51441 .00456 .88696	PB/1 ppm .49644 .00035	PB/2 ppm .49528 .00104 .21106	PB ppm .49567 .00058 .11698	SB/1 ppm .00135 .00063 46.557	SB/2 ppm 00060 .00156 259.81	SB ppm .00005 .00083 1687.6
#1	.51763	.49619	.49602	.49608	.00180	00170	00054
#2	.51118	.49669	.49454	.49526		.00050	.00064
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	2.1796	2.1775	2.1782	2.1975	.51248	.55067	
SDev	.0016	.0002	.0004	.0214	.00213	.00521	
%RSD	.07465	.00879	.01902	.97408	.41568	.94565	
#1	2.1807	2.1774	2.1785	2.2126	.51399	.55435	
#2	2.1784	2.1776	2.1779	2.1824	.51097	.54699	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

IntStd Mode Elem Wavlen	1 Counts Y 371.030	2 NOTUSED 	3 NOTUSED	4 NOTUSED	5 NOTUSED 	6 NOTUSED 	7 NOTUSED 
Avge	15269						
SDev	57.45243			<del>-</del>			
%RSD	.3762715		<b>→</b>		<del>-</del> -		<b></b>
#1	15228						
#2	15310		<del></del> -				

Operator: RJG

Analysis Report 05/17/00 11:36:55 AM page 1

Method: METTRA Sample Name: CCV3-4

Run Time: 05/17/00 11:32:49

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.0068	23.527	.50734	1.9302	1.9469	48.584	.47261
SDev	.0076	.212	.00819	.0085	.0196	.463	.00602
%RSD	.75848	.90247	1.6149	.44117	1.0050	.95291	1.2736
#1	1.0122	23.677	.51313	1.9362	1.9607	48.912	.47687
#2	1.0014	23.377	.50155	1.9241	1.9330	48.257	.46835
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	27.500	.55000	2.2000	2.2000	55.000	.55000
Low	.90000	22.500	.45000	1.8000	1.8000	45.000	.45000
Elem Units Avge SDev %RSD	CO ppm 1.9396 .0204 1.0523	CR ppm 1.9474 .0203 1.0428	CU ppm 1.9478 .0174 .89309	FE ppm 23.754 .299 1.2575	MG ppm 47.802 .494 1.0340	MN ppm 1.9144 .0177	MO ppm 1.9687 .0087 .44119
#1	1.9541	1.9618	1.9601	23.965	48.151	1.9269	1.9748
#2	1.9252	1.9331	1.9355	23.543	47.452	1.9018	1.9625
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.2000	2.2000	2.2000	27.500	55.000	2.2000	2.2000
Low	1.8000	1.8000	1.8000	22.500	45.000	1.8000	1.8000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.9909	.47814	.48776	.48455	.50293	.50055	.50134
SDev	.0272	.00562	.00544	.00550	.00800	.00956	.00904
%RSD	1.3650	1.1761	1.1161	1.1358	1.5908	1.9106	1.8038
#1	2.0101	.48212	.49161	.48845	.50858	.50731	.50774
#2	1.9716	.47416	.48391	.48066	.49727	.49379	.49495
Errors High Low	LC Pass 2.2000 1.8000	NOCHECK	NOCHECK	LC Pass .55000 .45000	NOCHECK	NOCHECK	LC Pass .55000 .45000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.50682	.51184	.51017	1.0060	1.9482	2.0046	
SDev	.01011	.00533	.00692	.0183	.0200	.0273	
%RSD	1.9949	1.0409	1.3565	1.8236	1.0266	1.3642	
#1	.51397	.51561	.51506	1.0189	1.9623	2.0239	
#2	.49967	.50807	.50528	.99298	1.9341	1.9852	
Errors High Low	NOCHECK	NOCHECK	LC Pass .55000 .45000	LC Pass 1.1000 .90000	LC Pass 2.2000 1.8000	LC Pass 2.2000 1.8000	

05/17/00 11:36:55 AM

page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16764 157.8257 .9414532	2 NOTUSED    	3 NOTUSED	4 NOTUSED    	5 NOTUSED   	6 NOTUSED	7 NOTUSED   	
#1	16652					<del></del>		
#2	16876	· ·						

page 1

Method: METTRA Sample Name: CCB4 Run Time: 05/17/00 11:36:59 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00001	.04003	00006	.00055	.00135	.05749	00000
SDev	.00044	.00213	.00148	.00020	.00010	.02707	.00011
%RSD	4115.5	5.3298	2300.2	37.217	7.5159	47.083	170530.
#1	00030	.04154	00111	.00040	.00142	.03835	.00008
#2	.00033	.03852	.00098	.00069	.00128	.07663	00008
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.01000	.20000	.00500	5.0000	.00500
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00054	.00000	00148	.00457	.01346	.00051	.00599
SDev	.00012	.00008	.00008	.00593	.00415	.00016	.00191
%RSD	22.553	1815.6	5.5052	129.65	30.821	30.777	31.960
#1	00045	.00006	00153	.00038	.01052	.00040	.00734
#2	00062	00005	00142	.00877	.01639	.00062	
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.05000	.01000	.02500	.10000	5.0000	.01500	.04000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00074	.00026	.00103	.00077	.00180	.00083	.00115
SDev	.00095	.00024	.00090	.00052	.00065	.00123	.00104
%RSD	128.18	91.914	87.328	67.129	36.045	149.22	90.206
#1	00007	.00009	.00167	.00114	.00226	.00170	.00189
#2	00142		.00039	.00041	.00134	00005	.00042
Errors High Low	LC Pass .04000 04000	NOCHECK	NOCHECK	LC Pass .00300 00300	NOCHECK	NOCHECK	LC Pass .06000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00511	.00342	.00398	.00594	.00077	.00098	
SDev	.00215	.00014	.00081	.00029	.00036	.00019	
%RSD	42.124	4.1322	20.361	4.9010	46.905	19.851	
#1	.00359	.00332	.00341	.00573	.00103	.00084	
#2	.00663	.00352	.00456	.00614	.00052	.00111	
Errors High Low	NOCHECK	NOCHECK	LC Pass .00500 00500	LC Pass .01000 01000	LC Pass .05000 05000	LC Pass .02000 02000	

page 2

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Wavlen	371.030				w =		
Avge	16824					<del>-</del> -	
SDev	7.354186		<b></b>				
%RSD	.0437131	<del></del>					
#1	16819	<b></b> -		<b></b>			
#2	16829			~ ~			

page 1

Method: METTRA Sample Name: DCVMK Run Time: 05/17/00 11:41:08 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00039	.68116	.00057	.08003	.00063	176.00	00009
SDev	.00038	.01246	.00042	.00045	.00008	1.46	.00003
%RSD	96.740	1.8295	73.428	.55839	12.480	.82775	35.376
#1	00066	.68997	.00087	.08035	.00068	177.03	00006
#2	00012	.67235	.00028	.07972	.00057	174.97	00011
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00070	.00204	.00012	7.1650	5.0511	.05463	.00347
SDev	.00018	.00011	.00038	.0413	.0373	.00040	.00011
%RSD	25.836	5.2248	328.60	.57637	.73762	.73869	3.1542
#1 #2	00083 00057	.00196	00015 .00039	7.1942 7.1358	5.0774 5.0247	.05492 .05435	.00339
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00011	.00214	.00026	.00089	.00054	.00104	.00088
SDev	.00031	.00096	.00065	.00012	.00316	.00010	.00099
%RSD	287.20	44.722	248.31	13.089	581.62	9.6189	112.58
#1	00032	.00146	.00072	.00097	00169	.00111	.00018
#2	.00011	.00282	00020	.00081	.00278	.00097	.00157
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm .	
Avge	00209	.00158	.00036	.00079	.00293	.01447	
SDev	.00069	.00564	.00399	.00126	.00002	.00004	
%RSD	32.949	357.66	1115.9	158.70	.61201	.29366	
#1	00160	.00557	.00318	.00168	.00294	.01450	
#2	00257	00241	00247	00010	.00292	.01444	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

05/17/00 11:45:14 AM page 2

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Wavlen	371.030						
Avge	16317		<b></b>				
SDev	110.1319			<b></b>			<del>-</del> -
%RSD	.6749404				<del>-</del> -		
#1	16239						
#2	16395				<b></b>		

page 1

Method: METTRA Sample Name: DCVML Run Time: 05/17/00 11:45:18 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00024	.14363	.00152	.15260	.00004	402.42	.00001
SDev	.00022	.01289	.00007	.00005	.00021	2.33	.00002
%RSD	91.098	8.9711	4.4938	.03440	510.93	.57809	246.26
#1	00009	.15274	.00147	.15256	.00019	404.07	00001
#2	00040	.13451	.00157	.15264	00011	400.78	.00002
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00082	.00224	00076	1.8869	38.876	1.0827	.00179
SDev	.00003	.00029	.00057	.0099	.103	.0019	.00023
%RSD	3.2463	13.116	74.408	.52256	.26616	.17171	13.127
#1	.00084	.00204	00116	1.8939	38.949	1.0841	.00162
#2		.00245	00036	1.8799	38.803	1.0814	.00195
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00199	00038	.00268	.00166	.00301	.00012	.00108
SDev	.00006	.00254	.00078	.00136	.00107	.00066	.00080
%RSD	3.1585	664.20	29.001	82.350	35.657	550.48	73.537
#1	.00194	.00142	.00322	.00262	.00225	00034	.00052
#2	.00203	00218	.00213	.00069	.00376	.00058	.00164
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00077	.00267	.00153	00042	.00312	.00863	
SDev	.00262	.00057	.00049	.00527	.00059	.00003	
%RSD	341.38	21.261	32.393	1268.8	18.929	.38331	
#1 #2	00262 .00109	.00308	.00118 .00188	.00331 00414	.00353 .00270	.00866 .00861	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Report

05/17/00 11:49:23 AM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 15637 8.414847 .0538132	2 NOTUSED   	3 NOTUSED    	4 NOTUSED    	5 NOTUSED   	6 NOTUSED    	7 NOTUSED    
#1	15631						
#2	15643						

Method: METTRA Sample Name: DCVMM Run Time: 05/17/00 11:49:27 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00001	.78762	.00107	.06784	.00042	114.96	00029
SDev	.00022	.01692	.00066	.00047	.00004	1.38	.00001
%RSD	1993.3	2.1485	61.779	.69456	8.7969	1.2004	1.8596
#1	.00015	.79958	.00060	.06818	.00044	115.94	00029
#2	00017	.77565	.00153	.06751	.00039	113.98	00030
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00091	.00264	.00060	3.3573	4.7846	.03918	.00245
SDev	.00036	.00001	.00019	.0296	.0537	.00038	.00005
%RSD	40.160	.18158	31.673	.88225	1.1232	.97303	2.0186
#1	00117	.00264	.00046	3.3783	4.8226	.03945	.00249
#2	00065	.00263		3.3364	4.7466	.03891	.00242
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00209	.00119	.00143	.00135	.00120	.00126	.00124
SDev	.00041	.00027	.00088	.00050	.00159	.00001	.00052
%RSD	19.705	22.659	61.403	36.721	132.93	.63550	42.281
#1 #2	.00180	.00138 .00100	.00081 .00205	.00100 .00170	.00007	.00127 .00126	.00087 .00161
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00798	.00073	00217	00379	.00408	.02289	
SDev	.00253	.00076	.00135	.00126	.00004	.00047	
%RSD	31.659	103.28	62.045	33.103	.93837	2.0698	
#1	00977	.00020	00312	00291	.00411	.02322	
#2	00620	.00127	00122	00468	.00406	.02255	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	Y						
Wavlen	371.030						
Avge	16160						
SDev	154.1134			900 HA			
%RSD	.9536969						
#1	16051			<del>-</del> -			
#2	16269	<b></b>			<del>-</del> -		

Method: METTRA Sample Name: DCVMN Run Time: 05/17/00 11:53:36 Method: METTRA Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00010	.46389	.00108	.07426	.00002	354.44	.00005
SDev	.00007	.01720	.00169	.00006	.00019	1.57	.00001
%RSD	69.790	3.7069	156.66	.08517	859.62	.44357	26.966
#1	00005	.47605	00012	.07421	.00015	355. <b>55</b>	.00004
#2	00015	.45173	.00228	.07430	00011	353.33	.00006
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00129	.00115	.00019	2.0278	47.936	.87434	.00192
SDev	.00010	.00031	.00022	.0065	.117	.00124	.00007
%RSD	7.9241	26.535	118.88	.31898	.24376	.14169	3.4331
#1	.00122	.00094	.00003	2.0233	48.019	.87522	.00187
#2	.00136	.00137		2.0324	47.854	.87346	.00197
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00195	00046	.00065	.00028	.00273	.00067	.00136
SDev	.00128	.00171	.00074	.00106	.00012	.00313	.00213
%RSD	65.603	369.86	115.03	385.28	4.3695	465.26	156.46
#1	.00105	.00075	.00117	.00103	.00265	00154	00014
#2	.00286	00167	.00012	00048	.00282	.00288	.00286
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00030	00047	00021	.00069	.00580	.01159	
SDev	.00128	.00076	.00093	.00443	.00038	.00004	
%RSD	420.34	162.88	445.16	639.51	6.5325	.34538	
#1	00060	00100	00087	00244	.00553	.01162	
#2	.00121	.00007	.00045	.00383	.00607	.01156	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Report

05/17/00 11:57:42 AM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem Wavlen	Y 371.030		- ··				
Avge	15561						
SDev	44.83015		<b></b>				
%RSD	.2881004			<del></del>			
	15500				<del></del>		
#1	15529 15592				<del>-</del> -		
#2	10024						

page 1

Method: METTRA Sample Name: DCVMP Run Time: 05/17/00 11:57:46 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00046	.05453	.00039	.07369	.00042	93.203	00018
SDev	.00004	.00909	.00041	.00002	.00002	.654	.00021
%RSD	9.1752	16.662	104.64	.03088	3.9252	.70163	113.89
#1	00049	.06096	.00010	.07371	.00041	93.665	00004
#2	00043	.04811	.00068	.07368		92.740	00033
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00114	.00167	00007	1.9055	3.5397	.05770	.00349
SDev	.00069	.00029	.00004	.0169	.0173	.00032	.00011
%RSD	60.912	17.634	58.199	.88893	.48779	.55899	3.1413
#1	00065	.00187	00004	1.9175	3.5519	.05793	.00357
#2	00163	.00146	00009	1.8935	3.5275	.05747	.00341
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem Units Avge SDev %RSD	NI ppm .00132 .00062 46.982	PB/1 ppm .00289 .00003 1.0943	PB/2 ppm .00090 .00018 20.468	PB ppm .00157 .00011 7.2004	SB/1 ppm .00082 .00014 17.005	SB/2 ppm .00168 .00008 4.6378	SB ppm .00139 .00001
#1	.00176	.00287	.00103	.00165	.00092	.00162	.00139
#2	.00088	.00292	.00077	.00149	.00072	.00173	.00139
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem							
Units Avge SDev %RSD	SE/1 ppm 00528 .00080 15.091	SE/2 ppm .00197 .00089 44.871	SE ppm 00044 .00086 194.19	TL ppm 00260 .00237 91.339	V_ ppm .00050 .00000 .88892	ZN ppm .01175 .00017 1.4625	
Avge SDev	ppm 00528 .00080	ppm .00197 .00089	ppm 00044 .00086	ppm 00260 .00237	ppm .00050 .00000	ppm .01175 .00017	

657 730

Analysis Report

05/17/00 12:01:51 PM

page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	Y						
Wavlen	371.030						
Avge	16415						
SDev	90.50967						
%RSD	.5513957				· ·		
#1	16351						
#2	16479						

Method: METTRA Sample Name: DD8DGBF Operator: RJG

Run Time: 05/17/00 12:01:55

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00016	.04059	.00020	.00004	.00030	.11829	00008
SDev	.00049	.01620	.00177	.00004	.00011	.00288	.00009
%RSD	312.94	39.916	909.35	96.579	35.062	2.4386	102.03
#1	.00019	.05205	.00145	.00001	.00037	.12033	00002
#2	00051	.02913	00106	.00007		.11625	00015
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.01000	.20000	.00500	5.0000	.00500
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00094	.00007	00107	.02213	00034	.00001	.00005
SDev	.00000	.00004	.00021	.00464	.00116	.00001	.00029
%RSD	.43352	53.139	19.532	20.963	339.99	119.76	553.46
#1	00094	.00010	00121	.01885	.00048	.00000	00015
#2	00093	.00005	00092	.02541	00116		.00026
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.05000	.01000	.02500	.10000	5.0000	.01500	.04000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00111	.00095	.00177	.00149	.00193	.00097	.00129
SDev	.00050	.00088	.00039	.00003	.00010	.00289	.00197
%RSD	44.757	92.654	21.903	2.2518	5.3801	297.51	152.08
#1	00146	.00157	.00149	.00152	.00201	.00302	.00268
#2	00076	.00033	.00204	.00147	.00186	00107	00010
Errors High Low	LC Pass .04000 04000	NOCHECK	NOCHECK	LC Pass .00300 00300	NOCHECK	NOCHECK	LC Pass .06000 06000
Elem	SE/1	SE/2	SE	TL	V <u></u>	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00448	.00135	00059	00337	.00067	.00240	
SDev	.00414	.00136	.00228	.00059	.00019	.00021	
%RSD	92.364	100.21	388.19	17.616	27.860	8.6873	
#1	00155	.00231	.00103	00379	.00054	.00254	
#2	00740	.00039	00220	00295	.00080	.00225	
Errors High Low	NOCHECK	NOCHECK	LC Pass .00500 00500	LC Pass .01000 01000	LC Pass .05000 05000	LC Pass .02000 02000	

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16273 61.76550 .3795553	NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED    	6 NOTUSED	7 NOTUSED    
#1	16229						ru
#2	16317						

5130

Sample Name: DD8DGCF Method: METTRA Method: METTRA Sample Na Run Time: 05/17/00 12:06:05 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04877	2.0099	2.0111	1.9691	.04972	47.938	.04721
SDev	.00007	.0550	.0471	.0383	.00112	1.063	.00109
%RSD	.15257	2.7344	2.3412	1.9434	2.2561	2.2176	2.3109
#1	.04882	2.0488	2.0443	1.9962	.05051	48.690	.04798
#2	.04871	1.9710	1.9778	1.9421	.04892	47.186	.04644
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.06000	2.4000	2.4000	2.4000	.06000	60.000	.06000
Low	.04000	1.6000	1.6000	1.6000	.04000	40.000	.04000
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.48628	.19764	.24638	.89662	47.285	.48257	.99408
SDev	.00988	.00354	.00398	.02694	.961	.00931	.01886
%RSD	2.0322	1.7937	1.6174	3.0049	2.0331	1.9288	1.8974
#1	.49327	.20014	.24919	.91567	47.965	.48915	1.0074
#2	.47929	.19513	.24356	.87757	46.605	.47599	.98074
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.60000	.24000	.30000	1.2000	60.000	.60000	1.2000
Low	.40000	.16000	.20000	.80000	40.000	.40000	.80000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.49350	.47724	.47510	.47581	.47762	.49181	.48708
SDev	.01322	.00549	.00108	.00110	.00419	.00909	.00467
%RSD	2.6796	1.1496	.22806	.23208	.87785	1.8489	.95853
#1	.50285	.48112	.47433	.47659	.47465	.49824	.49038
#2	.48415	.47336	.47587	.47503	.48058	.48538	.48378
Errors High Low	LC Pass .60000 .40000	NOCHECK	NOCHECK	LC Pass .60000 .40000	NOCHECK	NOCHECK	LC Pass .60000 .40000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	2.0080	1.9429	1.9646	2.0319	.49279	.51023	
SDev	.0208	.0219	.0077	.0509	.01131	.01120	
%RSD	1.0366	1.1286	.39165	2.5066	2.2955	2.1955	
#1	2.0228	1.9273	1.9591	2.0679	.50079	.51815	
#2	1.9933	1.9584	1.9700	1.9959	.48479	.50231	
Errors High Low	NOCHECK	NOCHECK	LC Pass 2.4000 1.6000	LC Pass 2.4000 1.6000	LC Pass .60000 .40000	LC Pass .60000 .40000	

05/17/00 12:10:11 PM page 2

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED 	7 NOTUSED
Wavlen	371.030						
Avge	16323					- <del>-</del>	
SDev	238.2239						
%RSD	1.459423			<del></del>			
#1	16155				•• ••		
#1 #2	16492						

Method: METTRA Sample Name: DCQF8F Operator: RJG

Run Time: 05/17/00 12:10:14

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units Avge	ppm 00005	ppm .03836	ppm .00063	ppm .03030	ppm .00023	ppm 104.85	ppm 00011
SDev %RSD	.00083	.01198	.00085	.00005	.00001	.40	.00029
	1640.5	-	134.62	.17506	3.6804	.38436	257.61
#1 #2	.00053 00064	.04683 .02989	.00124 .00003	.03034 .03026	.00023 .00024	105.14 104.57	.00009 00031
Errors High Low	LC Pass 2.0000 01000	LC Pass 600.00 20000	LC Pass 10.000 01000	LC Pass 10.000 20000	LC Pass 10.000 00500	LC Pass 600.00 -5.0000	LC Pass 5.0000 00500
Elem Units	CO	CR	CU	FE	MG	MN	МО
Avge	ppm 00091	ppm .00013	ppm 00043	ppm 4.1638	ppm 4.1692	ppm .19935	ppm .00472
SDev %RSD	.00020 21.571	.00013 100.26	.00008 17.392	.0139 .33489	.0147 .35298	.00040 .20009	.00189 39.945
#1	~.00077	.00022	00038	4,1737	4.1796	.19964	.00606
#2	00104	.00004	00049	4.1540	4.1588	.19907	.00339
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High Low	100.00 05000	20.000 01000	10.000 02500	500.00 10000	600.00 -5.0000	10.000 01500	20.000 04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Units Avge SDev	ppm .00090 .00012	ppm .00140 .00179	ppm .00194 .00108	ppm .00176 .00131	ppm .00298 .00262	ppm .00024 .00080	ppm .00115 .00140
Units Avge SDev %RSD	ppm .00090 .00012 12.737	ppm .00140 .00179 128.09	ppm .00194 .00108 55.427	ppm .00176 .00131 74.648	ppm .00298 .00262 88.081	ppm .00024 .00080 336.02	ppm .00115 .00140 122.23
Units Avge SDev	ppm .00090 .00012	ppm .00140 .00179	ppm .00194 .00108	ppm .00176 .00131	ppm .00298 .00262	ppm .00024 .00080	ppm .00115 .00140
Units Avge SDev %RSD #1 #2 Errors	ppm .00090 .00012 12.737 .00082 .00098	ppm .00140 .00179 128.09	ppm .00194 .00108 55.427	ppm .00176 .00131 74.648 .00269 .00083	ppm .00298 .00262 88.081	ppm .00024 .00080 336.02	ppm .00115 .00140 122.23 .00214 .00016
Units Avge SDev %RSD #1 #2	ppm .00090 .00012 12.737 .00082 .00098	ppm .00140 .00179 128.09 .00266 .00013	ppm .00194 .00108 55.427 .00270 .00118	ppm .00176 .00131 74.648 .00269 .00083	ppm .00298 .00262 88.081 .00483 .00112	ppm .00024 .00080 336.02 .00080 00033	ppm .00115 .00140 122.23 .00214 .00016
Units Avge SDev %RSD #1 #2 Errors High	ppm .00090 .00012 12.737 .00082 .00098 LC Pass 100.00	ppm .00140 .00179 128.09 .00266 .00013	ppm .00194 .00108 55.427 .00270 .00118	ppm .00176 .00131 74.648 .00269 .00083 LC Pass 5.0000	ppm .00298 .00262 88.081 .00483 .00112	ppm .00024 .00080 336.02 .00080 00033	ppm .00115 .00140 122.23 .00214 .00016 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low Elem Units	ppm .00090 .00012 12.737 .00082 .00098 LC Pass 100.00 04000 SE/1 ppm	ppm .00140 .00179 128.09 .00266 .00013 NOCHECK	ppm .00194 .00108 55.427 .00270 .00118 NOCHECK	ppm .00176 .00131 74.648 .00269 .00083 LC Pass 5.0000 00300	ppm .00298 .00262 88.081 .00483 .00112 NOCHECK	ppm .00024 .00080 336.02 .00080 00033 NOCHECK	ppm .00115 .00140 122.23 .00214 .00016 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge SDev	ppm .00090 .00012 12.737 .00082 .00098 LC Pass 100.00 04000 SE/1 ppm 00387 .00278	ppm .00140 .00179 128.09 .00266 .00013 NOCHECK SE/2 ppm .00479 .00076	ppm .00194 .00108 55.427 .00270 .00118 NOCHECK SE ppm .00191 .00042	ppm .00176 .00131 74.648 .00269 .00083 LC Pass 5.0000 00300 TL ppm .00055 .00137	ppm .00298 .00262 88.081 .00483 .00112 NOCHECK V_ppm .00069 .00057	ppm .00024 .00080 336.02 .00080 00033 NOCHECK ZN ppm .00510 .00009	ppm .00115 .00140 122.23 .00214 .00016 LC Pass 10.000
Units Avge SDev *RSD #1 #2 Errors High Low Elem Units Avge SDev *RSD	ppm .00090 .00012 12.737 .00082 .00098 LC Pass 100.00 04000 SE/1 ppm 00387	ppm .00140 .00179 128.09 .00266 .00013 NOCHECK SE/2 ppm .00479	ppm .00194 .00108 55.427 .00270 .00118 NOCHECK	ppm .00176 .00131 74.648 .00269 .00083 LC Pass 5.0000 00300 TL ppm .00055	ppm .00298 .00262 88.081 .00483 .00112 NOCHECK	ppm .00024 .00080 336.02 .00080 00033 NOCHECK	ppm .00115 .00140 122.23 .00214 .00016 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge SDev	ppm .00090 .00012 12.737 .00082 .00098 LC Pass 100.00 04000 SE/1 ppm 00387 .00278	ppm .00140 .00179 128.09 .00266 .00013 NOCHECK SE/2 ppm .00479 .00076	ppm .00194 .00108 55.427 .00270 .00118 NOCHECK SE ppm .00191 .00042	ppm .00176 .00131 74.648 .00269 .00083 LC Pass 5.0000 00300 TL ppm .00055 .00137	ppm .00298 .00262 88.081 .00483 .00112 NOCHECK V_ppm .00069 .00057	ppm .00024 .00080 336.02 .00080 00033 NOCHECK ZN ppm .00510 .00009	ppm .00115 .00140 122.23 .00214 .00016 LC Pass 10.000
Units Avge SDev RSD #1 #2 Errors High Low Elem Units Avge SDev RSD #1	ppm .00090 .00012 12.737 .00082 .00098 LC Pass 100.00 04000 SE/1 ppm 00387 .00278 71.952 00190	ppm .00140 .00179 128.09 .00266 .00013 NOCHECK SE/2 ppm .00479 .00076 15.959	ppm .00194 .00108 55.427 .00270 .00118 NOCHECK SE ppm .00191 .00042 21.865	ppm .00176 .00131 74.648 .00269 .00083 LC Pass 5.0000 00300 TL ppm .00055 .00137 246.92	ppm .00298 .00262 88.081 .00483 .00112 NOCHECK V_ppm .00069 .00057 82.466	ppm .00024 .00080 336.02 .00080 00033 NOCHECK ZN ppm .00510 .00009 1.8420	ppm .00115 .00140 122.23 .00214 .00016 LC Pass 10.000
Units Avge SDev RSD #1 #2 Errors High Low Elem Units Avge SDev RSD #1 #2	ppm .00090 .00012 12.737 .00082 .00098 LC Pass 100.00 04000 SE/1 ppm 00387 .00278 71.952 00190 00584	ppm .00140 .00179 128.09 .00266 .00013 NOCHECK SE/2 ppm .00479 .00076 15.959 .00425 .00533	ppm .00194 .00108 55.427 .00270 .00118 NOCHECK SE ppm .00191 .00042 21.865 .00220	ppm .00176 .00131 74.648 .00269 .00083 LC Pass 5.0000 00300 TL ppm .00055 .00137 246.92 .00152 00041	ppm .00298 .00262 88.081 .00483 .00112 NOCHECK V_ppm .00069 .00057 82.466 .00109 .00029	ppm .00024 .00080 336.02 .00080 00033 NOCHECK ZN ppm .00510 .00009 1.8420 .00517 .00504	ppm .00115 .00140 122.23 .00214 .00016 LC Pass 10.000

IntSto Mode Elem Wavler Avge SDev %RSD	Counts Y	2 NOTUSED    	3 NOTUSED    	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    	
#1 #2	16093 16256		· · ·		~ ~			

Method: METTRA Sample Name: DCQFLF Run Time: 05/17/00 12:14:24 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG ppm00021 .00018 82.691	AL	AS	BA	BE	CA	CD
Units		ppm	ppm	ppm	ppm	ppm	ppm
Avge		.04111	.00131	.02204	.00010	89.989	00031
SDev		.00154	.00081	.00037	.00019	1.097	.00004
%RSD		3.7539	61.696	1.6775	195.65	1.2193	12.527
#1	00009	.04221	.00074	.02178	.00024	89.213	00033
#2	00034	.04002	.00188	.02230	00004	90.765	00028
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00081	00022	00064	1.9577	4.7103	.18299	.00255
SDev	.00030	.00024	.00019	.0319	.0628	.00265	.00014
%RSD	36.812	107.43	28.965	1.6303	1.3341	1.4477	5.3031
#1	00102	00005	00078	1.9351	4.6659	.18112	.00245
#2	00060	00039	00051	1.9803	4.7548	.18486	.00264
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00055	.00082	.00163	.00136	.00090	.00188	.00155
SDev	.00149	.00307	.00067	.00057	.00140	.00040	.00073
%RSD	269.41	372.77	41.278	42.041	155.45	21.536	47.422
#1	.00161	.00299	.00116	.00177	.00189	.00216	.00207
#2	00050	00135	.00211	.00096	00009	.00159	.00103
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00447	.00153	00047	00129	.00105	.02417	
SDev	.00025	.00004	.00006	.00218	.00001	.00007	
%RSD	5.6214	2.6288	12.193	169.38	1.3574	.28579	
#1	00429	.00150	00043	00283	.00104	.02412	
#2	00465	.00156	00051	.00025	.00106	.02422	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

IntStd Mode Elem Wavlen Avge	1 Counts Y 371.030 16032 176.1407	2 NOTUSED  	3 NOTUSED  	4 NOTUSED	5 NOTUSED  	6 NOTUSED  	7 NOTUSED  
SDev %RSD	1.098710				- ·-		
#1 #2	16156 15907	<b></b>			<del>-</del> -		<b></b>
#4	13307	<b></b>					

page 1

Method: METTRA Sample Name: DCQFRF Run Time: 05/17/00 12:18:34 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00030	.02713	00058	.04075	.00038	118.02	00019
SDev	.00005	.01577	.00195	.00136	.00022	4.00	.00004
%RSD	17.732	58.125	337.42	3.3317	58.341	3.3909	20.759
#1	00033	.03829	00195	.04171	.00022	120.85	00022
#2	00026	.01598	.00080	.03979	.00054	115.19	00016
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00106	00006	00037	5.5379	4.7809	.10579	.00223
SDev	.00027	.00016	.00019	.1605	.1462	.00342	.00018
%RSD	25.425	261.49	51.982	2.8987	3.0573	3.2307	7.9673
#1	00087	00017	00023	5.6514	4.8843	.10820	.00211
#2	00125	.00005	00050	5.4244	4.6776	.10337	.00236
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00041	.00145	.00182	.00170	.00236	00082	.00024
SDev	.00012	.00042	.00034	.00009	.00138	.00122	.00036
%RSD	30.092	28.580	18.888	5.3701	58.544	149.35	147.67
#1	00050	.00116	.00207	.00177	.00334	00168	00001
#2	00032	.00175	.00158	.00164	.00139	.00005	.00049
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00717	.00265	00062	00035	.00066	.01010	
SDev	.00156	.00338	.00174	.00223	.00002	.00030	
%RSD	21.691	127.74	279.46	628.92	3.4423	2.9223	
#1	00827	.00504	.00061	00193	.00068	.01030	
#2	00607	.00026	00185	.00122	.00064	.00989	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

05/17/00 12:22:39 PM page 2

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED 	4 NOTUSED	5 NOTUSED 	6 NOTUSED	7 NOTUSED
Wavlen	371.030	<b>-</b>			m		
Avge	16512	<b></b>					
SDev	403.1572						
%RSD	2.441590					·· ·	
#1	16227						
#2	16797						

Operator: RJG

Method: METTRA Sample Name: CCV3-5

Run Time: 05/17/00 12:22:43

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

			-				
Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.0396	24.645	.52236	2.0086	1.9973	49.771	.48674
SDev	.0007	.027	.00241	.0011	.0063	.246	.00139
%RSD	.06807	.11070	.46211	.05590	.31323	.49360	.28517
#1	1.0391	24.626	.52407	2.0093	2.0017	49.945	.48772
#2	1.0401	24.664	.52066	2.0078	1.9928	49.597	.48576
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	27.500	.55000	2.2000	2.2000	55.000	.55000
Low	.90000	22.500	.45000	1.8000	1.8000	45.000	.45000
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.9998	2.0101	2.0185	24.683	49.242	1.9802	2.0230
SDev	.0045	.0032	.0030	.018	.073	.0008	.0114
%RSD	.22637	.16164	.15079	.07482	.14799	.03844	.56274
#1	2.0030	2.0124	2.0206	24.670	49.294	1.9808	2.0149
#2	1.9966	2.0078	2.0163	24.696	49.191	1.9797	2.0310
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.2000	2.2000	2.2000	27.500	55.000	2.2000	2.2000
Low	1.8000	1.8000	1.8000	22.500	45.000	1.8000	1.8000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	2.0367	.49350	.50180	.49904	.51536	.52015	.51856
SDev	.0234	.00331	.00096	.00047	.00136	.00324	.00171
%RSD	1.1492	.67110	.19037	.09331	.26386	.62275	.32933
#1	2.0532	.49116	.50248	.49871	.51633	.51786	.51735
#2	2.0201	.49584	.50113	.49937	.51440	.52244	.51977
Errors High Low	LC Pass 2.2000 1.8000	NOCHECK	NOCHECK	LC Pass .55000 .45000	NOCHECK	NOCHECK	LC Pass .55000 .45000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.52855	.52893	.52880	1.0257	2.0012	2.0658	
SDev	.00113	.00123	.00044	.0081	.0124	.0139	
%RSD	.21321	.23258	.08421	.78617	.61800	.67203	
#1	.52775	·.52980	.52912	1.0314	2.0100	2.0756	
#2	.52934	.52806	.52849	1.0200	1.9925	2.0559	
Errors High Low	NOCHECK	NOCHECK	LC Pass .55000 .45000	LC Pass 1.1000 .90000	LC Pass 2.2000 1.8000	LC Pass 2.2000 1.8000	

05/17/00 12:26:49 PM page 2

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED 	6 NOTUSED 	7 NOTUSED
Wavlen	371.030						
Avge	16341						
SDev	80.85808						
%RSD	.4948104						
#1	16398						
#2	16284						

Method: METTRA Sample Name: CCB5 Operator: Run Time: 05/17/00 12:26:53 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP Operator: RJG

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00004	.03260	00031	.00036	.00064	.02959	.00008
SDev	.00011	.01196	.00192	.00007	.00018	.00349	.00005
%RSD	279.58	36.692	614.49	18.328	28.026	11.784	56.756
#1	00011	.04106	.00104	.00031	.00077	.02712	.00005
#2	.00004	.02414	00167	.00041	.00051	.03205	.00012
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.01000	.20000	.00500	5.0000	.00500
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00043	.00015	00043	.01109	.01163	.00042	.00644
SDev	.00010	.00006	.00007	.00324	.00342	.00006	.00249
%RSD	23.375	40.559	17.131	29.200	29.439	14.302	38.689
#1	00050	.00020	00048	.00880	.01405	.00047	.00820
#2	00036	.00011	00038	.01337	.00921	.00038	.00468
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.05000	.01000	.02500	.10000	5.0000	.01500	.04000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00075	00259	.00099	00021	.00123	.00050	.00074
SDev	.00049	.00325	.00042	.00080	.00191	.00169	.00176
%RSD	64.606	125.31	42.578	389.74	155.13	336.82	236.81
#1	00041	00030	.00069	.00036	.00258	.00170	.00199
#2	00110	00489	.00128	00077	00012	00069	00050
Errors High Low	LC Pass .04000 04000	NOCHECK	NOCHECK	LC Pass .00300 00300	NOCHECK	NOCHECK	LC Pass .06000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00044	.00267	.00193	.00279	.00066	.00084	
SDev	.00100	.00073	.00016	.00006	.00019	.00008	
%RSD	224.84	27.438	8.1241	2.3035	28.204	9.9138	
#1	00026	.00319	.00204	.00284	.00079	.00089	
#2	.00115	.00215	.00182	.00275	.00053	.00078	
Errors High Low	NOCHECK	NOCHECK	LC Pass .00500 00500	LC Pass .01000 01000	LC Pass .05000 05000	LC Pass .02000 02000	

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Wavlen	371.030						
Avge	16467						
SDev	29.66258						
%RSD	.1801343	~ **					way wa
#1	16446	~ ~					
#2	16488						

Method: METTRA Sample Name: DCQFXF Operator: RJG

Run Time: 05/17/00 12:31:03

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00014	.04095	00027	.06033	.00009	158.01	00030
SDev	.00038	.00745	.00034	.00010	.00002	.56	.00009
%RSD	264.82	18.200	125.76	.16388	22.661	.35287	29.277
#1	00012	.04623	00050	.06026	.00011	158.40	00036
#2	.00041	.03568	00003	.06040		157.61	00024
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00102	00002	.00010	11.137	8.8984	.05492	.00347
SDev	.00005	.00014	.00045	.015	.0162	.00006	.00048
%RSD	5.3504	789.84	449.45	.13607	.18246	.11606	13.871
#1	00106	.00008	00022	11.148	8.9099	.05488	.00382
#2	00099	00012	.00041	11.126	8.8869	.05497	.00313
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00043	.00147	.00153	.00151	.00023	.00095	.00071
SDev	.00169	.00195	.00042	.00037	.00467	.00257	.00016
%RSD	395.57	132.76	27.571	24.396	2059.2	269.95	22.459
#1	00077	.00009	.00183	.00125	.00353	00087	.00060
#2	.00162	.00285	.00123	.00177	00307	.00277	.00082
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00415	.00387	.00120	00173	.00291	.00618	
SDev	.00114	.00097	.00103	.00012	.00000	.00006	
%RSD	27.549	25.091	85.615	7.1487	.13606	.91998	
#1	00334	.00456	.00193	00164	.00291	.00614	
#2	00496	.00319	.00047	00182	.00290	.00622	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16031 50.69970	2 NOTUSED   	3 NOTUSED   	4 NOTUSED   	5 NOTUSED   	6 NOTUSED   	7 NOTUSED   
#1 #2	15995 16067					 	

Operator: RJG

Method: METTRA Sample Name: DCQG5F Run Time: 05/17/00 12:35:12

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00068	.06391	.00201	.14392	00058	H691.20	00010
SDev	.00094	.00763	.00009	.00013	.00001	1.99	.00006
%RSD	138.49	11.931	4.3344	.09243	1.3227	.28724	65.347
#1	.00134	.06931	.00207	.14401	00059	H692.60	00015
#2		.05852	.00195	.14382	00058	H689.80	00005
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem Units Avge SDev %RSD	CO ppm 00046 .00008 18.221	CR ppm .00096 .00051 52.935	CU ppm .00130 .00041 31.507	FE ppm 6.9193 .0001	MG ppm 43.450 .084 .19360	MN ppm .33273 .00030 .09132	MO ppm .00234 .00100 42.843
#1	00052	.00133	.00159	6.9193	43.509	.33294	.00304
#2	00040	.00060	.00101	6.9194	43.390	.33251	.00163
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00202	.00432	00184	.00021	00177	.00334	.00164
SDev	.00094	.00303	.00073	.00052	.00268	.00056	.00127
%RSD	46.408	70.191	39.807	247.87	151.11	16.753	77.416
#1	.00268	.00647	00236	.00058	00367	.00294	.00074
#2	.00136	.00218	00132	00016	.00012	.00373	
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00337	00069	.00067	.00220	.00181	.01089	
SDev	.00249	.00097	.00018	.00122	.00000	.00016	
%RSD	73.965	141.34	27.651	55.506	.00113	1.5022	
#1	.00513	00137	.00080	.00133	.00181	.01077	
#2	.00161	00000	.00054	.00306	.00181	.01100	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

657 748

Analysis Report

05/17/00 12:39:18 PM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 15155 68.58935 .4525901	2 NOTUSED    	3 NOTUSED    	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1 #2	15106 15203						

Method: METTRA Sample Name: DCQGKF Operator: RJG

Run Time: 05/17/00 12:39:22

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00005	.02371	.00072	.06587	00013	121.93	00020
SDev	.00045	.00337	.00153	.00050	.00008	.11	.00008
%RSD	850.60	14.196	210.88	.75766	64.001	.08975	38.376
#1	.00037	.02609	00036	.06552	00007	121.85	00026
#2	00026	.02133	.00180	.06622	00019	122.01	00015
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem Units Avge SDev %RSD	CO ppm .00681 .00028 4.0955	CR ppm .00020 .00009 43.440	CU ppm .00021 .00026 122.85	FE ppm 10.294 .026 .25189	MG ppm 9.1114 .0153 .16819	MN ppm .33014 .00075 .22815	MO ppm 00024 .00000
#1	.00661	.00026	.00003	10.276	9.1005	.32960	00024
#2	.00701	.00014		10.313	9.1222	.33067	00024
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01094	.00183	.00190	.00188	.00119	.00032	.00061
SDev	.00000	.00342	.00028	.00095	.00006	.00036	.00026
%RSD	.01532	186.96	14.827	50.771	5.1021	111.35	42.504
#1	.01095	.00425	.00170	.00255	.00115	.00007	.00043
#2	.01094	00059	.00210	.00120	.00123	.00058	
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00534	.00385	.00079	00287	.00094	.03057	
SDev	.00138	.00364	.00289	.00147	.00246	.00004	
%RSD	25.870	94.609	365.17	51.266	260.82	.11941	
#1	00631	.00127	00125	00183	.00268	.03059	
#2	00436	.00642	.00283	00392	00080	.03054	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

	Ana	lysis	Re	port
--	-----	-------	----	------

05/17/00 12:43:27 PM

page 2

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Wavlen	371.030						
Avge	16103		<b>-</b> -				
SDev	2.580525	<del>-</del> -				***	
%RSD	.0160254						
11 -	16101						
#1	16101				<b></b>		- <del>-</del>
#2	16104	** **					

05/17/00 12:47:37 PM

page 1

Method: METTRA Sample Name: DCQGMF Operator: RJG

Run Time: 05/17/00 12:43:31

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG ppm00018 .00030 165.16	AL	AS	BA	BE	CA	CD
Units		ppm	ppm	ppm	ppm	ppm	ppm
Avge		.03499	.01650	.07071	00028	153.98	~.00120
SDev		.00866	.00090	.00089	.00006	1.44	.00016
%RSD		24.760	5.4853	1.2579	20.950	.93225	13.620
#1	.00003	.04111	.017 <b>14</b>	.07008	00024	152.96	00132
#2	00040	.02886	.01586	.07134	00032	154.99	00109
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04378	.00051	00094	85.296	20.089	.51592	00065
SDev	.00054	.00000	.00012	.900	.210	.00539	.00045
%RSD	1.2307	.54994	13.200	1.0556	1.0471	1.0456	70.110
#1	.04340	.00051	00103	84.659	19.940	.51211	00033
#2	.04416	.00050	00085	85.933	20.238	.51973	00097
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.07364	00021	.00017	.00004	.00020	.00025	.00023
SDev	.00075	.00208	.00055	.00032	.00042	.00040	.00013
%RSD	1.0151	980.43	328.70	782.57	208.14	159.91	54.296
#1	.07311	00168	.00056	00019	00009	.00053	.00032
#2	.07417	.00126	00022	.00027	.00049	00003	.00014
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00481	.00384	.00096	00084	00032	.07055	
SDev	.00368	.00042	.00150	.00249	.00049	.00042	
%RSD	76.422	10.864	157.16	294.52	151.05	.59593	
#1	00741	.00354	00011	.00091	00067	.07025	
#2	00221	.00413	.00202	00260	.00002	.07085	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Report

05/17/00 12:47:37 PM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	Y						
Wavlen	371.030	<b></b>					
Avge	15877						~ _
SDev	92.70142						
%RSD	.5838632	<b>-</b> -					
		,		•			
#1	15943					<del>-</del> -	
#2	15812	<del>~</del> -	***				

page 1

Method: METTRA Sample Name: DCQGNF Run Time: 05/17/00 12:47:41 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00002	.02054	.00446	.03214	.00006	104.00	00043
SDev	.00002	.00346	.00009	.00028	.00007	.68	.00015
%RSD	79.589	16.828	2.0091	.88367	109.97	.65630	33.695
#1	00001	.02298	.00440	.03194	.00011	103.52	00053
#2	00003	.01809	.00453	.03234		104.49	00033
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01891	.00047	00071	18.040	7.1990	.23436	.00012
SDev	.00094	.00001	.00044	.128	.0520	.00133	.00030
%RSD	4.9638	2.5736	63.089	.70793	.72191	.56613	250.98
#1 #2	.01824 .01957	.00048	00102 00039	17.949 18.130	7.1623 7.2358	.23343	.00034 00009
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.02455	.00338	00022	.00098	00233	.00206	.00060
SDev	.00014	.00297	.00305	.00105	.00260	.00172	.00028
%RSD	.58112	87.967	1411.1	106.87	111.47	83.307	46.781
#1	.02445	.00548	00237	.00024	00417	.00328	.00080
#2	.02465	.00128	.00194	.00172	00049	.00085	
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00074	.00250	.00142	00247	.00039	.04233	
SDev	.00286	.00191	.00032	.00183	.00052	.00025	
%RSD	384.51	76.324	22.525	74.293	132.64	.58184	
#1 #2	.00128 00276	.00115	.00119 .00164	00376 00117	.00075 .00002	.04215 .04250	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Ana]	vsis	Report

05/17/00 12:51:46 PM

page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16501 26.51650 .1607002	2 NOTUSED    	3 NOTUSED    	4 NOTUSED	5 NOTUSED    	6 NOTUSED	7 NOTUSED    
#1	16482					<del></del>	
#2	16519				- <del>-</del>	<b>→</b> -	

Method: METTRA Sample Name: DCQGQF Run Time: 05/17/00 12:51:50 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD				
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm				
Avge	00127	2.3885	.00338	.01780	.00741	99.061	.00335				
SDev	.00019	.0826	.00123	.00042	.00006	2.513	.00011				
%RSD	15.259	3.4602	36.555	2.3305	.80259	2.5372	3.1319				
#1	00113	2.4469	.00250	.01810	.00737	100.84	.00328				
#2	00141	2.3300	.00425	.01751	.00746	97.284	.00342				
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass				
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000				
Low	01000	20000	01000	20000	00500	-5.0000	00500				
Elem	CO	CR	CU	FE	MG	MN	MO				
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm				
Avge	.20985	.00369	00087	37.930	34.426	.36830	00134				
SDev	.00559	.00019	.00004	.936	.812	.00854	.00017				
%RSD	2.6654	5.0454	4.6720	2.4670	2.3591	2.3193	12.997				
#1	.21380	.00356	00090	38.591	35.000	.37434	00122				
#2	.20589		00084	37.268	33.851	.36226	00146				
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass				
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000				
Low	05000	01000	02500	10000	-5.0000	01500	04000				
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB				
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm				
Avge	.28048	.00010	.00187	.00128	.00139	00047	.00015				
SDev	.00880	.00042	.00183	.00108	.00190	.00178	.00055				
%RSD	3.1356	434.50	97.920	84.453	137.04	374.88	378.08				
#1	.28670	00020	.00316	.00204	.00004	.00078	.00054				
#2	.27426	.00040	.00057	.00052	.00273	00173	00024				
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000				
Elem	SE/1	SE/2	SE	TL	V_	ZN					
Units	ppm	ppm	ppm	ppm	ppm	ppm					
Avge	00341	.00206	.00023	00040	.00153	.58684					
SDev	.00186	.00208	.00077	.00086	.00039	.01602					
%RSD	54.392	101.06	327.05	213.07	25.635	2.7305					
#1	00472	.00352	.00078	.00020	.00180	.59817					
#2	00210	.00059	00031	00101	.00125	.57551					
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000					

05/17/00 12:55:56 PM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 17543 420.9757 2.399717	2 NOTUSED   	3 NOTUSED    	4 NOTUSED    	5 NOTUSED   	6 NOTUSED    	7 NOTUSED    
#1 #2	17245 17840			<del>-</del> -	- <del>-</del>	- <b>-</b>	- <del>-</del>

05/17/00 01:00:06 PM

Method: METTRA Sample Name: DCQGVF Run Time: 05/17/00 12:56:00 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00035	.02025	.00743	.02114	00020	86.423	00012
SDev	.00011	.01042	.00060	.00001	.00003	.441	.00007
%RSD	30.873	51.458	8.1119	.03456	14.477	.50975	64.246
#1	00043	.02762	.00701	.02114	00022	86.735	00017
#2	00027	.01288	.00786	.02113	00018	86.112	00006
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00044	.00005	.00122	8.6577	17.622	.54538	.00038
SDev	.00020	.00027	.00047	.0331	.081	.00176	.00063
%RSD	45.344	568.98	38.226	.38287	.46204	.32294	165.36
#1	00030	00014	.00089	8.6812	17.680	.54663	.00082
#2	00058	.00024	.00155	8.6343	17.564	.54414	00006
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00038	.00147	.00061	.00090	00172	.00154	.00045
SDev	.00099	.00398	.00011	.00125	.00059	.00093	.00042
%RSD	262.11	269.88	17.929	138.83	34.431	60.539	93.409
#1	.00108	00134	.00069	.00002	00130	.00088	.00015
#2	00032	.00428	.00054	.00178	00214	.00220	.00075
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00777	.00272	00077	00191	.00226	.01624	
SDev	.00070	.00102	.00045	.00301	.00001	.00002	
%RSD	9.0635	37.465	57.760	157.10	.38286	.10050	
#1	00728	.00200	00109	.00021	.00227	.01625	
#2	00827	.00345	00046	00404	.00225	.01623	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

05/17/00 01:00:06 PM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED	
Elem	Y					<del>-</del> -		
Wavlen	371.030							
Avge	16263				<del></del>			
SDev	103.9447							
%RSD	.6391503					<b>-</b> -	<del></del>	
#1	16189		<del></del>					
#2	16336		<del>-</del> -					

Method: METTRA Sample Name: DCVMDF Run Time: 05/17/00 13:00:10 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00006	.08445	00067	.21981	00094	552.08	.00193
SDev	.00032	.01902	.00022	.00160	.00005	6.22	.00009
%RSD	546.43	22.518	32.510	.72861	5.3837	1.1258	4.8278
#1	.00017	.09790	00083	.22094	00097	556.48	.00186
#2	00028	.0 <b>71</b> 01	00052	.21868	00090	547.69	.00199
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01571	.00073	.00205	.58627	83.280	1.5385	00077
SDev	.00045	.00012	.00019	.00449	.904	.0150	.00035
%RSD	2.8615	16.558	9.3602	.76655	1.0860	.97374	45.076
#1	.01602	.00065	.00192	.58309	83.919	1.5491	00052
#2	.01539	.00082	.00219	.58945	82.640	1.5279	00101
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00978	00001	.00034	.00023	.00168	00117	00022
SDev	.00056	.00019	.00032	.00015	.00087	.00127	.00056
%RSD	5.7528	2861.6	93.271	66.642	51.587	108.20	251.02
#1	.01018	00014	.00057	.00033	.00107	00027	.00017
#2	.00938	.00013	.00012		.00229	00207	00062
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00006	00127	00083	.00208	.00244	.02854	
SDev	.00238	.00171	.00035	.00595	.00043	.00049	
%RSD	4149.8	134.74	42.257	285.50	17.416	1.7027	
#1	.00174	00249	00108	00212	.00275	.02888	
#2	00163	00006	00058	.00629	.00214	.02819	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Report 657		657 76	5 0	05/17/00 01:04:15 PM			page 2	
IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED	
Elem Wavlen	Y 371.030							
Avge	15095							
SDev	146.1239						<b></b>	
%RSD	.9680557							
#1	14991							
#2	15198			<del></del>				

Method: METTRA Sample Name: DCVMEF Run Time: 05/17/00 13:04:19 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem Units Avge SDev %RSD	AG ppm 00029 .00015 53.037	AL ppm .13145 .00773 5.8831	AS ppm .00055 .00012 21.107	BA ppm .10568 .00034 .32655	BE ppm 00014 .00014 103.63	CA ppm 166.05 .09	CD ppm 00013 .00023 175.94
#1	00040	.13692	.00064	.10543	00004	165.99	00029
#2	00018	.12598	.00047	.10592	00024	166.12	.00003
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00112	.00006	.00091	5.3524	4.8520	.06388	.00063
SDev	.00002	.00018	.00055	.0170	.0092	.00005	.00025
%RSD	1.4124	271.67	59.634	.31687	.18918	.07069	40.010
#1	00114	00006	.00053	5.3404	4.8455	.06385	.00045
#2	00111	.00019	.00130	5.3643	4.8585	.06392	.00081
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00019	.00157	.00145	.00149	.00152	00121	00030
SDev	.00073	.00019	.00068	.00052	.00208	.00034	.00046
%RSD	376.13	12.168	46.914	34.714	136.74	28.345	155.03
#1	.00071	.00170	.00193	.00185	.00005	00097	00063
#2	00032	.00143	.00097	.00112	.00299	00145	.00003
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00518	.00210	00032	00009	.00140	.01466	
SDev	.00028	.00211	.00131	.00136	.00000	.00009	
%RSD	5.4891	100.31	405.48	1525.7	.31687	.57673	
#1	00538	.00359	.00060	.00087	.00139	.01472	
#2	00498	.00061	00125	00105	.00140	.01461	
Errors High	NOCHECK	NOCHECK	LC Pass	LC Pass	LC Pass	LC Pass	

657 762

Analysis kepor	ysis Report
----------------	-------------

05/17/00 01:08:25 PM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16357 119.2179 .7288629	2 NOTUSED    	3 NOTUSED    	4 NOTUSED	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1	16441						
#2	16272	- <b>-</b>					

Method: METTRA Sample Name: DCVMFF Run Time: 05/17/00 13:08:29 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00027	.07545	00030	.08150	00121	598.77	.00024
SDev	.00030	.00560	.00268	.00005	.00002	.38	.00010
%RSD	110.98	7.4218	895.51	.05678	1.3022	.06282	42.812
#1	.00048	.07941	.00160	.08153	00122	599.04	.00032
#2	.00006	.07149	00219	.08146	00119	598.50	.00017
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem Units Avge SDev %RSD	CO ppm .00957 .00015 1.6059	CR ppm .00134 .00004 3.2761	CU ppm .00199 .00030 15.092	FE ppm .14180 .00155 1.0922	MG ppm 306.13 .15	MN ppm 3.6181 .0027 .07381	MO ppm 00048 .00000 .69497
#1	.00968	.00131	.00178	.14290	306.03	3.6162	00048
#2	.00946	.00137	.00221	.14071	306.24	3.6200	00047
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00846	00121	.00096	.00024	.00058	00045	00011
SDev	.00081	.00161	.00060	.00013	.00206	.00140	.00025
%RSD	9.5239	133.13	63.006	56.686	355.43	313.25	236.42
#1	.00903	00234	.00138	.00014	00088	.00054	.00007
#2	.00789	00007	.00053		.00204	00144	00028
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00291	00104	00166	00132	.00886	.01380	
SDev	.00322	.00062	.00149	.00246	.00043	.00004	
%RSD	110.63	60.227	89.656	186.51	4.8357	.27928	
#1	00063	00060	00061	.00042	.00856	.01383	
#2	00519	00148	00271	00306	.00917	.01377	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Repor	s Report
----------------	----------

05/17/00 01:12:35 PM

page 2

Mode	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	371.030				<b></b>		
Wavlen							
Avge SDev	14683			~-			
	21.53085			<b>→</b>	<del></del>		
%RSD	.1466422						
#1	14698					~ ~	
#2	14667				<del></del>		

05/17/00 01:16:45 PM

page 1

Method: METTRA Sample Name: CCV3-6 Run Time: 05/17/00 13:12:39 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.0217	24.493	.50774	1.9947	1.9549	48.689	.47818
SDev	.0031	.103	.00492	.0039	.0087	.259	.00181
%RSD	.30487	.42009	.96969	.19500	.44567	.53255	.37866
#1	1.0239	24.566	.51122	1.9975	1.9611	48.873	.47946
#2	1.0195	24.420	.50426	1.9920	1.9488	48.506	.47690
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	27.500	.55000	2.2000	2.2000	55.000	.55000
Low	.90000	22.500	.45000	1.8000	1.8000	45.000	.45000
Elem Units Avge SDev %RSD	CO ppm 1.9645 .0112 .56924	CR ppm 1.9732 .0081 .41096	CU ppm 1.9876 .0107 .53844	FE ppm 24.429 .075 .30721	MG ppm 48.165 .186 .38594	MN ppm 1.9449 .0060	MO ppm 1.9825 .0031 .15730
#1	1.9724	1.9790	1.9952	24.482	48.296	1.9491	1.9803
#2	1.9566	1.9675	1.9800	24.376	48.033	1.9406	1.9847
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.2000	2.2000	2.2000	27.500	55.000	2.2000	2.2000
Low	1.8000	1.8000	1.8000	22.500	45.000	1.8000	1.8000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.9854	.48421	.49073	.48856	.49983	.50753	.50497
SDev	.0159	.00048	.00321	.00230	.00270	.00032	.00111
%RSD	.80141	.09897	.65453	.47118	.54095	.06219	.22000
#1	1.9966	.48455	.49300	.49019	.50174	.50776	.50575
#2	1.9741	.48387	.48846	.48693	.49792	.50731	.50418
Errors High Low	LC Pass 2.2000 1.8000	NOCHECK	NOCHECK	LC Pass .55000 .45000	NOCHECK	NOCHECK	LC Pass .55000 .45000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.51990	.51704	.51799	.99821	1.9511	2.0208	
SDev	.00170	.00201	.00191	.01063	.0110	.0156	
%RSD	.32733	.38915	.36848	1.0649	.56381	.77044	
#1	.52110	.51847	.51934	1.0057	1.9588	2.0318	
#2	.51870	.51562	.51664	.99069	1.9433	2.0098	
Errors High Low	NOCHECK	NOCHECK	LC Pass .55000 .45000	LC Pass 1.1000 .90000	LC Pass 2.2000 1.8000	LC Pass 2.2000 1.8000	

657 766

Analysis Report

05/17/00 01:16:45 PM

page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16821 22.73376 .1351492	2 NOTUSED	3 NOTUSED    	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1	16805						
#2	16837		- **		<b>~</b> →		

Method: METTRA Sample Name: CCB6 Operator: RJG

Run Time: 05/17/00 13:16:49

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG ppm00003 .00006 186.29	AL	AS	BA	BE	CA	CD
Units		ppm	ppm	ppm	ppm	ppm	ppm
Avge		.00085	.00003	.00045	.00067	.04271	.00002
SDev		.00980	.00081	.00006	.00014	.01214	.00015
%RSD		1149.0	2379.1	14.213	21.343	28.421	661.44
#1	00007	.00778	.00061	.00050	.00077	.03413	.00013
#2	.00001	00608	00054	.00041	.00057	.05129	00008
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.01000	.20000	.00500	5.0000	.00500
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00040	.00021	.00008	.00014	.01946	.00057	.00558
SDev	.00027	.00002	.00023	.00942	.00414	.00001	.00246
%RSD	67.243	10.948	289.83	6838.8	21.275	2.1190	44.050
#1	00021	.00022	00008	.00680	.01654	.00058	.00731
#2	00058	.00019	.00024	00653	.02239	.00057	.00384
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.05000	.01000	.02500	.10000	5.0000	.01500	.04000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00097	00041	.00053	.00022	.00010	.00084	.00059
SDev	.00006	.00008	.00131	.00090	.00048	.00141	.00110
%RSD	5.9022	19.019	245.57	407.25	500.03	167.33	185.32
#1	.00101	00046	00039	00042	.00044	.00184	.00137
#2	.00093	00035	.00146	.00086	00024	00015	00018
Errors High Low	LC Pass .04000 04000	NOCHECK	NOCHECK	LC Pass .00300 00300	NOCHECK	NOCHECK	LC Pass .06000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00124	00054	.00006	.00308	.00063	.00085	
SDev	.00335	.00295	.00308	.00224	.00018	.00008	
%RSD	269.27	547.15	5575.6	72.831	28.532	9.2727	
#1	.00361	.00155	.00223	.00149	.00076	.00090	
#2	00113	00262	00212	.00467	.00051	.00079	
Errors High Low	NOCHECK	NOCHECK	LC Pass .00500 00500	LC Pass .01000 01000	LC Pass .05000 05000	LC Pass .02000 02000	

STL Pittsburgh

05/17/00 01:20:54 PM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 17018 24.43109 .1435643	2 NOTUSED   	3 NOTUSED   	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1	17035						
#2	17000						

page 1

Method: METTRA Sample Name: DCVMGF Run Time: 05/17/00 13:20:58 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00018	.02272	00043	.07373	00026	161.27	00004
SDev	.00022	.00810	.00195	.00039	.00006	.06	.00004
%RSD	127.53	35.642	457.66	.52951	24.504	.03875	93.234
#1	.00002	.02844	.00095	.07345	00022	161.22	00001
#2	.00033	.01699	00180	.07400	00031	161.31	00006
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00057	.00442	.00180	1.7115	5.8384	.08000	.00815
SDev	.00028	.00004	.00003	.0069	.0096	.00027	.00053
%RSD	49.278	.88928	1.4210	.40362	.16362	.33300	6.5278
#1	00077	.00444	.00179	1.7067	5.8317	.07981	.00777
#2	00037		.00182	1.7164	5.8452	.08019	.00852
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00240	.01001	.00838	.00892	00005	.00149	.00098
SDev	.00037	.00035	.00011	.00019	.00123	.00081	.00013
%RSD	15.563	3.4531	1.3623	2.1430	2373.5	54.639	13.737
#1	.00267	.00976	.00830	.00879	.00082	.00091	.00088
#2	.00214	.01025	.00847	.00906	00092	.00206	.00107
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00633	.00479	.00108	.00128	.00085	.01461	
SDev	.00065	.00192	.00107	.00199	.00019	.00009	
%RSD	10.278	40.178	98.304	155.97	22.127	.61635	
#1 #2	00679 00587	.00614 .00343	.00184	.00269 00013	.00098 .00072	.01467 .01454	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

657 770

Analysis Report

05/17/00 01:25:04 PM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	Y						
Wavlen	371.030	<b></b>					
Avge	16130					- <del>-</del>	
SDev	6.328744	₩ ₩			<b>~</b> ~		
%RSD	.0392352						
#1	16126				w 4		
#2	16135					<del></del>	

Operator: RJG

page 1

Method: METTRA Sample Name: DCVMHF

Run Time: 05/17/00 13:25:08

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

				•			
Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00038	.04047	.00068	.06850	00061	H630.64	00018
SDev	.00025	.00710	.00181	.00023	.00014	.68	.00003
%RSD	64.568	17.550	264.72	.33018	23.494	.10861	18.981
#1	00056	.04550	00060	.06834	00050	H631.13	00020
#2	00021	.03545	.00196	.06866	00071	H630.16	00015
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00293	.00022	.00041	.06685	41.856	1.0109	.00062
SDev	.00025	.00009	.00087	.00085	.007	.0008	.00030
%RSD	8.4383	41.401	210.51	1.2790	.01594	.07531	49.078
#1	.00276	.00016	00020	.06746	41.851	1.0104	.00083
#2	.00311	.00029	.00103	.06625	41.860	1.0114	
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00381	00114	.00003	00036	.00133	.00194	.00174
SDev	.00072	.00139	.00196	.00177	.00014	.00039	.00030
%RSD	18.978	122.10	6406.2	494.78	10.364	19.979	17.525
#1 #2	.00330	00212 00016	00136 .00142	00161 .00089	.00143	.00222 .00167	.00196 .00153
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00361	.00077	00069	00082	.00168	.00893	
SDev	.00215	.00272	.00253	.00046	.00001	.00024	
%RSD	59.667	352.80	367.52	55.985	.39383	2.6675	
#1	00514	00115	00248	00114	.00167	.00876	
#2	00209	.00269	.00110	00050	.00168	.00909	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

657 772

		-	
Anal	vala	Ren	へかた
LITICIT	уртр	T.C.D.	

05/17/00 01:29:14 PM

page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 15623 64.27628 .4114222	2 NOTUSED   	3 NOTUSED    	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1 #2	15668 15578						

Method: METTRA Sample Name: DCVMHP5F Operator: RJG

Run Time: 05/17/00 13:29:18

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00035	.00842	00058	.01344	00025	128.56	00007
SDev	.00030	.01163	.00151	.00016	.00008	2.75	.00025
%RSD	86.556	138.15	259.70	1.1802	31.405	2.1407	364.18
#1	00056	.01664	00164	.01355	00030	130.50	00025
#2	00014	.00019	.00048	.01333	00019	126.61	.00011
Elem	CO	CR	CU	FE ppm00175 .00459 262.44	MG	MN	MO
Units	ppm	ppm	ppm		ppm	ppm	ppm
Avge	.00004	00020	00019		8.2087	.19772	.00006
SDev	.00047	.00026	.00039		.1678	.00405	.00026
%RSD	1112.2	129.21	211.35		2.0442	2.0500	428.53
#1	00029	00039	00046	.00150	8.3273	.20059	.00024
#2	.00038	00002	.00009	00499	8.0900	.19485	00012
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00105	00202	.00084	00011	.00201	.00105	.00137
SDev	.00002	.00069	.00045	.00053	.00287	.00038	.00121
%RSD	1.6829	34.139	53.373	476.36	142.93	36.524	88.495
#1	.00106	00251	.00053	00049	.00404	.00132	.00223
#2	.00104	00154	.00116	.00026	00002	.00078	.00051
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00046	00101	00083	00022	.00065	.00206	
SDev	.00012	.00085	.00060	.00337	.00017	.00008	
%RSD	25.740	83.821	73.020	1506.6	26.866	3.8477	
#1 #2	00038 00055	00041 00161	00040 00125	.00216 00260	.00053 .00077	.00211	

05/17/00 01:33:23 PM page 2

IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED 	6 NOTUSED	7 NOTUSED
Wavlen	371.030				- <del>-</del>		
Avge	16553						
SDev	215.6676					~ ~	
%RSD	1.302879						
#1	16401					<b>-</b> -	
#2	16706						

Method: METTRA Sample Name: DCVMHSF Operator: RJG

Run Time: 05/17/00 13:33:27

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.05448	2.3702	2.1854	2.1546	.04958	H682.93	.04793
SDev	.00018	.0008	.0123	.0016	.00030	2.92	.00040
%RSD	.32981	.03582	.56231	.07261	.59964	.42782	.82534
#1	.05435	2.3708	2. <b>1</b> 941	2.1557	.04979	H685.00	.04821
#2	.05461	2.3696	2.1767	2.1535	.04937	H680.87	.04765
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.51271	.20534	.27171	1.0594	92.793	1.5364	1.0424
SDev	.00120	.00101	.00066	.0038	.254	.0041	.0006
%RSD	.23340	.49168	.24219	.35634	.27402	.26719	.05878
#1	.51355	.20606	.27217	1.0620	92.973	1.5393	1.0428
#2	.51186	.20463	.27124	1.0567	92.613	1.5335	1.0420
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.50773	.51067	.50018	.50367	.51919	.53794	.53169
SDev	.00383	.00046	.00212	.00157	.00114	.00334	.00261
%RSD	.75401	.09085	.42451	.31186	.22048	.62085	.49067
#1	.51043	.51100	.50168	.50478	.52000	.54030	.53354
#2	.50502	.51034	.49868	.50256	.51838	.53557	.52985
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	2.2199	2.1720	2.1879	2.2130	.52005	.55649	
SDev	.0110	.0031	.0057	.0008	.00232	.00166	
%RSD	.49745	.14151	.26177	.03520	.44551	.29734	
#1.	2.2277	2.1742	2.1920	2.2135	.52169	.55766	
#2	2.2121	2.1698	2.1839	2.2124	.51841	.55532	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

05/17/00 01:37:33 PM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	Y			<b>-</b> -			
Wavlen	371.030	<del></del>					~-
Avge	15308						
SDev	91.71189						
%RSD	.5991069	<del></del>					~ ~
#1	15243						~ ~
#2	15373						

Method: METTRA Sample Name: DCVMHDF Run Time: 05/17/00 13:37:37 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	an.
Units Avge	ppm .05250	ppm 2.2651	ppm	ppm	ppm	ppm	CD ppm
SDev	.00040	.0132	2.1046 .0185	2.0759 .0056	.04787 .00035	H660.76 3.08	.04599 .00005
%RSD	.76543	.58230	.87721	.27082	.72318	.46639	.11666
#1 #2	.05222	2.2745	2.1177	2.0799	.04812	H662.93	.04595
#4	.05278	2.2558	2.0916	2.0719	.04763	H658.58	.04602
Errors High	LC Pass 2.0000	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
Low	01000	20000	10.000 01000	10.000 20000	10.000 00500	600.00 -5.0000	5.0000 00500
Elem	CO	CR	CU	FE	MG		
Units	ppm	ppm	ppm	ppm	ppm	MN ppm	MO mqq
Avge SDev	.49378 .00194	.19830 .00003	.26184 .00038	1.0092 .0057	89.557 .317	1.4845	1.0089
%RSD	.39283	.01464	.14490	.56844	.35434	.0049 .33276	.0015 .14602
#1	.49515	.19832	.26211	1.0051	89.781	1.4880	1.0100
#2	.49240	.19828	.26157	1.0133	89.332	1.4810	1.0079
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High Low	100.00 05000	20.000 01000	10.000 02500	500.00 10000	600.00 -5.0000	10.000 01500	20.000 04000
					3.0000	.01500	04000
Elem	NT	ו/ מם	DD / 2	DD.	an /-	/-	
Elem Units	NI ppm	PB/1 ppm	PB/2 ppm	PB Rqq	SB/1	SB/2	SB
Units Avge	ppm .48815	ppm .49174	ppm .48323	ppm .48606	ppm .50533	ppm .51927	ppm .51463
Units	ppm	ppm	ppm	ppm	ppm	ppm .51927 .00078	ppm .51463 .00010
Units Avge SDev %RSD	ppm .48815 .00267 .54796	ppm .49174 .00158 .32033	ppm .48323 .00061 .12591	ppm .48606 .00012 .02443	ppm .50533 .00127 .25122	ppm .51927 .00078 .15074	ppm .51463 .00010 .01931
Units Avge SDev	ppm .48815 .00267	ppm .49174 .00158	ppm .48323 .00061	ppm .48606 .00012	ppm .50533 .00127	ppm .51927 .00078	ppm .51463 .00010
Units Avge SDev %RSD #1 #2 Errors	ppm .48815 .00267 .54796 .49004 .48626	ppm .49174 .00158 .32033	ppm .48323 .00061 .12591	ppm .48606 .00012 .02443	ppm .50533 .00127 .25122	ppm .51927 .00078 .15074	ppm .51463 .00010 .01931 .51456 .51470
Units Avge SDev %RSD #1 #2 Errors High	ppm .48815 .00267 .54796 .49004 .48626 LC Pass 100.00	ppm .49174 .00158 .32033 .49285 .49062	ppm .48323 .00061 .12591 .48280 .48366	ppm .48606 .00012 .02443 .48614 .48598 LC Pass 5.0000	ppm .50533 .00127 .25122 .50623 .50443	ppm .51927 .00078 .15074 .51872 .51983	ppm .51463 .00010 .01931 .51456 .51470 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low	ppm .48815 .00267 .54796 .49004 .48626 LC Pass 100.00 04000	ppm .49174 .00158 .32033 .49285 .49062 NOCHECK	ppm .48323 .00061 .12591 .48280 .48366 NOCHECK	ppm .48606 .00012 .02443 .48614 .48598 LC Pass 5.0000 00300	ppm .50533 .00127 .25122 .50623 .50443 NOCHECK	ppm .51927 .00078 .15074 .51872 .51983	ppm .51463 .00010 .01931 .51456 .51470
Units Avge SDev %RSD #1 #2 Errors High	ppm .48815 .00267 .54796 .49004 .48626 LC Pass 100.00 04000 SE/1	ppm .49174 .00158 .32033 .49285 .49062 NOCHECK	ppm .48323 .00061 .12591 .48280 .48366 NOCHECK	ppm .48606 .00012 .02443 .48614 .48598 LC Pass 5.0000 00300	ppm .50533 .00127 .25122 .50623 .50443 NOCHECK	ppm .51927 .00078 .15074 .51872 .51983 NOCHECK	ppm .51463 .00010 .01931 .51456 .51470 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge	ppm .48815 .00267 .54796 .49004 .48626 LC Pass 100.00 04000 SE/1 ppm 2.1417	ppm .49174 .00158 .32033 .49285 .49062 NOCHECK SE/2 ppm 2.1075	ppm .48323 .00061 .12591 .48280 .48366 NOCHECK	PPM .48606 .00012 .02443 .48614 .48598 LC Pass 5.000000300 TL ppm 2.1320	ppm .50533 .00127 .25122 .50623 .50443 NOCHECK	ppm .51927 .00078 .15074 .51872 .51983	ppm .51463 .00010 .01931 .51456 .51470 LC Pass 10.000
Units Avge SDev RSD #1 #2 Errors High Low Elem Units	ppm .48815 .00267 .54796 .49004 .48626 LC Pass 100.00 04000 SE/1 ppm	ppm .49174 .00158 .32033 .49285 .49062 NOCHECK	ppm .48323 .00061 .12591 .48280 .48366 NOCHECK SE ppm 2.1189 .0066	PPM .48606 .00012 .02443 .48614 .48598 LC Pass 5.000000300 TL ppm 2.1320 .0161	ppm .50533 .00127 .25122 .50623 .50443 NOCHECK V_ppm .50222 .00357	ppm .51927 .00078 .15074 .51872 .51983 NOCHECK ZN ppm .53412 .00367	ppm .51463 .00010 .01931 .51456 .51470 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge SDev %RSD	ppm .48815 .00267 .54796 .49004 .48626 LC Pass 100.00 04000 SE/1 ppm 2.1417 .0018 .08555	Ppm .49174 .00158 .32033 .49285 .49062 NOCHECK SE/2 Ppm 2.1075 .0089 .42322	PPM .48323 .00061 .12591 .48280 .48366 NOCHECK  SE PPM 2.1189 .0066 .30957	PPM .48606 .00012 .02443 .48614 .48598 LC Pass 5.000000300 TL PPM 2.1320 .0161 .75597	ppm .50533 .00127 .25122 .50623 .50443 NOCHECK V_ppm .50222 .00357 .71100	DPM .51927 .00078 .15074 .51872 .51983 NOCHECK ZN PPM .53412 .00367 .68703	ppm .51463 .00010 .01931 .51456 .51470 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge SDev	ppm .48815 .00267 .54796 .49004 .48626 LC Pass 100.00 04000 SE/1 ppm 2.1417 .0018	ppm .49174 .00158 .32033 .49285 .49062 NOCHECK SE/2 ppm 2.1075 .0089	ppm .48323 .00061 .12591 .48280 .48366 NOCHECK SE ppm 2.1189 .0066	PPM .48606 .00012 .02443 .48614 .48598 LC Pass 5.000000300 TL ppm 2.1320 .0161	ppm .50533 .00127 .25122 .50623 .50443 NOCHECK V_ppm .50222 .00357	ppm .51927 .00078 .15074 .51872 .51983 NOCHECK ZN ppm .53412 .00367	ppm .51463 .00010 .01931 .51456 .51470 LC Pass 10.000
Units Avge SDev RSD #1 #2 Errors High Low Elem Units Avge SDev RSD #1 #2 Errors	ppm .48815 .00267 .54796 .49004 .48626 LC Pass 100.00 04000 SE/1 ppm 2.1417 .0018 .08555	ppm .49174 .00158 .32033 .49285 .49062 NOCHECK SE/2 ppm 2.1075 .0089 .42322 2.1138	PPM .48323 .00061 .12591 .48280 .48366 NOCHECK  SE PPM 2.1189 .0066 .30957 2.1235	ppm .48606 .00012 .02443 .48614 .48598 LC Pass 5.0000 00300 TL ppm 2.1320 .0161 .75597 2.1433	ppm .50533 .00127 .25122 .50623 .50443 NOCHECK V_ppm .50222 .00357 .71100 .50474 .49969	DPM .51927 .00078 .15074 .51872 .51983 NOCHECK ZN PPM .53412 .00367 .68703 .53671 .53152	ppm .51463 .00010 .01931 .51456 .51470 LC Pass 10.000
Units Avge SDev RSD #1 #2 Errors High Low Elem Units Avge SDev RSD #1 #2	ppm .48815 .00267 .54796 .49004 .48626 LC Pass 100.00 04000 SE/1 ppm 2.1417 .0018 .08555 2.1430 2.1404	Ppm .49174 .00158 .32033 .49285 .49062 NOCHECK SE/2 ppm 2.1075 .0089 .42322 2.1138 2.1012	PPM .48323 .00061 .12591 .48280 .48366 NOCHECK  SE PPM 2.1189 .0066 .30957 2.1235 2.1142	PPM .48606 .00012 .02443 .48614 .48598 LC Pass 5.000000300 TL ppm 2.1320 .0161 .75597 2.1433 2.1206	ppm .50533 .00127 .25122 .50623 .50443 NOCHECK V_ppm .50222 .00357 .71100 .50474	Ppm .51927 .00078 .15074 .51872 .51983 NOCHECK ZN ppm .53412 .00367 .68703	ppm .51463 .00010 .01931 .51456 .51470 LC Pass 10.000

05/17/00 01:41:43 PM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 15701 32.63325 .2078442	2 NOTUSED   	3 NOTUSED    	4 NOTUSED    	5 NOTUSED    	6 NOTUSED    	7 NOTUSED   
#1 #2	15678 15724		- <del>-</del>		• -	~~	

Method: METTRA Sample Name: DCVMKF Run Time: 05/17/00 13:41:47 Operator: RJG

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00038	.04836	.00067	.08458	00049	181.62	00021
SDev	.00006	.01891	.00153	.00022	.00001	.34	.00002
%RSD	15.791	39.104	229.42	.25526	1.4261	.18538	11.481
#1	00042	.06173	00042	.08443	00050	181.86	00019
#2	00033	.03499	.00175	.08473	00049	181.38	00022
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00102	.00002	.00087	6.5926	5.1506	.05611	.00458
SDev	.00045	.00059	.00019	.0106	.0095	.00008	.00119
%RSD	44.424	2486.8	21.636	.16027	.18451	.14098	25.941
#1 #2	00070 00134	.00044 00039	.00100	6.5851 6.6000	5.1573 5.1439	.05617 .05606	.00542
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00131	.00038	.00143	.00108	.00031	00017	00001
SDev	.00181	.00022	.00160	.00100	.00285	.00142	.00001
%RSD	138.44	57.854	112.16	92.198	931.96	821.99	44.162
#1	.00259	.00054	.00030	.00038	00171	.00083	00002
#2	.00003		.00256	.00178	.00232	00117	00001
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00299	.00274	.00083	.00042	.00051	.03099	
SDev	.00341	.00075	.00063	.00167	.00020	.00005	
%RSD	114.05	27.360	75.884	397.81	38.685	.15348	
#1	00058	.00221	.00128	.00160	.00037	.03103	
#2	00539	.00327	.00039	00076	.00065	.03096	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

Analysis Repor	Ana	lvsi	ls F	epc	ort
----------------	-----	------	------	-----	-----

05/17/00 01:45:53 PM

page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16116 85.77191 .5322076	2 NOTUSED    	3 NOTUSED    	4 NOTUSED	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    
#1 #2	16056 16177						

Operator: RJG

Method: METTRA Sample Name: DCVMLF

Run Time: 05/17/00 13:45:57

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG ppm00002 .00056 2792.9	AL	AS	BA	BE	CA	CD
Units		ppm	ppm	ppm	ppm	ppm	ppm
Avge		.04216	.00035	.16156	00099	418.68	.00007
SDev		.00186	.00033	.00039	.00005	.80	.00007
%RSD		4.4183	96.020	.24371	5.4415	.19180	100.98
#1	.00037	.04084	.00058	.16184	00103	419.25	.00002
#2	00041	.04347	.00011	.16128	00095	418.12	.00011
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00104	.03484	.00296	1.8539	40.448	1.1390	.00981
SDev	.00018	.00004	.00033	.0021	.057	.0015	.00015
%RSD	17.600	.11495	11.089	.11350	.14099	.13528	1.5738
#1	.00091	.03482	.00272	1.8554	40.488	1.1400	.00970
#2	.00117	.03487	.00319	1.8524	40.408	1.1379	.00992
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
m1	177		/-				
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00449	.00141	.00177	.00165	00011	.00038	.00022
SDev	.00052	.00032	.00107	.00082	.00041	.00021	.00028
%RSD	11.470	22.788	60.624	49.840	373.32	56.206	128.62
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00449	.00141	.00177	.00165	00011	.00038	.00022
SDev	.00052	.00032	.00107	.00082	.00041	.00021	.00028
Units Avge SDev %RSD #1	ppm .00449 .00052 11.470	ppm .00141 .00032 22.788	ppm .00177 .00107 60.624	ppm .00165 .00082 49.840	ppm 00011 .00041 373.32	ppm .00038 .00021 56.206	ppm .00022 .00028 128.62
Units Avge SDev %RSD #1 #2 Errors High	ppm .00449 .00052 11.470 .00486 .00413 LC Pass 100.00	ppm .00141 .00032 22.788 .00119 .00164	ppm .00177 .00107 60.624 .00101 .00253	ppm .00165 .00082 49.840 .00107 .00224 LC Pass 5.0000	ppm 00011 .00041 373.32 .00018 00040	ppm .00038 .00021 56.206 .00053 .00023	ppm .00022 .00028 128.62 .00042 .00002 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge SDev	ppm .00449 .00052 11.470 .00486 .00413 LC Pass 100.00 04000 SE/1 ppm 00102 .00445	ppm .00141 .00032 22.788 .00119 .00164 NOCHECK SE/2 ppm .00015 .00201	ppm .00177 .00107 60.624 .00101 .00253 NOCHECK SE ppm 00024 .00282	ppm .00165 .00082 49.840 .00107 .00224 LC Pass 5.0000 00300 TL ppm 00016 .00042	ppm00011 .00041 373.32 .0001800040 NOCHECK  V_ ppm .01004 .00059	ppm .00038 .00021 56.206 .00053 .00023 NOCHECK ZN ppm .01112 .00006	ppm .00022 .00028 128.62 .00042 .00002 LC Pass 10.000

page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 15599 5.056090 .0324139	2 NOTUSED	3 NOTUSED    	4 NOTUSED	5 NOTUSED    	6 NOTUSED    	7 NOTUSED    	
#1	15595					<b></b>		
#2	15602							

Operator: RJG

Method: METTRA Sample Name: DCVMMF

Run Time: 05/17/00 13:50:06

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

			-				
Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00042	.08169	.00055	.06866	00024	114.40	00019
SDev	.00044	.00747	.00004	.00006	.00002	.42	.00001
%RSD	103.93	9.1376	6.7598	.09288	7.0284	.36686	6.2499
#1	00011	.08697	.00053	.06861	00023	114.70	00019
#2	00074	.07642	.00058	.06870	00025	114.10	00018
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00096	00016	.00130	2.4450	4.6361	.03725	.00236
SDev	.00012	.00053	.00021	.0006	.0092	.00008	.00114
%RSD	12.221	332.51	16.451	.02251	.19808	.22656	48.287
#1	00088	.00022	.00115	2.4446	4.6426	.03731	.00316
#2	00105	00054	.00145	2.4454	4.6296	.03719	.00155
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00023	00062	.00209	.00119	.00198	.00036	.00090
SDev	.00067	.00397	.00087	.00190	.00096	.00065	.00012
%RSD	285.96	640.99	41.682	160.12	48.454	180.49	12.984
#1	.00070	.00219	.00271	.00253	.00130	.00083	.00099
#2	00024	00342	.00147	00016	.00266	00010	
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00637	.00161	00105	00046	.00129	.00982	
SDev	.00324	.00146	.00010	.00228	.00055	.00012	
%RSD	50.818	90.877	9.9188	490.51	42.983	1.2177	
#1	00408	.00057	00098	.00115	.00168	.00974	
#2	00865	.00264	00112	00208	.00090	.00991	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

05/17/00 01:54:12 PM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUS <b>E</b> D	7 NOTUSED
Elem	Y						
Wavlen	371.030	<b></b>					
Avge	16636			<del>-</del> -			
SDev	81.14050					<b></b>	
%RSD	.4877309						
#1	16579						
#2	16694						<b></b>

Operator: RJG

page 1

Method: METTRA Sample Name: DCVMNF Run Time: 05/17/00 13:54:16

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

					`		
Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00004	.03072	.00074	.07231	00082	338.08	00003
SDev	.00001	.00575	.00001	.00010	.00004	1.30	.00003
%RSD	23.801	18.729	1.7261	.14113	4.9990	.38506	101.94
#1 #2	.00004	.03479 .02665	.00075	.07238 .07223	00085 00080	339.00 337.15	00001 00006
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	10.000	10.000	10.000	600.00	5.0000
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00094	00014	.00225	1.3643	45.742	.84049	.00085
SDev	.00006	.00020	.00033	.0015	.138	.00228	.00040
%RSD	5.8287	144.72	14.660	.11122	.30227	.27090	47.325
#1	.00098	00028	.00249	1.3653	45.840	.84210	.00114
#2	.00090	.00000	.00202	1.3632	45.644	.83888	.00057
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	10.000	500.00	600.00	10.000	20.000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units.	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00125	.00041	.00023	.00029	.00019	.00097	.00071
SDev	.00006	.00059	.00151	.00081	.00139	.00234	.00110
%RSD	4.6259	142.75	650.41	277.58	713.95	242.09	154.73
#1	.00121	.00083	00084	00028	00079	.00262	.00148
#2	.00129	00000	.00130	.00087	.00118	00069	00007
Errors High Low	LC Pass 100.00 04000	NOCHECK	NOCHECK	LC Pass 5.0000 00300	NOCHECK	NOCHECK	LC Pass 10.000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	00227	.00276	.00109	.00016	.00391	.01124	
SDev	.00317	.00069	.00151	.00142	.00001	.00013	
%RSD	139.72	24.865	139.25	905.73	.33986	1.1828	
#1 #2	00451 00003	.00227 .00324	.00002 .00216	.00116 00085	.00392 .00390	.01114	
Errors High Low	NOCHECK	NOCHECK	LC Pass 10.000 00500	LC Pass 10.000 01000	LC Pass 50.000 05000	LC Pass 5.0000 02000	

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 15817 57.38130 .3627842	2 NOTUSED	3 NOTUSED	4 NOTUSED   	5 NOTUSED   	6 NOTUSED   	7 NOTUSED    
#1	15776						
#2	15858				- <del>-</del>		

Method: METTRA Sample Name: DCVMPF Operator: RJG

Run Time: 05/17/00 13:58:26

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	BA	D.D.	G.P.	<b></b>
Units	ppm	ppm	ppm	ppm	BE ppm	CA ppm	CD ppm
Avge SDev	.00004 .00030	.00493 .00682	.00124 .00081	.07901 .00008	00049 .00006	98.562	.00001
%RSD	686.55	138.20	65.450	.09610	12.920	.139 .14081	.00008 872.04
#1 #2	00017 .00026	.00976 .00011	.00066	.07906	00054	98.660	.00007
πΔ	.00026		.00181	.07896	00045	98.464	00005
Errors High	LC Pass 2.0000	LC Pass 600.00	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
Low	01000	20000	10.000 01000	10.000 20000	10.000 00500	600.00 ~5.0000	5.0000 00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge SDev	00127 .00048	00021 .00019	.00203 .00048	1.6326 .0072	3.7751 .0045	.05676 .00009	.00407 .00026
%RSD	37.994	89.995	23.766	.44297	.11834	.15279	6.2763
#1	00162	00008	.00169	1.6377	3.7782	.05670	.00426
#2	00093	00034	.00237	1.6275	3.7719	.05683	.00389
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High Low	100.00 05000	20.000 01000	10.000 02500	500.00 10000	600.00 -5.0000	10.000 01500	20.000
	.03000	.0100	.02500	10000	-5.0000	01500	04000
<b>777</b>	***	/-	n= /-				
Elem Units	IM	PB/1	PB/2	PB ppm	SB/1	SB/2	SB
Units Avge	ppm .00233	ppm 00028	ppm .00115	ppm .00068	ppm 00016	ppm 00057	SB ppm 00043
Units Avge SDev	ppm .00233 .00032	ppm 00028 .00049	ppm .00115 .00087	ppm .00068 .00074	ppm 00016 .00074	ppm 00057 .00075	ppm 00043 .00074
Units Avge SDev %RSD	ppm .00233 .00032 13.655	ppm 00028 .00049 177.00	ppm .00115 .00087 75.256	ppm .00068 .00074 109.86	ppm 00016 .00074 462.72	ppm 00057 .00075 131.43	ppm 00043 .00074 172.26
Units Avge SDev	ppm .00233 .00032	ppm 00028 .00049	ppm .00115 .00087	ppm .00068 .00074 109.86	ppm 00016 .00074 462.72	ppm 00057 .00075 131.43	ppm 00043 .00074 172.26
Units Avge SDev %RSD #1 #2	ppm .00233 .00032 13.655 .00256	ppm 00028 .00049 177.00 .00007 00063	ppm .00115 .00087 75.256 .00177 .00054	ppm .00068 .00074 109.86 .00120 .00015	ppm 00016 .00074 462.72 .00036 00068	ppm 00057 .00075 131.43 00004 00109	ppm 00043 .00074 172.26 .00009 00096
Units Avge SDev %RSD #1 #2 Errors High	ppm .00233 .00032 13.655 .00256 .00211 LC Pass 100.00	ppm 00028 .00049 177.00	ppm .00115 .00087 75.256	ppm .00068 .00074 109.86	ppm 00016 .00074 462.72	ppm 00057 .00075 131.43	ppm 00043 .00074 172.26 .00009 00096
Units Avge SDev %RSD #1 #2 Errors	ppm .00233 .00032 13.655 .00256 .00211	ppm 00028 .00049 177.00 .00007 00063	ppm .00115 .00087 75.256 .00177 .00054	ppm .00068 .00074 109.86 .00120 .00015	ppm 00016 .00074 462.72 .00036 00068	ppm 00057 .00075 131.43 00004 00109	ppm 00043 .00074 172.26 .00009 00096
Units Avge SDev %RSD #1 #2 Errors High Low Elem	ppm .00233 .00032 13.655 .00256 .00211 LC Pass 100.00 04000	ppm00028 .00049 177.00 .0000700063 NOCHECK	ppm .00115 .00087 75.256 .00177 .00054	ppm .00068 .00074 109.86 .00120 .00015 LC Pass 5.0000	ppm 00016 .00074 462.72 .00036 00068	ppm 00057 .00075 131.43 00004 00109	ppm 00043 .00074 172.26 .00009 00096 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low Elem Units	ppm .00233 .00032 13.655 .00256 .00211 LC Pass 100.00 04000 SE/1 ppm	ppm 00028 .00049 177.00 .00007 00063 NOCHECK	ppm .00115 .00087 75.256 .00177 .00054 NOCHECK	ppm .00068 .00074 109.86 .00120 .00015 LC Pass 5.0000 00300 TL ppm	ppm 00016 .00074 462.72 .00036 00068 NOCHECK	ppm 00057 .00075 131.43 00004 00109 NOCHECK	ppm 00043 .00074 172.26 .00009 00096 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low Elem	ppm .00233 .00032 13.655 .00256 .00211 LC Pass 100.00 04000	ppm00028 .00049 177.00 .0000700063 NOCHECK	ppm .00115 .00087 75.256 .00177 .00054 NOCHECK	ppm .00068 .00074 109.86 .00120 .00015 LC Pass 5.0000 00300	ppm00016 .00074 462.72 .0003600068 NOCHECK	ppm00057 .00075 131.430000400109 NOCHECK	ppm 00043 .00074 172.26 .00009 00096 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge	ppm .00233 .00032 13.655 .00256 .00211 LC Pass 100.00 04000 SE/1 ppm 00331	ppm 00028 .00049 177.00 .00007 00063 NOCHECK SE/2 ppm .00267	ppm .00115 .00087 75.256 .00177 .00054 NOCHECK SE ppm .00068	ppm .00068 .00074 109.86 .00120 .00015 LC Pass 5.0000 00300 TL ppm 00159	ppm 00016 .00074 462.72 .00036 00068 NOCHECK V_ppm .00109	ppm00057 .00075 131.430000400109 NOCHECK  ZN ppm .01048	ppm 00043 .00074 172.26 .00009 00096 LC Pass 10.000
Units Avge SDev RSD #1 #2 Errors High Low Elem Units Avge SDev RSD #1	ppm .00233 .00032 13.655 .00256 .00211 LC Pass 100.00 04000 SE/1 ppm 00331 .00023 6.8017	ppm00028 .00049 177.00 .0000700063  NOCHECK  SE/2 ppm .00267 .00177 66.219 .00142	ppm .00115 .00087 75.256 .00177 .00054 NOCHECK SE ppm .00068 .00111 162.27	ppm .00068 .00074 109.86 .00120 .00015 LC Pass 5.0000 00300 TL ppm 00159 .00065 41.153	ppm00016 .00074 462.72 .0003600068  NOCHECK  V_ ppm .00109 .00056 51.292 .00069	ppm00057 .00075 131.430000400109 NOCHECK  ZN ppm .01048 .00010 .92909 .01055	ppm 00043 .00074 172.26 .00009 00096 LC Pass 10.000
Units Avge SDev RSD #1 #2 Errors High Low Elem Units Avge SDev RSD	ppm .00233 .00032 13.655 .00256 .00211 LC Pass 100.00 04000 SE/1 ppm 00331 .00023 6.8017	ppm00028 .00049 177.00 .0000700063 NOCHECK  SE/2 ppm .00267 .00177 66.219	Dpm .00115 .00087 75.256 .00177 .00054 NOCHECK SE ppm .00068 .00111 162.27	ppm .00068 .00074 109.86 .00120 .00015 LC Pass 5.0000 00300 TL ppm 00159 .00065 41.153	Ppm00016 .00074 462.72 .0003600068 NOCHECK  V_ ppm .00109 .00056 51.292	ppm00057 .00075 131.430000400109 NOCHECK  ZN ppm .01048 .00010 .92909	ppm 00043 .00074 172.26 .00009 00096 LC Pass 10.000
Units Avge SDev RSD #1 #2 Errors High Low Elem Units Avge SDev RSD #1 #2 Errors	ppm .00233 .00032 13.655 .00256 .00211 LC Pass 100.00 04000 SE/1 ppm 00331 .00023 6.8017	ppm00028 .00049 177.00 .0000700063  NOCHECK  SE/2 ppm .00267 .00177 66.219 .00142	ppm .00115 .00087 75.256 .00177 .00054 NOCHECK SE ppm .00068 .00111 162.27 00010 .00146 LC Pass	ppm .00068 .00074 109.86 .00120 .00015 LC Pass 5.0000 00300 TL ppm 00159 .00065 41.153 00113 00205 LC Pass	Ppm00016 .00074 462.72 .0003600068 NOCHECK  V_ ppm .00109 .00056 51.292 .00069 .00148 LC Pass	ppm00057 .00075 131.430000400109  NOCHECK  ZN ppm .01048 .00010 .92909 .01055 .01041  LC Pass	ppm 00043 .00074 172.26 .00009 00096 LC Pass 10.000
Units Avge SDev %RSD #1 #2 Errors High Low Elem Units Avge SDev %RSD #1 #2	ppm .00233 .00032 13.655 .00256 .00211 LC Pass 100.00 04000 SE/1 ppm 00331 .00023 6.8017 00315 00347	ppm00028 .00049 177.00 .0000700063 NOCHECK  SE/2 ppm .00267 .00177 66.219 .00142 .00393	ppm .00115 .00087 75.256 .00177 .00054 NOCHECK SE ppm .00068 .00111 162.27 00010 .00146	ppm .00068 .00074 109.86 .00120 .00015 LC Pass 5.0000 00300 TL ppm 00159 .00065 41.153 00113 00205	ppm00016 .00074 462.72 .0003600068 NOCHECK  V_ppm .00109 .00056 51.292 .00069 .00148	ppm00057 .00075 131.430000400109  NOCHECK  ZN ppm .01048 .00010 .92909 .01055 .01041	ppm 00043 .00074 172.26 .00009 00096 LC Pass 10.000

Analysis Report		657 788		05/17/00 02:02:32 PM			page 2
IntStd Mode Elem	1 Counts Y	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Wavlen	371.030						<del>-</del> -

- -

--

--

- --

Avge SDev

SDev %RSD

#1

#2

16286 67.35192

16238

16333

.4135654

Operator: RJG

Method: METTRA Sample Name: CCV3-7 Run Time: 05/17/00 14:02:36

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.0394	25.100	.51538	2.0368	1.9885	49.581	.48838
SDev	.0011	.014	.00175	.0009	.0040	.073	.00097
%RSD	.10351	.05729	.33941	.04504	.20330	.14712	.19855
#1	1.0402	25.110	.51661	2.0374	1.9913	49.632	.48907
#2	1.0386	25.089	.51414	2.0361	1.9856	49.529	.48769
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	27.500	.55000	2.2000	2.2000	55.000	.55000
Low	.90000	22.500	.45000	1.8000	1.8000	45.000	.45000
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	2.0051	2.0123	2.0197	24.996	49.093	1.9851	2.0150
SDev	.0035	.0054	.0015	.031	.098	.0037	.0073
%RSD	.17547	.26855	.07653	.12402	.19923	.18484	.36295
#1	2.0076	2.0161	2.0207	25.018	49.162	1.9877	2.0098
#2	2.0026	2.0085	2.0186	24.974	49.024	1.9825	· 2.0202
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.2000	2.2000	2.2000	27.500	55.000	2.2000	2.2000
Low	1.8000	1.8000	1.8000	22.500	45.000	1.8000	1.8000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	2.0144	.49696	.50074	.49948	.50787	.52012	.51604
SDev	.0032	.00161	.00663	.00496	.00747	.00027	.00267
%RSD	.15636	.32472	1.3237	.99270	1.4712	.05215	.51719
#1	2.0166	.49810	.50543	.50299	.51316	.52031	.51793
#2	2.0122	.49582	.49606	.49598	.50259	.51993	.51416
Errors High Low	LC Pass 2.2000 1.8000	NOCHECK	NOCHECK	LC Pass .55000 .45000	NOCHECK	NOCHECK	LC Pass .55000 .45000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.53392	.52714	.52940	1.0150	1.9853	2.0566	
SDev	.00120	.00803	.00575	.0001	.0034	.0051	
%RSD	.22472	1.5231	1.0870	.00512	.17019	.24924	
#1	.53476	.53282	.5334 <b>7</b>	1.0149	1.9876	2.0603	
#2	.53307	.52146	.52533	1.0150	1.9829	2.0530	
Errors High Low	NOCHECK	NOCHECK	LC Pass .55000 .45000	LC Pass 1.1000 .90000	LC Pass 2.2000 1.8000	LC Pass 2.2000 1.8000	

05/17/00 02:06:42 PM page 2

IntStd Mode	1 Counts	2 NOTUSED	3 NOTUSED	4 NOTUSED	5 NOTUSED	6 NOTUSED	7 NOTUSED
Elem	Y			<del></del>		<del>-</del> -	
Wavlen	371.030						
Avge	16572						~ ~
SDev	48.29594				<b>-</b> -		
%RSD	.2914284	₩ ₩		<b>40</b> 40			
#1	16538		<del>-</del> -		<del></del>		
#2	16606		<b></b>	<del>-</del> -			

5188

page 1

Method: METTRA Sample Name: CCB7 Operator: RJG

Run Time: 05/17/00 14:06:46

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT TRACEICP

Elem	AG	AL	AS	BA	BE	CA	CD
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00004	.00733	00070	.00062	.00027	.04215	.00024
SDev	.00028	.00665	.00088	.00041	.00013	.01807	.00001
%RSD	765.31	90.725	125.58	66.231	47.819	42.859	5.8348
#1	00024	.00263	00132	.00033	.00018	.02938	.00025
#2	.00016	.01204	00008		.00037	.05493	.00023
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.01000	.20000	.00500	5.0000	.00500
Low	01000	20000	01000	20000	00500	-5.0000	00500
Elem	CO	CR	CU	FE	MG	MN	MO
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00008	.00050	.00129	.00482	.01578	.00068	.00613
SDev	.00025	.00035	.00011	.00373	.00943	.00032	.00204
%RSD	321.96	68.728	8.4295	77.368	59.715	46.479	33.204
#1	00025	.00026	.00121	.00746	.00912	.00045	.00757
#2	.00010	.00075	.00137	.00218	.02245		.00469
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.05000	.01000	.02500	.10000	5.0000	.01500	.04000
Low	05000	01000	02500	10000	-5.0000	01500	04000
Elem	NI	PB/1	PB/2	PB	SB/1	SB/2	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00028	00036	.00168	.00100	.00070	.00012	.00032
SDev	.00006	.00089	.00051	.00005	.00107	.00160	.00071
%RSD	21.548	243.65	30.671	4.8241	151.95	1300.1	225.35
#1	00024	.00026	.00131	.00096	00005	.00126	.00082
#2	00032	00099	.00204	.00103	.00146	00101	00019
Errors High Low	LC Pass .04000 04000	NOCHECK	NOCHECK	LC Pass .00300 00300	NOCHECK	NOCHECK	LC Pass .06000 06000
Elem	SE/1	SE/2	SE	TL	V_	ZN	
Units	ppm	ppm	ppm	ppm	ppm	ppm	
Avge	.00378	00037	.00101	.00334	.00052	.00099	
SDev	.00199	.00226	.00085	.00127	.00000	.00037	
%RSD	52.607	616.54	83.473	37.887	.44291	36.992	
#1	.00518	00196	.00042	.00245	.00052	.00073	
#2	.00237	.00123	.00161	.00424	.00053	.00125	
Errors High Low	NOCHECK	NOCHECK	LC Pass .00500 00500	LC Pass .01000 01000	LC Pass .05000 05000	LC Pass .02000 02000	

05/17/00 02:10:52 PM page 2

IntStd Mode Elem Wavlen Avge SDev %RSD	1 Counts Y 371.030 16530 104.1574 .6301245	2 NOTUSED   	3 NOTUSED   	4 NOTUSED    	5 NOTUSED   	6 NOTUSED	7 NOTUSED    
#1	16603			<del>-</del> -			
#2	16456						

5190

# Log Méfals Preparation

Analyst Method:

Matrix:

Start Time: Lot Number:

020

SDG:

Pate: 5-16-00

COE 1930194

Lab Lo! No. (book, page, line)
96101-177-15

Soul Concl

5240N49A07 6623U3 508



STL Pittsburgh 450 William Pitt Way 412-820-8380 Pittsburgh, PA 15238

All Mark 5-16-00	Digestate(s)	24.	23	22	21.	20	19	18	17.	16.	15	14	13	12.	11	10.		8 TOWARC	<b>沙</b>	6 50,47	75/7	77.0	3 776447	2 77.445	1. 772°B	( Sample in	Sample ID	men / / / now	Reviewed Bu:
0140 01110 A	icestate																*	<b>*</b>		_				6	3	g/mL	Init WIVVol	S=16-00	2
West y	Received							d	Jost 2	July 1							•	•							1,774	mL .	Final Vol	·	
MCHON 2-1/10 10:05															\			13818-511 3m1+		2008	0.0	115	N. 6	1 mg NSB12EV	(20 c) 4 = 17 6	Comments	Comments		の方になっている
Many apple	Divestate(s) Relinauished								-																		Color	MYBIR	MS
WE'TO	7																						-			Post Pre		eV	
Texture F=Fine M=Medium C=Coarse	Hot plate													176	1											Post	Clarity	27	
Clarity C=Cloudy C=Cloudy O=Opaque														5-16 PD												Pre Post	Texture	ional Conc A	)Me 30% 1120
R=Red BL=Blue BR=Brown BLK=Black Y=Yellow O=Orange	ion																									-	Artifacts		8
ad V=Violet Iue P=Pink Own W=White lack GY=Gray Ow GN=Green nge C=Colorless	Color															l=Insects	P=Paper	C=Cloth	R=Rubber/Plastic	M=Metal Frgmts	G=Glass	W=free H2O	(plant mat'l)	O=Organic	S=Stones		Artifact Codes	Nallanchadt 5587 N41 NO7	Chrost 52401

PAGE 95 of 100

N'\QA\FORMS\METALS\_PREP

LOGBOOK # IP-00-0026

PSR024 5/16/00

8:40:23 MT

SAMPLE CUSTODIAN REMOVAL REQUEST

PAGE 001

REQUESTED BY: FAUSTE

METHÓD: QM

Inductively Coupled Plasma (6010B Trace)

STORAGE LOCATION	WORK ORDER #	PICKED CNTR#	CONTROL #	CLIENT #	ANALYSIS	LOTID	SMP#	<u>sfx</u>	MATRIX DESCRIPTION		D REG	
2F CLP1	DD6A4		233877	399411	A-46-QM	COE130194	001	SOLID		0	3	1
2F CLP1	DD6A5		233878	399411	A-46-QM	COE <b>13019</b> 4	002	SOLID		0	3	1
2F CLP1	DD6A6		233879	39941 <b>1</b>	A-46-QM	COE130194	003	SOLID	1	0	3	1
2F CLP1	DD6A7		233880	<b>399</b> 411	A-46-QM	c0E130194	004	SOLID		0	3	1

RELINQUISHED BY	RECEIVED BY RECEIVED BY	5-1600 1050 5-1600 1315
EwiM Faust	t Cen	<u>51600 1315</u>

#### TCLP METALS DATA

## Cover Page - Inorganic Analysis Data Package

	Client ID	Lab Sample ID:
	DF/S1/0133/SDC/001 A	DD6A4T
	DF/S1/0133/SDC/001 AD	DD6A4DT
	DF/S1/0133/SDC/001 AS	DD6A4ST
	DF/\$1/0133/SDC/001 B	DD6A5T
	DF/S1/0133/SDC/001 C	DD6A6T
	DF/S1/0133/SDC/001 D	DD6A7T
technically at	data package and in the computer-readable of	detailed above. Release of the data combined in data submitted on diskette has been authorized by
the Laborator	ry Manager or the Manager's designee, as ver	ified by the following signature.
Signature:		Name:
Date:		Title:

REVIEWED BY:
DATE:

Version 3.63.3

Cover Page Equivalent

TCLP METALS RESULTS

١

STL Pittsburgh 6003

#### Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A4T Client ID: DF/S1/0133/SDC/001 A

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139097

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.000045	0.00020	0 000045	U	1	CVAA	5/18/00	11:03

Comments:

#### Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A5T Client ID: DF/S1/0133/SDC/001 B

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139097

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.000045	0.00020	0.000045	บ	1	CVAA	5/18/00	11:09

Version 3.63.3

657 800

#### STL-Pittsburgh

## Metals Data Reporting Form

Sample Results

DD6A6T Lab Sample ID:

Client ID:

DF/S1/0133/SDC/001 C

Matrix:

Water

mg/L Units:

Prep Date:

5/18/00

Prep Batch:

0139097

Weight:

NA

Volume:

100

Percent Moisture:

NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.000045	0.00020	0.000045	U	1	CVAA	5/18/00	11:10

Version 3 63 3

657 801

#### Metals Data Reporting Form

 Sample Results

 Lab Sample ID:
 DD6A7T
 Client ID:
 DF/S1/0133/SDC/001 D

 Matrix:
 Water
 Units:
 mg/L
 Prep Date:
 5/18/00
 Prep Batch:
 0139097

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0 000045	0 00020	0.000045	U	1	CVAA	5/18/00	11:13

657 802

## Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A4T Client ID: DF/S1/0133/SDC/001 A

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	193.7	0.030	0.50	0.15	В	1	ICP	5/19/00	9:14
Barium	493.41	0.00041	10.0	1.2	BE	1	ICP	5/19/00	9:14
Cadmium	228.80	0.0028	0.10	0.0028	U	1	ICP	5/19/00	9:14
Chromium	267.72	0.0038	0.50	0.0038	U	1	ICP	5/19/00	9:14
Lead	220.35	0.025	0.50	0.025	Ų	1	ICP	5/19/00	9:14
Selenium	196.03	0.067	0.25	0.067	U	1	ICP	5/19/00	15:29
Silver	328.07	0.0031	0.50	0.0031	U	1	ICP	5/19/00	9:14

Comments: C0E130194001

U Result is less than the MDL

Form I Equivalent

#### Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A5T Client ID: DF/S1/0133/SDC/001 B

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Cone	Q	DF	Instr	Anal Date	Anal Time
Arsenic	193.7	0.030	0.50	0.21	В	1	ICP	5/19/00	9:27
Barium	493.41	0.00041	10.0	1.2	BE	1	ICP	5/19/00	9:27
Cadmium	228.80	0.0028	0.10	0.0028	U	1	ICP	5/19/00	9:27
Chromium	267.72	0.0038	0.50	0.0038	U	1	ICP	5/19/00	9:27
Lead	220.35	0.025	0 50	0.025	U	1	ICP	5/19/00	9:27
Selenium	196.03	0.067	0 25	0.067	บ	1	ICP	5/19/00	15:42
Silver	328.07	0.0031	0.50	0.0031	U	1	ICP	5/19/00	9:27

657 804

#### Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A6T Client ID: DF/S1/0133/SDC/001 C

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	193.7	0.030	0.50	0.20	В	1	ICP	5/19/00	9:36
Barium	493.41	0.00041	10.0	1.3	BE	1	ICP	5/19/00	9:36
Cadmium	228.80	0.0028	0.10	0.0028	υ	1	ICP	5/19/00	9:36
Chromium	267.72	0.0038	0.50	0.0038	บ	1	ICP	5/19/00	9:36
Lead	220.35	0.025	0.50	0.025	U	1	ICP	5/19/00	9:36
Selenium	196.03	0 067	0.25	0.067	U	1	ICP	5/19/00	15:51
Silver	328.07	0.0031	0.50	0.0031	U	1	ICP	5/19/00	9:36

Comments: <u>C0E130194003</u>

## Metals Data Reporting Form

Sample Results

Lab Sample ID: DD6A7T Client ID: DF/S1/0133/SDC/001 D

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF_	Instr	Anal Date	Anal Time
Arsenic	193,7	0.030	0.50	0.15	В	1	ICP	5/19/00	9:39
Barium	493.41	0.00041	10.0	1.2	BE	1	ICP	5/19/00	9:39
Cadmium	228 80	0 0028	0.10	0.0028	U	1	ICP	5/19/00	9:39
Chromium	267 72	0 0038	0.50	0.0038	U	ì	ICP	5/19/00	9:39
Lead	220.35	0.025	0.50	0.025	U	1	ICP	5/19/00	9:39
Selenium	196.03	0 067	0.25	0.067	U	1	ICP	5/19/00	15:55
Silver	328.07	0.0031	0.50	0.0031	υ	1	ICP	5/19/00	9:39

657 806

## Metals Data Reporting Form

Initial Calibration Verification Standard

Instrument: CVAA Units: ug/L

Chart Number: 0518HGA.PRN Acceptable Range: 90% - 110%

Standard Source: Ultra Standard ID: 0014-094-5

	WL/		ICV5-1 5/18/00 10:49 AM									
Element	Mass	True Conc	Found	% Rec	Found	% Rec	Found	% Rec	Found	% Rec	Found	% Rec
Mercury	253.7	2.5	2 54	101.6								1

## Metals Data Reporting Form

Initial Calibration Verification Standard

Instrument: ICP Units: ug/L

Chart Number: J00519A.ARC Acceptable Range: 90% - 110%

Standard Source: Inorganic Ventures Standard ID: 0014-061-7

	33/1 /		ICV2- 5/19/00 8:50 Al	0								±4
Element	WL/ Mass	True Conc	Found	% Rec	Found	% Rec	Found	% Rec	Found	% Rec	Found	% Rec
Arsenic	193.696	1000.0	1014.30	101.4								
Barium	493.409	1000.0	1019.97	102.0				I				
Cadmium	228.802	1000.0	1021.96	102.2		Į						
Chromium	267.716	1000.0	1017.40	101.7		1		!				
Lead	220 353	1000.0	1044.87	104.5								
Silver	328.068	500.0	510.86	102.2								

657 808

## Metals Data Reporting Form

Initial Calil	oration \	erifica	tion Stand	dard								
Instrument:	IC	CP					Units:	u	ıg/L	_		
Chart Numb	er: <u>J00</u>	519B.A	RC_				Accepta	ble Ra	nge: 9	0% - 11	10%	
Standard So	ource:	In	organic Ve	ntures	<del></del>		Standar	d ID: _	0	014-06	1-7	<del></del>
	WL/	True	ICV2- 5/19/00 3:05 PM	0 M %		%	~	%	Found	% Rec	Found	% Rec
Element	Mass	Conc	Found	Rec	Found	Rec	Found	Rec	Found	Rec	round	Nec
Selenium	196,026	1000.0	1015.86	1015.86 101.6						1		

## Metals Data Reporting Form

Continuing Calibration Verification Units: ug/L **CVAA** Instrument: Acceptable Range: 80% - 120% Chart Number: 0518HGA.PRN Standard ID: 0014-094-6 Inorganic Ventures Standard Source: CCV5-2 CCV5-1 5/18/00 5/18/00 10:53 AM 11:17 AM WL/ % % % % % True Found Rec **Found** Rec Rec Found Rec Rec Found Conc Found Element Mass 5.19 103.8 5.21 104.2 5.0 253.7 Mercury

657 810

## Metals Data Reporting Form

Continuing Calibration Verification

Instrument: ICP Units: ug/L

Chart Number: J00519A.ARC Acceptable Range: 90% - 110%

Standard Source: Inorganic Ventures Standard ID: 0014-087-7

			CCV2-1 5/19/00 9.30 AN	, [	CCV2-7 5/19/00 9·43 AM	)						
	WL/	True		%		%		%	T	%	Tournel	% Rec
Element	Mass	Conc	Found	Rec	Found	Rec	Found_	Rec	Found	Rec	Found	Rec
Arsenic	193.696	5000.0	4996 70	99.9	5021.14	100.4						
Barium	493.409	5000.0	4763.98	95.3	4802.08	96.0		ļ				
Cadmium	228.802	5000.0	4980.88	99.6	5023.14	100.5						
Chromium	267.716	5000.0	4856.58	97.1	4902.29	98.0		Ì		- 1		
Lead	220.353	5000.0	4889.67	97.8	4941.05	98.8		-		l l		
Silver	328.068	1000 0	977.94	97.8	984.69	98.5						

## Metals Data Reporting Form

Continuing Calibration Verification Units: ug/L Instrument: ICP Acceptable Range: 90% - 110% Chart Number: J00519B.ARC **Standard ID:** 0014-087-7 Inorganic Ventures Standard Source: CCV2-1 CCV2-2 5/19/00 5/19/00 3:58 PM 3:45 PM WL/ % % % % % True Found Rec Found Rec Rec Found Rec Found Rec Found Element Mass Conc 5249.79 **105.0** 5218.77 104.4 196.026 5000.0 Selenium

# 657 812

## STL-Pittsburgh

## Metals Data Reporting Form

Initial Calil	bration B	lank R	esults									
Instrument:	CV	AA	_				Units:	·	ug/L	_		
Chart Numb	er: 0518	HGA.P	RN_									
Standard So	urce:						Standar	d ID:				_
			ICB1 5/18/0 10:51 A	0								
WL/ Report Limit Found Q Fo			Found	Q	Found	Q	Found	Q	Found	Q		
Mercury	253.7	0.2	0.0	В		-					<u> </u>	

Form 3 Equivalent

U Result is less than the MDL

## Metals Data Reporting Form

Initial Calib	ration B	lank Re	esults		<del></del>		<del></del> ·				·							
Instrument:	IC	P	•				Units:		ug/L									
Chart Numb	er: <u>J005</u>	19A.AI	RC															
Standard So	urce:						Standar	d ID:										
	ICB1 5/19/00 8:53 AM																	
Element	WL/ Mass	Report Limit	Found	Q	Found	Q_	Found	Q	Found	Q	Found	Q						
Arsenic	193.696	500	30.3	υ														
Barium	493.409	10000	0.4	U														
Cadmium	228.802	100	2.8	U					}		<u> </u>							
Chromium	267.716	500	3.8	TI	!		I				İ							

500

500

220.353

328.068

Lead

Silver

U

U

24.6

3.1

U Result is less than the MDL

657 814

# Metals Data Reporting Form

Initial Calib	oration B	lank R	esults									
Instrument:	IC	P	<del>-</del>				Units:	· · · · · · · · · · · · · · · · · · ·	ug/L	<b>-</b>		
Chart Numb	er: <u>J00</u> 5	519B.AI	RC_									
Standard So	urce:	<del></del>		<del></del>			Standar	rd ID:				
Element	WL/ Mass	Report Limit	ICB1 5/19/00 3:08 PM		Found	Q	Found	Q_	Found	Q	Found	Q_
Selenium	196.026			U								

B Result is between MDL and RL

## Metals Data Reporting Form

Continuin	g Calibra	tion B	ank Res	ults			<del></del>		<del></del> -			
Instrument	t: <u>C</u>	7AA	<u>.</u>				Units:		ug/L			
Chart Nun	nber: 051	8HGA.	PRN									
Standard S	Source:				<u> </u>	Standa	ard ID	);			<u></u>	
				1 00 AM	CCB: 5/18/0 11:18 /	00						
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	Q	Found	Q	Found	Q
Maroury	253.7	0.2		) []	-0.1	l B			[		l .	

U Result is less than the MDL

B Result is between MDL and RL

## Metals Data Reporting Form

			CCB1 5/19/00 9:33 AM		CCB2 5/19/00 9:46 AM							
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	Q	Found	Q	Found	Q
Arsenic	193.696	500	32.7	В	30.3	Ü						
Barium	493.409	10000	0.8	В	0.6	В			ļ			
Cadmium	228.802	100	2.8	U	2.8	Ü			ļ			
Chromium	267.716	500	3.8	U	3.8	U	}		]			
Lead	220.353	500	24.6	IJ	24.6	Ų						
Silver	328.068	500	3.1	U	3.1	U			L			

Form 3 Equivalent

Version 3.63.3

U Result is less than the MDL

## Metals Data Reporting Form

Continuing	g Calibra	tion B	ank Resu	lts								
Instrument	: <u>I</u> (	CP	_				Units:		ug/L	<del></del>		
Chart Num	ber: J00	)519B.A	RC_									
Standard S	ource:						Standa	rd ID	:			
			CCB1 5/19/00 3:48 PM		CCB2 5/19/00 4:01 PM	)		·				
Element	WL/ Mass	Report Limit	Found	Q	Found	Q	Found	Q	Found	Q	Found	Q
Selenium	196.026	250	67.4	U	67.4	U						

U Result is less than the MDL

## Metals Data Reporting Form

Preparation Blank Results

Lab Sample ID: DDA9GBT

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139097

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.000045	0.00020	0.000063	В	1	CVAA	5/18/00	11:01

#### Metals Data Reporting Form

Preparation Blank Results

Lab Sample ID:

DDC2GBT

Matrix: Water

Units:

mg/L

Prep Date:

5/18/00

Prep Batch:

0139097

Weight:

NA

Volume:

100

Percent Moisture:

NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.000045	0.00020	-0.000090	В	1	CVAA	5/18/00	10:57

# 657 820

## Metals Data Reporting Form

Preparation Blank Results

Lab Sample ID: DDA9GBT

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Cone	Q	DF	Instr	Anal Date	Anal Time
Arsenic	193.696	0.030	0.50	0.18	В	1	ICP	5/19/00	9:08
Barium	493.409	0.00041	10.0	0.037	В	1	ICP	5/19/00	9:08
Cadmium	228.802	0.0028	0.10	0.0028	υ	1	ICP	5/19/00	9:08
Chromium	267.716	0.0038	0.50	0.0038	U	1	ICP	5/19/00	9:08
Lead	220.353	0.025	0.50	0.025	U	1	ICP	5/19/00	9:08
Selenium	196.026	0.067	0.25	0.067	U	1	ICP	5/19/00	15:23
Silver	328.068	0.0031	0.50	0.0031	U	1	ICP	5/19/00	9.08

## Metals Data Reporting Form

Preparation Blank Results

Lab Sample ID: DDC39BT

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MDL	Report Limit	Conc	Q	DF	Instr	Anal Date	Anal Time
Arsenic	193.696	0.030	0 50	0.030	U	1	ICP	5/19/00	9:05
Barium	493.409	0.00041	10.0	0.00041	U	1	ICP	5/19/00	9:05
Cadmium	228.802	0.0028	0.10	0.0028	U	1	ICP	5/19/00	9:05
Chromium	267.716	0.0038	0.50	0.0038	U	1	ICP	5/19/00	9:05
Lead	220.353	0.025	0.50	0.025	υ	1	ICP	5/19/00	9:05
Selenium	196.026	0.067	0.25	0.067	U	1	ICP	5/19/00	15:20
Silver	328.068	0.0031	0.50	0.0031	<b>U</b> _	1	ICP	5/19/00	9:05

657 822

## Metals Data Reporting Form

Interference Check Standard A

Instrument: ICP Units: ug/L

Chart Number: J00519A.ARC Acceptable Range: 0% - 0%

Standard Source: Inorganic Ventures Standard ID: 0014-088-12

				ICSA 5/19/00 8:56 AM				
Element	WL/ Mass	Reporting Limit	True Conc	Found	Found	Found	Found	Found
Arsenic	193.696	500		30			·	
Barium	493.409	10000		2				
Cadmium	228.802	100		3		<u> </u>	1	
Chromium	267 716	500		4				
Lead	220.353	500		128				1
Silver	328.068	500		3		<u> </u>		<u> </u>

U Result is less than the MDL

B Result is between MDL and RL

657 823

# Metals Data Reporting Form

Interferenc	e Check S	tandard A				<u> </u>		
Instrument:	ICI	•			Units:	ug/L		
Chart Numl	ber: <u>J005</u>	19B.ARC			Accep	table Range:	0% - 0%	<del>_</del>
Standard Sc	ource:	Inorgai	nic Ventur	es	Standa	ard ID:	0014-088-1	2
				ICSA 5/19/00 3:11 PM				
Element	WL/ Mass	Reporting Limit	True Conc	Found	Found	Found	Found	Found
Selenium	196.026	250		89				

U Result is less than the MDL

B Result is between MDL and RL

657 824

## Metals Data Reporting Form

Interference Check Standard AB
Units: ug/L

Instrument: ICP Units: ug/L

Chart Number: J00519A.ARC Acceptable Range: 80% - 120%

Standard Source: Inorganic Ventures Standard ID: 0014-043-1

			ICSAB 5/19/00 8:59 AN							2		
Element	WL/ Mass	True Conc	Found	% Rec	Found_	% Rec	Found	% Rec	Found	% Rec	Found	% Rec
Arsenic	193.696	1000	871.4	87.1						1		
Barium	493 409	500	456.1	91.2				]				
Cadmium	228.802	1000	949.9	95.0								
Chromium	267.716	500	434.8	87.0				į		1		
Lead	220.353	1000	1039.8	104.0								
Silver	328.068	1000	944 4	94.4								

U Result is less than the MDL

B Result is between MDL and RL

N Spike recovery failed

## Metals Data Reporting Form

Interference	ce Check	Standa	rd AB									
Instrument	: <u>IC</u>	CP	_				Units:	l	ug/L	_		
Chart Num	ber: <u>J00</u>	519B.A	RC_				Accepta	ble Ra	nge:8	0% - 12	20%	
Standard S	ource:	In	organic V	entures			Standar	d ID:	0	014-04	3-1	_
		<del></del>	ICSA 5/19/0 3:14 P	00								
Element	WL/ Mass	True Conc	Found	% Rec	Found	% Rec	Found	% Rec	Found	% Rec	Found	% Rec
Selenium	196 026	1000	1089.8	3 109.0			- <u></u>				<u> </u>	

Version 3.63.3

U Result is less than the MDL

B Result is between MDL and RL

N Spike recovery failed

657 826

## Metals Data Reporting Form

Matrix Spike Sam	ple Results	<u> </u>				<del></del> -
Spike Sample ID:	DD6A4ST					
Original Sample ID:	DD6A4T	Client ID:	DF/S1/0133/	SDC/001 AS		
Matrix: Water	Units: mg/L	Prep Date:	5/18/00	Prep Batch:	0139097	<b></b>
Weight: NA	Volume: 100	Percent Mois	ture: NA	<u> </u>		
			T T T	OS	OS MS	MS

Element	WL/ Mass	OS Conc	0	MS Conc	Q	Spike Level	% Rec	OS DF	MS DF	Instr	OS Anal Date	OS Anal Time	MS Anal Date	MS Anal Time
Mercury	253.7	0.000045	U	0.0053		0.005	106.6	1	1	CVAA	5/18/00	11:03	5/18/00	11:04

Comments: \_

Version 3.63 3

U Result is less than the MDL

Form 5A Equivalent

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

Duplicate analysis RPD was not within limits

## Metals Data Reporting Form

Matrix Spike Duplicate Sample Result	Matrix	ix Spike	Duplicate	Sample	Result
--------------------------------------	--------	----------	-----------	--------	--------

Spike Sample ID:	DD6A4DT				
Original Sample ID:	DD6A4T	Client ID:	DF/S1/013	3/SDC/001 AD	
Matrix: Water	Units: mg/L	Prep Date:	5/18/00	Prep Batch:_	0139097
			STA		

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	OS Conc	Q	MSD Conc	Q	Spike Level	% Rec	OS DF	MSD DF	Instr	OS Anal Date	OS Anal Time	MSD Anal Date	MSD Anal Time
Mercury	253.7	0.000045	U	0.0052		0.005	104.8	l	1	CVAA	5/18/00	11:03	5/18/00	11:07

Comments: \_\_\_\_\_
Version 3.63.3

U Result is less than the MDL

Form 5A Equivalent

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

\* Duplicate analysis RPD was not within limits

#### 657 828

#### Metals Data Reporting Form

Matrix Spike Sample Results

Spike Sample ID: DD6A4ST\_

Original Sample ID: DD6A4T Client ID: DF/S1/0133/SDC/001 AS

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	OS Conc	Q	MS Conc	Q	Spike Level	% Rec	OS DF	MS DF	Instr	OS Anal Date	OS Anal Time	MS Anal Date	MS Anal Time
Arsenic	193.7	0.15	В	4.9		5	94.8	1	1	ICP	5/19/00	9:14	5/19/00	9:20
Barium	493.4	1.2	_	43.3	l	50	84.1	1	1	ICP	5/19/00	9:14	5/19/00	9:20
Cadmium	228.8	0.0028		0.94		1	94.4	1	1	ICP	5/19/00	9:14	5/19/00	9:20
Chromium	267.7	0.0038		4.3	ì	5	85.3	1	1	ICP	5/19/00	9:14	5/19/00	9:20
	220.4	0.025	Ψ.	4.4	l	5	88.7	1	1	ICP	5/19/00	9:14	5/19/00	9:20
Lead	196.0	0.023		1.1		1	104.8	1	1	ICP	5/19/00	15:29	5/19/00	15:36
Selenium Silver	328.1	0.0031	i .	0.90		1	90.4	1	1	ICP	5/19/00	9:14	5/19/00	9:20

Comments:

Version 3.63.3

Duplicate analysis RPD was not within limits

U Result is less than the MDL

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

#### Metals Data Reporting Form

Matrix Spike Duplicate Sample Results

Spike Sample ID: DD6A4DT

Original Sample ID: DD6A4T Client ID: DF/S1/0133/SDC/001 AD

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	OS Cone	Q	MSD Conc	Q	Spike Level	% Rec	OS DF	MSD DF	Instr	OS Anal Date	OS Anal Time	MSD Anal Date	MSD Anal Time
Arsenic	193.7	0.15	В	5.0		5	96.2	1	1	ICP	5/19/00	9:14	5/19/00	9:24
Barium	493.4	1.2		43.8		50	85.2	i	1	ICP	5/19/00	9:14	5/19/00	9:24
Cadmium	228.8	0.0028		0.96	1	1	95.5	1	1	ICP	5/19/00	9:14	5/19/00	9:24
Chromium	267.7	0.0038		4.3	ļ	5	86.3	1	i	ICP	5/19/00	9:14	5/19/00	9:24
Lead	220.4	0.025		4.5		5	89.8	1	1	ICP	5/19/00	9:14	5/19/00	9:24
Selenium	196.0	0.067	L	1.1	ł	1	105 2	1	1	ICP	5/19/00	15:29	5/19/00	15:39
Silver	328.1	0.0031	L	0.91		1	91.2	1	1	ICP	5/19/00	9:14	5/19/00	9:24

Comments:

Version 3.63.3

\* Duplicate analysis RPD was not within limits

U Result is less than the MDL

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

657 830

#### Metals Data Reporting Form

Matrix S	pike Duplic	cate RPL	) Re	eport									
Matrix Sp	ike Duplicate	Sample II	D:	)	DD6	A4DT		_					
Matrix Sp	ike Sample II	):	D	D6A4ST		Cli	ent III	):	DF/S1/0	133/SDC/	001 AD		
Matrix:	Water	Units:	n	ng/L		Prep Da	te: _	5/18	8/00	Prep	Batch:	01390	97
Weight: _	NA	Volume:		100		Percent l	Moist	ure: _	NA	<u> </u>			
Element	WL/ Mass	MS Conc	O	MSD Conc	o	RPD	MS DF	MSD DF	Instr	MS Anal Date	MS Anal Time	MSD Anal Date	MSD Anal Time
Mercury	253.7			0.0052		1.7%	1	1	CVAA	5/18/00	11:04	5/18/00	11:07

Comments.

Version 3.63.3

U Result is less than the MDL

Form 6 Equivalent

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

Duplicate analysis RPD was not within limits

#### Metals Data Reporting Form

Matrix Spike Duplicate RPD Report

Matrix Spike Duplicate Sample ID: DD6A4DT

Matrix Spike Sample ID: DD6A4ST Client ID: DF/S1/0133/SDC/001 AD

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	MS Conc	Q	MSD Conc	Q	RPD	MS DF	MSD DF	Instr	MS Anal Date	MS Anal Time	MSD Anal Date	MSD Anal Time
Arsenic	193,696	4.9		5 0		1.5 %	1	1	ICP	5/19/00	9:20	5/19/00	9:24
Barium	493.409	43.3		43.8		1.4%	1	1	ICP	5/19/00	9:20	5/19/00	9:24
Cadmium	228.802	0.94		0.96		1.1 %	1	1	ICP	5/19/00	9:20	5/19/00	9:24
Chromium	267.716		1	4.3		1.2 %	1	1	ICP	5/19/00	9:20	5/19/00	9:24
Lead	220.353		1	4.5		1.2%	1	i	ICP	5/19/00	9:20	5/19/00	9:24
Selenium	196.026	1.1	ļ	1.1		10.2 %	1	1	ICP	5/19/00	15:36	5/19/00	15:39
Silver	328.068	0.90		0.91		1.0 %	1	1	ICP	5/19/00	9:20	5/19/00	9:24

Comments:

Version 3.63.3

U Result is less than the MDL

B Result is between MDL and RL

N Spike recovery failed

NC Percent recovery was not calculated

\* Duplicate analysis RPD was not within limits

## Metals Data Reporting Form

Laboratory Control Sample Results

Lab Sample ID: DDC2GCT

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139097

Weight: NA Volume: 100 Percent Moisture: NA

Element	WL/ Mass	Spike Level	Conc	Percent Recovery	Q	Range	DF	Instr	Anal Date	Anal Time
Mercury	253.7	0.0025	0.0028	110.0		80-120	1	CVAA	5/18/00	10:58

#### Metals Data Reporting Form

Laboratory Control Sample Results

Lab Sample ID: DDC39CT

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	Spike Level	Cone	Percent Recovery	Q	Range	DF	Instr	Anal Date	Anal Time
Arsenic	193.696	2.0	2.2	110.1		80-120	1	ICP	5/19/00	9:11
Barium	493,409	2.0	1.9	93.7	В	80-120	1	ICP	5/19/00	9:11
Cadmium	228.802	0.050	0.046	91.3	В	80-120	1	ICP	5/19/00	9:11
Chromium	267.716	0.20	0.19	94.7	В	80-120	1	ICP	5/19/00	9:11
Lead	220.353	0.50	0.47	94.5	В	80-120	1	ICP	5/19/00	9:11
Selenium	196.026	2.0	2.2	107.8		80-120	1	ICP	5/19/00	15:26
Silver	328.068	0.050	0.047	94.3	В	80-120	1	ICP	5/19/00	9:11

657 834

#### Metals Data Reporting Form

Serial Dilution RPD Report

Serial Dilution Sample ID: DD6A4PT

Original Sample ID: DD6A4T Client ID: DF/S1/0133/SDC/001 A

Matrix: Water Units: mg/L Prep Date: 5/18/00 Prep Batch: 0139114

Weight: NA Volume: 50 Percent Moisture: NA

Element	WL/ Mass	OS Conc	Q	Serial Dilution Conc	Q	Percent Diff	OS DF	Ser Dil DF	Instr	OS Anal Date	OS Anal Time	Ser Dil Anal Date	Ser Dil Anal Time
Arsenic	193.696	0.15	В	0.17	В		1	5	ICP	5/19/00	9:14	5/19/00	9:17
Barium	493.409				ВΕ	11.4%	1	5	ICP	5/19/00	9:14	5/19/00	9:17
Cadmium	228.802	i		0.014	1		1	5	ICP	5/19/00	9:14	5/19/00	9:17
Chromium	267.716			0.019			1	5	ICP	5/19/00	9:14	5/19/00	9:17
Lead	220.353			0.13	1		1	5	ICP	5/19/00	9:14	5/19/00	9:17
Selenium	196.026		i	0.34	l		1	5	ICP	5/19/00	15:29	5/19/00	15:32
Silver	328.068			0.016	1		1	5	ICP	5/19/00	9:14	5/19/00	9:17

Comments:

Version 3.63.3

E Serial dilution percent difference not within limits

Form 9 Equivalent

U Result is less than the MDL

B Result is between MDL and RL

## Metals Data Reporting Form

Instrument Detection Limits

Instrument: CVAA

Units: ppb

Element	Wavelength /Mass	Reporting Limit	MDL	Date of MDL
Mercury	253.70	0.2	0.045	3/20/00

## Metals Data Reporting Form

Instrument Detection Limits

Instrument: ICP Units: ppb

Element	Wavelength /Mass	Reporting Limit	MDL	Date of MDL
Arsenic	193.70	500	30.3	4/1/00
Barium	493.41	10000	0.41	4/1/00
Cadmium	228.80	100	2.8	4/1/00
Chromium	267.72	500	3.8	4/1/00
Lead	220.35	500	24.6	4/1/00
Selenium	196.03	250	67.4	4/1/00
Silver	328.07	500	3.1	4/1/00

# Metals Data Reporting Form

Inter-Element Correction Factors

nstrument:	ICP	Date of IEC's: 3/27/00
Interfering Element	Wavelength /Mass	Correction Factor(s)
Aluminum	308.215	As(0.008365), Mn(0.00002), Pb(0.000527)
Antimony	206.838	Ni(-0.000449), Pb(-0.001338), Sn(-0.004668)
Arsenic	193.696	Cd(0.011196)
Barium	493.409	Co(0.000506)
Beryllium	313.042	Cd(0.008625)
Cadmium	228.802	Co(0.002633)
Chromium	267.716	Pb(-0.000686), Sb(0.008213), V(-0.001979)
Cobalt	228.616	Al(-0.014067), B(0.00201), Cd(-0.004523), Cu(-0.00091), Pb(-0.027395), Sb(-0.003935), Tl(0.007862)
Copper	324.754	Zn(0.00466)
Iron	259.94	Ag(-0.000239), As(0.001314), B(-0.001921), Cd(-0.000034), Cu(-0.00008), Mn(-0.000288), Mo(-0.00015), Pb(0.000225), Se(-0.003656), Sn(-0.00019), Tl(0.010326) Zn(0.000098)
Manganese	257.61	Ag(0.000201), Tl(-0.005634)
Molybdenum	202.03	Al(0.008699), Cr(-0.000292), Mn(-0.00033), Sb(0.005808), V(-0.019318)
Nickel	231.604	Cd(-0.000409), Sb(-0.009092), Zn(0.003263)
Tin	189.989	Sb(0.002262)
Titanium	334.941	Co(0.001637), Fe(-0.003475), Sb(0.001696), Sn(0.003624)
Vanadium	292.402	Ag(-0.005069), Al(0.012877), As(0.017242), Be(0.00265), Cd(0.000094), Cr(0.000568), Sb(-0.003793), Si(-0.012762), Tl(0.007148), Zn(-0.004494)

#### Metals Data Reporting Form

Linear Dynamic Ranges

Instrument: CVAA

Units: ppb

Element	Wavelength	Linear	Date of Linear
	/Mass	Range	Range
Mercury	253.70	10	1/15/00

## Metals Data Reporting Form

Linear Dynamic Ranges

Instrument: ICP Units: ppb

Element	Wavelength /Mass	Linear Range	Date of Linear Range
Arsenic	193 70	100000	4/5/00
Barium	493.41	100000	4/5/00
Cadmium	228.80	100000	4/5/00
Chromium	267.72	100000	4/5/00
Lead	220.35	100000	4/5/00
Selenium	196.03	100000	4/5/00
Silver	328.07	2000	4/5/00

657 840

#### Metals Data Reporting Form

Preparation Log

Preparation Batch: 0139097 Instrument: CVAA Matrix: Water

Sample ID	Prep Date	Weight (g)	Volume (ml)	% Moisture
DDA9GBT	5/18/00	NA	100	NA
DDC2GBT	5/18/00	NA	100	NA
DDC2GCT	5/18/00	NA	100	NA
DD6A4DT	5/18/00	NA	100	NA
DD6A4ST	5/18/00	NA	100	NA
DD6A4T	5/18/00	NA	100	NA
DD6A5T	5/18/00	NA	100	NA
DD6A6T	5/18/00	NA	100	NA
DD6A7T	5/18/00	NA	100	NA

#### Metals Data Reporting Form

657 841

Preparation Log

Preparation Batch: 0139114 Instrument: ICP Matrix: Water

		,		<del></del>
Sample ID	Prep Date	Weight (g)	Volume (ml)	% Moisture
DDA9GBT	5/18/00	NA	50	NA
DDC39BT	5/18/00	NA	50	NA
DDC39CT	5/18/00	NA	50	NA
DD6A4DT	5/18/00	NA	50	NA
DD6A4ST	5/18/00	NA	50	NA
DD6A4T	5/18/00	NA	50	NA
DD6A5T	5/18/00	NA	50	NA
DD6A6T	5/18/00	NA	50	NA
DD6A7T	5/18/00	NA	50	NA

657 842

## Metals Data Reporting Form

Instrument Runlog

Instrument: CVAA

Chart Number: 0518HGA.PRN

<del></del>		
<del> </del>	Date of	Time of
Sample Name	Analysis	Analysis
Std1Rep1	5/18/00	10.34 AM
Std2Rep1	5/18/00	10:37 AM
Std3Rep1	5/18/00	10:39 AM
Std4Rep1	5/18/00	10:41 AM
Std5Rep1	5/18/00	10:43 AM
Std6Rep1	5/18/00	10:46 AM
ICV5-1	5/18/00	10:49 AM
ICB1	5/18/00	10:51 AM
CCV5-1	5/18/00	10:53 AM
CCB1	5/18/00	10:55 AM
DDC2GBT	5/18/00	10.57 AM
DDC2GCT	5/18/00	10:58 AM
DDA9GBT	5/18/00	11:01 AM
DD6A4T	5/18/00	11:03 AM
DD6A4ST	5/18/00	11:04 AM
DD6A4DT	5/18/00	11:07 AM
DD6A5T	5/18/00	11:09 AM
DD6A6T	5/18/00	11:10 AM
DD6A7T	5/18/00	11:13 AM
ZZZZZZ	5/18/00	11:15 AM
CCV5-2	5/18/00	11:17 AM
CCB2	5/18/00	11:18 AM
ZZZZZZ	5/18/00	11:21 AM
ZZZZZZ	5/18/00	11:23 AM
ZZZZZZ	5/18/00	11:25 AM
ZZZZZZ	5/18/00	11:27 AM
ZZZZZZ	5/18/00	11:28 AM
ZZZZZZ	5/18/00	11:30 AM
ZZZZZZ	5/18/00	11:32 AM
ZZZZZZ	5/18/00	11:34 AM
ZZZZZZ	5/18/00	11:36 AM
ZZZZZZ	5/18/00	11:38 AM
ZZZZZZ	5/18/00	11:40 AM
ZZZZZZ	5/18/00	11:42 AM
ZZZZZZ	5/18/00	11.44 AM
ZZZZZZ	5/18/00	11:46 AM
ZZZZZZ	5/18/00	11:48 AM
ZZZZZZ	5/18/00	11:50 AM
ZZZZZZ	5/18/00	11:52 AM
ZZZZZZ	5/18/00	11:54 AM
ZZZZZZ	5/18/00	11:56 AM

#### Metals Data Reporting Form

657 843

Instrument Runlog

Instrument: CVAA Chart Number: 0518HGA.PRN

	Date of	Time of
Sample Name	Analysis	Analysis
ZZZZZZ	5/18/00	11:58 AM
ZZZZZZ	5/18/00	12:00 PM
ZZZZZZ	5/18/00	12:02 PM
ZZZZZZ	5/18/00	12:04 PM
ZZZZZZ	5/18/00	12:06 PM
ZZZZZZ	5/18/00	12:10 PM
ZZZZZZ	5/18/00	12:12 PM
ZZZZZZ	5/18/00	12:14 PM
ZZZZZZ	5/18/00	12:16 PM
ZZZZZZ	5/18/00	12:19 PM
ZZZZZZ	5/18/00	12:21 PM
ZZZZZZ	5/18/00	12:23 PM
ZZZZZZ	5/18/00	12:25 PM
ZZZZZZ	5/18/00	12:28 PM
ZZZZZZ	5/18/00	12:30 PM
ZZZZZZ	5/18/00	12:32 PM
ZZZZZZ	5/18/00	12:34 PM
ZZZZZZ	5/18/00	12:36 PM
ZZZZZZ	5/18/00	12:38 PM
ZZZZZZ	5/18/00	12:40 PM
ZZZZZZ	5/18/00	12:43 PM
ZZZZZZ	5/18/00	12:45 PM
ZZZZZZ	5/18/00	2:07 PM
ZZZZZZ	5/18/00	2:09 PM
ZZZZZZ	5/18/00	2:11 PM
ZZZZZZ	5/18/00	2:13 PM
ZZZZZZ	5/18/00	2:16 PM
ZZZZZZ	5/18/00	2:18 PM
ZZZZZZ	5/18/00	2:20 PM
ZZZZZZ	5/18/00	2:22 PM
ZZZZZZ	5/18/00	2:24 PM
ZZZZZZ	5/18/00	2:26 PM

#### Metals Data Reporting Form

657 844 Instrument Runlog

Instrument: ICP Chart Number: J00519A.ARC

	Date of	Time of
Sample Name	Analysis	Analysis
ZZZZZZ	5/19/00	7:48 AM
ZZZZZZ	5/19/00	7:52 AM
ZZZZZZ	5/19/00	7:55 AM
ZZZZZZ	5/19/00	7:58 AM
ZZZZZZ	5/19/00	8:11 AM
ZZZZZZ	5/19/00	8:14 AM
ZZZZZZ	5/19/00	8:17 AM
ZZZZZZ	5/19/00	8:20 AM
ZZZZZZ	5/19/00	8:25 AM
ZZZZZZ	5/19/00	8:28 AM
ZZZZZZ	5/19/00	8:32 AM
ZZZZZZ	5/19/00	8:35 AM
STD1	5/19/00	8:40 AM
STD5A	5/19/00	8:43 AM
STD5B	5/19/00	8:47 AM
ICV2-1	5/19/00	8·50 AM
ICB1	5/19/00	8·53 AM
ICSA	5/19/00	8:56 AM
ICSAB	5/19/00	8:59 AM
DDC39BT	5/19/00	9:05 AM
DDA9GBT	5/19/00	9:08 AM
DDC39CT	5/19/00	9:11 AM
DD6A4T	5/19/00	9:14 AM
DD6A4PT	5/19/00	9:17 AM
DD6A4ST	5/19/00	9:20 AM
DD6A4DT	5/19/00	9:24 AM
DD6A5T	5/19/00	9:27 AM
CCV2-1	5/19/00	9:30 AM
CCB1	5/19/00	9:33 AM
DD6A6T	5/19/00	9:36 AM
DD6A7T	5/19/00	9:39 AM
CCV2-2	5/19/00	9:43 AM
CCB2	5/19/00	9:46 AM
ZZZZZZ	5/19/00	9:49 AM
ZZZZZZ	5/19/00	9:52 AM
ZZZZZZ	5/19/00	9:55 AM
ZZZZZZ	5/19/00	9:58 AM
ZZZZZZ	5/19/00	10:02 AM
ZZZZZZ	5/19/00	10:05 AM
ZZZZZZ	5/19/00	10:08 AM
ZZZZZZ	5/19/00	10:11 AM

## Metals Data Reporting Form

657 845

Instrument Runlog

ICP Instrument:

Chart Number: J00519A.ARC

Sample Name	Date of Analysis	Time of Analysis
Sample Name		
ZZZZZZ	5/19/00	10:14 AM
ZZZZZZ	5/19/00	10:17 AM
ZZZZZZ	5/19/00	10:20 AM
ZZZZZZ	5/19/00	10:23 AM
ZZZZZZ	5/19/00	10:27 AM
ZZZZZZ	5/19/00	10:30 AM
ZZZZZZ	5/19/00	10:33 AM
ZZZZZZ	5/19/00	10:36 AM
ZZZZZZ	5/19/00	10:39 AM
ZZZZZZ	5/19/00	10:42 AM
ZZZZZZ	5/19/00	10:46 AM
ZZZZZZ	5/19/00	10:49 AM
ZZZZZZ	5/19/00	10:52 AM
ZZZZZZ	5/19/00	10:59 AM
ZZZZZZ	5/19/00	11·02 AM
ZZZZZZ	5/19/00	11:05 AM
ZZZZZZ	5/19/00	11:08 AM
ZZZZZZ	5/19/00	11:11 AM
ZZZZZZ	5/19/00	11:15 AM
ZZZZZZ	5/19/00	11:18 AM
ZZZZZZ	5/19/00	11:21 AM
ZZZZZZ	5/19/00	11:24 AM
ZZZZZZ	5/19/00	11:32 AM
ZZZZZZ	5/19/00	11:35 AM
ZZZZZZ	5/19/00	11:38 AM
ZZZZZZ	5/19/00	11:41 AM
ZZZZZZ	5/19/00	11:44 AM
ZZZZZZ	5/19/00	11:48 AM
ZZZZZZ	5/19/00	11:54 AM
ZZZZZZ	5/19/00	11:57 AM
ZZZZZZ	5/19/00	12:00 PM
ZZZZZZ	5/19/00	12:03 PM
ZZZZZZ	5/19/00	12:06 PM
ZZZZZZ	5/19/00	12:10 PM
ZZZZZZ	5/19/00	12:13 PM
ZZZZZZ	5/19/00	12:16 PM
ZZZZZZ	5/19/00	12:19 PM
ZZZZZZ	5/19/00	12:22 PM
ZZZZZZ	5/19/00	12:26 PM
ZZZZZZ	5/19/00	12:29 PM
ZZZZZZ	5/19/00	12:32 PM

Form 14 Equivalent

657 846

## Metals Data Reporting Form

Instrument Runlog

Instrument: ICP Chart Number: J00519A.ARC

Sample Name	Date of Analysis	Time of Analysis		
7.Z.Z.Z.Z	5/19/00	12:35 PM		
ZZZZZZ ZZZZZZ	5/19/00	12:38 PM		
7.7.7.7.Z.Z	5/19/00	12:42 PM		
 ZZZZZZ	5/19/00	12:45 PM		
ZZZZZZ	5/19/00	12:48 PM		

## Metals Data Reporting Form

Instrument Runlog

Instrument: ICP Chart Number: J00519B.ARC

Sample Name	Date of Analysis	Time of Analysis
ZZZZZZ	5/19/00	2.41 PM
ZZZZZZ	5/19/00	2:44 PM
ZZZZZZ	5/19/00	2:47 PM
ZZZZZZ	5/19/00	2:51 PM
STD1	5/19/00	2:55 PM
STD5A	5/19/00	2:59 PM
STD5B	5/19/00	3:02 PM
ICV2-1	5/19/00	3:05 PM
ICB1	5/19/00	3.08 PM
ICSA	5/19/00	3:11 PM
ICSAB	5/19/00	3:14 PM
DDC39BT	5/19/00	3:20 PM
DDA9GBT	5/19/00	3:23 PM
DDC39CT	5/19/00	3:26 PM
DD6A4T	5/19/00	3:29 PM
DD6A4PT	5/19/00	3:32 PM
DD6A4ST	5/19/00	3:36 PM
DD6A4DT	5/19/00	3·39 PM
DD6A5T	5/19/00	3:42 PM
CCV2-1	5/19/00	3:45 PM
CCB1	5/19/00	3:48 PM
DD6A6T	5/19/00	3:51 PM
DD6A7T	5/19/00	3:55 PM
CCV2-2	5/19/00	3:58 PM
CCB2	5/19/00	4:01 PM
ZZZZZZ	5/19/00	4:06 PM
ZZZZZZ	5/19/00	4:09 PM
ZZZZZZ	5/19/00	4:12 PM
ZZZZZZ	5/19/00	4:15 PM
ZZZZZZ	5/19/00	4:18 PM
ZZZZZZ	5/19/00	4:22 PM
ZZZZZZ	5/19/00	4:25 PM
ZZZZZZ	5/19/00	4:28 PM
ZZZZZZ	5/19/00	4:31 PM
ZZZZZZ	5/19/00	4:34 PM
ZZZZZZ	5/19/00	4:37 PM
ZZZZZZ	5/19/00	4:40 PM
ZZZZZZ	5/19/00	4:44 PM
ZZZZZZ	5/19/00	4:47 PM
ZZZZZZ	5/19/00	4:50 PM

#### TCLP METALS RAW DATA

STL Pittsburgh

657 849 page 3

	muchael Mendorho	5-19-00					
#	Sample Name	AG	AS	BA	CD	CR	PB
1	STD1	001	00049			.0013	.0006
2	STD5A	.25694	.5115	3.88785	.88959	3.51219	.25249
	STD5B					2 0254	1 0440
4	ICV2-1 0014-061-7		1.0143	1.0200	1.0220	1.0174	1.0449
	ICB1	.00075	00799	.00025	00082	00341	01353
6	ICSA 0014-088-12	00032	11064	.00192	00251	00640	.12813
7	ICSAB 0014-043-1	.94437	.87142	.45609	.94991	.43477	1.0398
8	DDC39BT	.00038	00403	.00025	.00107	00128	.00189
	DDA9GBT	.00263	.17865	.03691	00177	00213	01164
10	DDC39CT	.04714	2.2012	1.8741	.04566	.18933	.47228
	DD6A4T	00265	.14949	1.2123	.00241	00427	.02117
	DD6A4TP5	00039		.27021	00049	00384 4.2631	.02501 4.4358
	DD6A4ST	.90355	4.8882	43.256	.94443		4.4898
	DD6A4DT	.91225	4.9588	43.834	.95485	4.3155	.01154
	DD6A5T	00302		1.2211	00302	00584	4.8897
	CCV2-1 0014-087-7	.97794	4.9967	4.7640	4.9809	4.8566 .00000	02124
	CCB1	.00075	.03266	.00082	.00066 .00189	00484	01347
	DD6A6T	00266		1.2934		00555	.00759
	DD6A7T	00077		1.2356	00077 5.0231	4.9023	4.9410
	CCV2-2	.98469	5.0211	4.8021	.00260	00241	00381
	CCB2	.00188	.02176	.00059	00055	00241	01156
	DDDFQBF	.00112	00198	.00025		L00171	
	DDDFQCF	L00037		L.00036	.00007	00171	.00365
	DC7M7F	00056		.09306	00012	00199	
	DC7MAF	00177			.00109	00327	03232
	DC7MAFP5	.00007	00499		00154	00398	.00258
	DC7MASF	00074			.00134	00313	.02382
	DC7MADF	00151			00029	00313	
	DC7MCF	00022		.68327	00226	00469	
	DC7MDF	.00010	.01866 4.9237	4.6919	5.0384	4.8757	
	. CCV2-3	.97412	00406		.00050	00156	
_	CCB3	.00150	.01586	.00037	.00075	00128	00772
	DDCAWB	.00113	1.9425	1.8926	.04496	.19360	.47144
	DDCAWC	.04873 .00113	.01064	.00488	00059		
	DD9PJ	.00000	00896		00175		
	DD9PJP5	.04912	1.9412	1.9181	.04715	.19389	.62546
	7 DD9PJS 3 DD9PJD	.04879	2.0281	1.9491	.04761	.19830	.65662
	DD5V9	.00266	.01827	.02413	.00002	.00142	.04429
	DD9NN DD9NN	.00113	.03108	.00329	.00207	.00128	.72810
	L CCV2-4	.98042	5.0265	4.7074	5.0791	4.9195	4.9365
	CCB4	.00150	.02173	.00069	00068		.01924
	DDCAEB	.00150	00303		00077		
	DDCAEC	.04987	1.9582	1.8438	.04594	.19389	.50239
	DD7WE	.00616	.11347	.08185	.00042	.06296	.03259
	5 DD7WEP5	.00166	.03949	.01636	.00014	.00968	.00146
	DD7WES	.05936	2.0936	1.8400	.04515	.24988	.51465
	DD7WED	.05817	2.0937	1.8105	.04334	.24617	.51863
	9 DD7NT	.00120	.02757	.00239	00030	.06823	.00032
	CCV2-5	.97243	4.9489	4.7232	5.0463	4.8989	4.9188
	L CCB5	.00188	00505		.00084	00070	
	2 DD7WE/2 K	.00231	.02759	.04115	.00102	.03048	.01054
	DD7WEP10 K	.00158	01566	.00874	.00008	.00612	.00167

	- <u> </u>	-		. ,			
#	Sample Name	AG	AS	BA	CD	CR	PB
							•
	DD7WES/2 K	.02968	1.0049	.92912	.02275	.12579	.24564
55	DD7WED/2 K	.02815	1.0190	.92526	.02317	.12322	.25734
56	CCV2-6	.97724	5.0282	4.7509	5.1080	4.9376	4.9445
57	CCB6	.00339	.04355	.00097	~.00083	00013	.00384
58	DDDCMB	.00038	04580	.00038	.00029	00270	01356
59	DDDCMC	.04599	1.8347	1.8753	.04242	.18733	.45749
60	DD5VC	.00465	.09241	.10077	.00149	.11396	.21142
	DD5VE	.13956	.03916	1.5337	.02092	1.5403	.91382
62	DD9NT	.01888	.08065	.21910	.01033	.18276	.55759
	DD9NTP5	.00505	.03770	.04491	.00212	.03590	.10873
	DD9NTS	.06486	1.9511	2.0596	.04787	.36955	.96673
	DD9NTD	.06827	1.9407	2.0672	.05225	.37353	.98456
	DD9NW	.00359	.05674	2.1494	00028	.09913	.37483
	DD9NWP5	.00123	.00381	.42976	00103	.01837	.07262
	CCV2-7	.98883	5.1467	4.8056	5.1936	5.0340	4.9952
	CCB7	.00036	00606	.00064	.00166	00127	01157
	DD9NWS	.04629	1.7822	3.6160	.04038	.27378	.81275
	DD9NWD	.04901	1.9057	3.7783	.04177	.29031	.85556
	DD9PE	.02216	.02932	.17488	.00343	.07777	.08419
	DD9PX	.01182	.09457	.23767	.00377	.41296	.38543
	CCV2-8	.99225	5.1507	4.8167	5.2091	5.0374	5.0436
	CCB8	.00150	.02667	.00051	.00117	.00114	.03275
, –							
#	Sample Name	SE					
		Upro-					
		'  '	WIN 14-00				
1	STD1	.00045	•				
2	STD5A	.503					
3	STD5B						
4	ICV2-1 0014-061-7	1.0340	se				
5	ICB1	<u>00695</u>	ישוטע איירי	The TKO +	ant LC	JEAL LINK	<u>!</u>
6	ICSA 0014-088-12	25834	>C MTW	THE SELECT SE	S OW OF CO	INTROL LIMIT WILL Need	<b>.</b>
7		1.1776	,,,,,	5-14-00 SAM	pres ton se	, WILL NEED	
	DDC39BT	01481	Ke	analyzed m			
	DDA9GBT	03378		m:	5-17-00		
	DDC39CT	2.1189					
	DD6A4T	<b>03</b> 965					
	DD6A4TP5	.00400					
	DD6A4ST	.95 27					
	DD6A4DT	.98310					
	DD6A5T	01684					
	CCV2-1 0014-087-7	5.0441					
	CCB1	01489					
	DD6A6T	00193					
	DD6A7T	03575					
	CCV2-2	5.0 <b>975</b>					
	CCB2	02385					
	DDDFQBF	01490					
	DDDFQCF	ь.00801					
	DC7M7F	00981					
	DC7MAF	.01257					
	DC7MAFP5	.00106					
	DC7MASF	.08425					
28	DC7MADF	.00028					
		X					

# Sample Name	SE VOIDE see previous page for explanation
	1 mru 5-19-00
29 DC7MCF	.02103
30 DC7MDF	01276
31 CCV2-3	4.9863
32 CCB3	0494
33 DDCAWB	04384
34 DDCAWC	1.8783
35 DD9PJ	.01409
36 DD9PJP5	04085
37 DD9PJS	1.9450
38 DD9PJD	1.9481
39 DD5V9	.02199
40 DD9NN	02/265
41 CCV2-4	5.02/33
42 CCB4	.02192
43 DDCAEB	.00599
44 DDCAEC	1.9837
45 DD7WE	00602
46 DD7WEP5	.01150
47 DD7WES	2.0083
48 DD7WED	1.9720
49 DD7NT	01445
50 CCV2-5	5.0276
51 CCB5	03877
52 DD7WE/2 K	.02038
53 DD7WEP10 K 54 DD7WES/2 K	01556 .95343
54 DD7WES/2 K 55 DD7WED/2 K	.94928
56 CCV2-6	4.9660
57 CCB6	00494
58 DDDCMB	00889
59 DDDCMC	1.8294
60 DD5VC	.02195
61 DD5VE	.01929
62 DD9NT	.015 <mark>9</mark> 2
63 DD9NTP5	01 <mark>5</mark> 60
64 DD9NTS	1.9873
65 DD9NTD	1.9452
66 DD9NW	. 02703
67 DD9NWP5	.01118
68 CCV2-7	5.0375
69 CCB7	04275
70 DD9NWS	1.8400
71 DD9NWD	1.8867
72 DD9PE	00 <u>4</u> 77
73 DD9PX 74 CCV2-8	01592 5.0194
74 CCV2-8 75 CCB8	
/5 CCB0	01887

#### MTW 5-19-00

	mrw 5-19-00							
#	Sample Name	File	Method	Date	Time	QΙαO	Type	Mode
1	STD1	J00519A	QUANMET	05/19/00	08:40		X	IR
	STD5A	J00519A	QUANMET	05/19/00	08:43		X	IR
3	STD5B	J00519A	QUANMET	05/19/00	08:47		X	IR
_						MITTER		
4	ICV2-1 0014-061-7	J00519A	QUANMET	05/19/00	08:50		S	CONC
5	ICB1	J00519A	QUANMET	05/19/00	08:53		S	CONC
	ICSA 0014-088-12	J00519A	QUANMET	05/19/00	08:56		Q	CONC
7	ICSAB 0014-043-1	J00519A	QUANMET	05/19/00	08:59		Q	CONC
8	DDC39BT	J00519A	QUANMET	05/19/00	09:05	MTW	S	CONC
9	DDA9GBT	J00519A	QUANMET	05/19/00	09:08	WTM	S	CONC
10	DDC39CT	J00519A	QUANMET	05/19/00	09:11	MTW	S	CONC
11	DD6A4T	J00519A	QUANMET	05/19/00	09:14	MTW	S	CONC
12	DD6A4TP5	J00519A	QUANMET	05/19/00	09:17	MTW	S	CONC
	DD6A4ST	J00519A	QUANMET	05/19/00	09:20		S	CONC
	DD6A4DT	J00519A	QUANMET	05/19/00	09:24		S	CONC
	DD6A5T	J00519A	QUANMET	05/19/00	09:27		S	CONC
	CCV2-1 0014-087-7	J00519A	QUANMET	05/19/00	09:30		ŝ	CONC
	CCB1	J00519A	QUANMET	05/19/00	09:33		Š	CONC
	DD6A6T	J00519A	QUANMET	05/19/00	09:36		S	CONC
				05/19/00	09:39		S	CONC
	DD6A7T	J00519A	QUANMET					
	CCV2-2	J00519A	QUANMET	05/19/00	09:43		S	CONC
	CCB2	J00519A	QUANMET	05/19/00	09:46		s	CONC
	DDDFQBF	J00519A	QUANMET	05/19/00	09:49		S	CONC
	DDDFQCF	J00519A	QUANMET	05/19/00	09:52		S	CONC
	DC7M7F	J00519A	QUANMET	05/19/00	09:55		S	CONC
	DC7MAF	J00519A	QUANMET	05/19/00	09:58		S	CONC
	DC7MAFP5	J00519A	QUANMET	05/19/00	10:02		S	CONC
27	DC7MASF	J00519A	QUANMET	05/19/00	10:05		S	CONC
28	DC7MADF	J00519A	QUANMET	05/19/00	10:08		S	CONC
29	DC7MCF	J00519A	QUANMET	05/19/00	10:11		S	CONC
30	DC7MDF	J00519A	QUANMET	05/19/00	10:14		S	CONC
31	CCV2-3	J00519A	QUANMET	05/19/00	10:17	MTW	S	CONC
32	CCB3	J00519A	QUANMET	05/19/00	10:20	WTM	S	CONC
33	DDCAWB	J00519A	QUANMET	05/19/00	10:23		S	CONC
34	DDCAWC	J00519A	QUANMET	05/19/00	10:27	MTW	S	CONC
35	DD9PJ	J00519A	QUANMET	05/19/00	10:30	MTW	S	CONC
36	DD9PJP5	J00519A	QUANMET	05/19/00	10:33	MTW	S	CONC
37	DD9PJS	J00519A	QUANMET	05/19/00	10:36	MTW	S	CONC
	DD9PJD	J00519A	QUANMET	05/19/00	10:39		S	CONC
	DD5V9	J00519A	QUANMET	05/19/00	10:42		S	CONC
	DD9NN	J00519A	QUANMET	05/19/00	10:46		s	CONC
	CCV2-4	J00519A	QUANMET	05/19/00	10:49		Š	CONC
	CCB4	J00519A	QUANMET	05/19/00	10:52		s	CONC
	DDCAEB	J00519A	QUANMET	05/19/00	10:59		S	CONC
	DDCAEC	J00519A	QUANMET	05/19/00	11:02		S	CONC
	DD7WE	J00519A	QUANMET	05/19/00	11:05		S	CONC
			• •		11:03		S	CONC
	DD7WEP5	J00519A	QUANMET	05/19/00	11.11	MOUNT	S	CONC
	DD7WES	J00519A	QUANMET	05/19/00	11:11			
	DD7WED	J00519A	QUANMET	05/19/00	11:15		S	CONC
	DD7NT	J00519A	QUANMET	05/19/00	11:18		S	CONC
	CCV2-5	J00519A	QUANMET	05/19/00	11:21		S	CONC
	CCB5	J00519A	QUANMET	05/19/00	11:24		S	CONC
	DD7WE/2 K	J00519A	QUANMET	05/19/00	11:32		S	CONC
53	DD7WEP10 K	J00519A	QUANMET	05/19/00	11:35	MTW	S	CONC

STL Pittsburgh 6058

05/19/00 12:56:24 PM

Alla	.yara kepore	Danmary		03/13/00 11				page
#	Sample Name	File	Method	Date	Time	OpID	Туре	Mode
54	DD7WES/2 K	J00519A	QUANMET	05/19/00	11:38	MTW	S	CONC
55	DD7WED/2 K	J00519A	QUANMET	05/19/00	11:41	MTW	S	CONC
56	CCV2-6	J00519A	QUANMET	05/19/00	11:44	$\mathtt{WTM}$	S	CONC
57	CCB6	J00519A	QUANMET	05/19/00	11:48	$\mathbf{W}\mathbf{T}\mathbf{M}$	S	CONC
58	DDDCMB	J00519A	QUANMET	05/19/00	11:54	MTW	S	CONC
59	DDDCMC	J00519A	QUANMET	05/19/00	11:57	MTW	S	CONC
60	DD5VC	J00519A	QUANMET	05/19/00	12:00	MTW	S	CONC
61	DD5VE	J00519A	QUANMET	05/19/00	12:03	MTW	S	CONC
62	DD9NT	J00519A	QUANMET	05/19/00	12:06	$\mathtt{WTM}$	S	CONC
63	DD9NTP5	J00519A	QUANMET	05/19/00	12:10	MTW	S	CONC
64	DD9NTS	J00519A	QUANMET	05/19/00	12:13	MTW	S	CONC
65	DD9NTD	J00519A	QUANMET	05/19/00	12:16	MTW	S	CONC
66	DD9NW	J00519A	QUANMET	05/19/00	12:19		S	CONC
67	DD9NWP5	J00519A	QUANMET	05/19/00	12:22	MTW	S	CONC
68	CCV2-7	J00519A	QUANMET	05/19/00	12:26	MTW	S	CONC
69	CCB7	J00519A	QUANMET	05/19/00	12:29	MTW	S	CONC
70	DD9NWS	J00519A	QUANMET	05/19/00	12:32	MTM	S	CONC
71	DD9NWD	J00519A	QUANMET	05/19/00	12:35		S	CONC
72	DD9PE	J00519A	QUANMET	05/19/00	12:38		S	CONC
73	DD9PX	J00519A	QUANMET	05/19/00	12:42		S	CONC
74	CCV2-8	J00519A	QUANMET	05/19/00	12:45		S	CONC
75	CCB8	J00519A	QUANMET	05/19/00	12:48	MTW	S	CONC

Method:	QUANMET	Standard:	STD1
·	0= /20 /00	00 40 40	

Run T	ime: 05	/19/00	08:40:49
-------	---------	--------	----------

.00035 .00024

#4

	•						
Elem	AG	AL	AS	B	BA	BE	CA
Avge	00100	.01005	00050	.00116	00008	.00295	00138
SDev	.00016	.00066	.00222	.00043	.00006	.00010	.00031
%RSD	16.330	6.5752	444.82	37.150	69.881	3.3898	22.386
#1	00120	.00960	.00160	.00180	00012	.00300	00169
#2	00100	.01100	.00040	.00094	00008	.00280	00156
#3	00080	.01000	00360	.00094	00012	.00300	00125
#4	00100	.00960	00040	.00094	.00000	.00300	00101
Elem	CD	CO	CR	CU	FE	K_	LI
Avge	.00003	.00020	.00130	.00025	.00020	03490	.00086
SDev	.00009	.00023	.00105	.00010	.00016	.00342	.00054
%RSD	303.76	115.47	80.922	40.000	81.650	9.7870	62.445
#1	00008	.00000	.00120	.00040	.00020	03520	.00140
#2	.00000	.00040	.00040	.00020	.00040	03280	.00012
#3	.00008	.00000	.00080	.00020	.00000	03960	.00098
#4	.00012	.00040	.00280	.00020	.00020	03200	.00094
Elem	MG	MN	MO	NA	NI	PB	SB
Avge	.00035	.00025	.00005	.05555	.00094	.00060	00010
SDev	.00087	.00010	.00019	.00101	.00076	.00046	.00053
%RSD	248.53	40.000	382.97	1.8211	80.742	76.980	529.15
#1	00040	.00020	.00000	.05680	.00154	.00020	.00040
#2	.00160	.00020	.00020	.05440	.00164	.00020	.00020
#3	.00000	.00020	00020	.05580	.00026	.00100	00080
#4	.00020	.00040	.00020	.05520	.00031	.00100	00020
Elem	SE	SI	SN	SR	TI	TL	V
Avge	.00045	.00050	.00070	00005	.00210	.00000	.00040
SDev	.00222	.00020	.00202	.00010	.00020	.00036	.00046
%RSD	494.08	40.000	288.09	200.00	9.5238	.00000	115.47
#1	00260	.00040	.00280	.00000	.00180	00040	.00000
#2	.00020	.00040	00140	.00000	.00220	00020	.00080
#3	.00220	.00040	.00200	.00000	.00220	.00020	.00000
#4	.00200	.00080	00060	00020	.00220	.00040	.00080
Elem Avge SDev %RSD	ZN .00029 .00015 51.640						
#1 #2 #3	.00012 .00047 .00035						

6060 STL Pittsburgh

Method: QUANMET Standard: STD5A 0014-072-2 Run Time: 05/19/00 08:43:59

Elem	AG	AS	B_	BA	BE	CD	CO
Avge	.25695	.51150	.55832	3.8879	14.163	.88960	1.3870
SDev	.00082	.00527	.00210	.0190	.072	.00124	.0047
%RSD	.32014	1.0298	.37583	.48775	.50536	.13899	.34150
#1	.25680	.50880	.55780	3.8809	14.150	.88912	1.3878
#2	.25700	.50900	.55956	3.9024	14.200	.88908	1.3884
#3	.25800	.51940	.56032	3.9040	14.234	.89144	1.3916
#4	.25600	.50880	.55560	3.8642	14.069	.88875	1.3804
Elem	CR	CU	LI	MN	MO	NI	PB
Avge	3.5122	2.1644	4.5413	1.9358	.40270	1.6493	.25250
SDev	.0161	.0098	.0126	.0094	.00167	.0053	.00035
%RSD	.45792	.45306	.27677	.48548	.41454	.32253	.13719
#1	3.5198	2.1586	4.5304	1.9390	.40240	1.6499	.25220
#2	3.5146	2.1706	4.5518	1.9382	.40240	1.6489	.25280
#3	3.5254	2.1746	4.5525	1.9438	.40500	1.6556	.25280
#4	3.4890	2.1538	4.5304	1.9222	.40100	1.6426	.25220
			O.T.	an.	an		mr
Elem	SB	SE	SI	SN	SR	TI	TL
Avge	.18995	.50300	.32920	.96635	7.4043	5.7691	.34705
SDev	.00103	.00629	.00356	.00975	.0360	.0235	.00567
%RSD	.53945	1.2502	1.0800	1.0086	.48592	.40818	1.6350
Avge	.18995	.50300	.32920	.96635	7.4043	5.7691	.34705
SDev	.00103	.00629	.00356	.00975	.0360	.0235	.00567
Avge	.18995	.50300	.32920	.96635	7.4043	5.7691	.34705
SDev	.00103	.00629	.00356	.00975	.0360	.0235	.00567
%RSD	.53945	1.2502	1.0800	1.0086	.48592	.40818	1.6350
#1	.19040	.50100	.32420	.96520	7.3922	5.7666	.34240
#2	.18980	.50960	.33020	.97400	7.4301	5.7800	.35080
#3	.19100	.50620	.33260	.97320	7.4362	5.7922	.35300

STL Pittsburgh 6061

05/19/00 08:50:13 AM page 1

Method: QUANMET Standard: STD5B 0014-072-3 Run Time: 05/19/00 08:47:08

Elem	AL	CA	FE	K_	MG	NA
Avge	10.921	23.167	28.472	2.4135	7.2390	47.165
SDev	.013	.077	.051	.0233	.0125	.136
%RSD	.12282	.33241	.17992	.96439	.17232	.28904
#1	10.931	23.257	28.539	2.4236	7.2544	47.178
#2	10.902	23.078	28.414	2.3830	7.2266	46.984
#3	10.929	23.135	28.470	2.4100	7.2436	47.180
#4	10.923	23.196	28.467	2.4376	7.2316	47.316

05/19/00 08:50:14 AM

Element	Wavelen	High std	Low std	Slope	Y-intercept	Date Standardized
AG	328.068	STD5A	STD1	7.56474	.007565	05/19/00 08:47:08
AL	308.215	STD5B	STD1	9.16502	092108	05/19/00 08:47:08
AS	193.696	STD5A	STD1	19.8680	.009934	05/19/00 08:47:08
В	249.600	STD5A	STD1	17.9841	020787	05/19/00 08:47:08
BĀ	493.409	STD5A	STD1	2.57206	.000204	05/19/00 08:47:08
BE	313.042	STD5A	STD1	.708081	002089	05/19/00 08:47:08
CA	317.933	STD5B	STD1	4.31630	.005954	05/19/00 08:47:08
CD	228.802	STD5A	STD1	11.4098	000336	05/19/00 08:47:08
CO	228.616	STD5A	STD1	7.24502	001449	05/19/00 08:47:08
CR	267.716	STD5A	STD1	2.84906	003704	05/19/00 08:47:08
CU	324.754	STD5A	STD1	4.61654	001154	05/19/00 08:47:08
FE	259.940	STD5B	STD1	3.51221	000702	05/19/00 08:47:08
ĸ	766.491	STD5B	STD1	40.8422	1.42539	05/19/00 08:47:08
${f r}_{f I}$	670.789	STD5A	STD1	2.20244	001892	05/19/00 08:47:08
MG	279.079	STD5B	STD1	13.8146	004835	05/19/00 08:47:08
MN	257.610	STD5A	STD1	5.16479	001291	05/19/00 08:47:08
MO	202.030	STD5A	STD1	24.8355	001242	05/19/00 08:47:08
NA	588.995	STD5B	STD1	2.12273	117918	05/19/00 08:47:08
NI	231.604	STD5A	STD1	6.06401	005672	05/19/00 08:47:08
PB	220.353	STD5A	STD1	38.5304	023118	05/19/00 08:47:08
SB	206.838	STD5A	STD1	52,6787	.005268	05/19/00 08:47:08
SE	196.026	STD5A	STD1	19.8985	008954	05/19/00 08:47:08
SI	288.158	STD5A	STD1	30.0346	015017	05/19/00 08:47:08
SN	189.989	STD5A	STD1	10.3449	007241	05/19/00 08:47:08
SR	409.552	STD5A	STD1	1.35055	.000068	05/19/00 08:47:08
$\mathtt{TI}$	334.941	STD5A	STD1	1.73402	003641	05/19/00 08:47:08
$\mathtt{TL}$	190.864	STD5A	STD1	57.8987	.000000	05/19/00 08:47:08
<b>V</b>	292.402	STD5A	STD1	6.44478	002578	05/19/00 08:47:08
$z\overline{N}$	213.856	STD5A	STD1	6.73386	001981	05/19/00 08:47:08

page 1

05/19/00 08:53:22 AM

Method: QUANMET Sample Name: ICV2-1 0014-061-7 Operator: MTW

Run Time: 05/19/00 08:50:17

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

•	icac. co.							
	Elem	AG	AL	AS	B_	BA	BE	CA
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	.51087	25.708	1.0143	1.0200	1.0200	.99958	25.697
	SDev	.00147	.113	.0427	.0018	.0041	.00260	.048
	%RSD	.28795	.44140	4.2126	.17906	.40521	.26039	.18524
	#1	.51201	25.599	.97944	1.0180	1.0170	.99944	25.724
	#2	.51052	25.826	.97740	1.0188	1.0249	1.0027	25.735
	#3	.50893	25.623	1.0629	1.0214	1.0162	.99634	25.629
	#4	.51202	25.783	1.0375	1.0216	1.0218	.99986	25.700
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	.55000	27.500	1.1000	1.1000	1.1000	1.1000	27.500
	Low	.45000	22.500	.90000	.90000	.90000	.90000	22.500
	Elem	CD	CO	CR	CU	FE	K_	LI
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	1.0220	1.0105	1.0174	1.0191	26.474	25.984	1.0438
	SDev	.0039	.0025	.0010	.0043	.090	.446	.0124
	%RSD	.38627	.24421	.09702	.42146	.34075	1.7183	1.1918
	#1	1.0220	1.0080	1.0171	1.0138	26.463	25.367	1.0272
	#2	1.0207	1.0095	1.0188	1.0231	26.573	26.429	1.0497
	#3	1.0179	1.0109	1.0165	1.0175	26.357	26.021	1.0420
	#4	1.0273	1.0138	1.0171	1.0221	26.501	26.119	1.0561
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	1.1000	1.1000	1.1000	1.1000	27.500	27.500	1.1000
	Low	.90000	.90000	.90000	.90000	22.500	22.500	.90000
	Elem	MG	MN	MO	NA	NI	PB	SB
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	25.938	1.0159	.99614	26.232	1.0114	1.0449	1.0867
	SDev	.096	.0018	.00407	.294	.0132	.0438	.0133
	%RSD	.36967	.17802	.40846	1.1197	1.3036	4.1894	1.2230
	#1	25.853	1.0174	.99613	25.831	1.0039	1.0892	1.0892
	#2	26.027	1.0164	1.0011	26.468	1.0122	1.0197	1.0998
	#3	25.856	1.0132	.99115	26.196	.99986	1.0738	1.0681
	#4	26.014	1.0164	.99614	26.434	1.0296	.99676	1.0896
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	27.500	1.1000	1.1000	27.500	1.1000	1.1000	1.1000
	Low	22.500	.90000	.90000	22.500	.90000	.90000	.90000
	Elem	SE	SI	SN	SR	TI	TL	V_
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	1.0340	1.0070	1.0146	1.0152	1.0052	4.9860	1.0073
	SDev	.0369	.0290	.0266	.0045	.0031	.0997	.0026
	%RSD	3.5710	2.8819	2.6186	.44403	.31242	1.9989	.26245
	#1	1.0071	.99500	1.0358	1.0107	1.0038	5.0093	1.0086
	#2	1.0553	.98299	1.0255	1.0200	1.0090	4.8808	1.0087

Analysis	Report			05/19,	/00 08:53:	22 AM	657 859 page 2
#3 #4	.99875 1.0749	1.0009 1.0491	1.0212 .97579	1.0121 1.0180	1.0017 1.0063	4.9409 5.1131	1.0033 1.0086
Errors High Low	LC Pass 1.1000 .90000	LC Pass 1.1000 .90000	LC Pass 1.1000 .90000	LC Pass 1.1000 .90000	LC Pass 1.1000 .90000	LC Pass 5.5000 4.5000	LC Pass 1.1000 .90000
Elem Units Avge SDev %RSD	ZN ppm 1.0152 .0047 .46630						
#1 #2 #3 #4	1.0127 1.0196 1.0188 1.0098						
Errors High Low	LC Pass 1.1000 .90000						

Analysis Report

05/19/00 08:56:30 AM page 1

Operator: MTW

Method: QUANMET Sample Name: ICB1 Run Time: 05/19/00 08:53:25

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00076	.00496	00799	.01276	.00026	.00007	.00001
SDev	.00196	.00685	.05448	.00746	.00034	.00007	.00220
%RSD	258.64	137.87	681.43	58.476	131.84	106.95	17339.
#1	.00001	.00484	08552	.01519	.00020	.00003	.00022
#2	00152	.00308	.04174	.01878	.00020	.00018	00303
#3	.00304	.01418	.00182	.00186	00010	.00003	.00067
#4	.00150	00225	.00999	.01519	.00072	.00004	.00220
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00083	00145	00342	00069	.00299	12661	00341
SDev	.00224	.00167	.00136	.00053	.00327	.02750	.00194
%RSD	270.98	115.50	39.678	76.849	109.33	21.719	56.862
#1	.00061	00290	00256	00116	.00070	11027	00334
#2	00310	00289	00484	00023	00000	16745	00352
#3	.00152	00000	00428	00023	.00421	11844	00101
#4	00233	.00000	00199	00115	.00703	11027	00575
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00829	.00026	.00621	.00297	00355	01353	00270
SDev	.01018	.00059	.00860	.00658	.00427	.01927	.02337
%RSD	122.85	227.69	138.55	221.43	120.25	142.40	866.52
#1	.00345	00025	.01366	.00138	00046	02320	.00515
#2	02141	00025	.01366	00541	00010	02317	.02631
#3	00760	.00078	00124	.00987	00439	02315	01578
#4	00760	.00078	00124	.00605	00925	.01538	02646
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00695	.00300	.00775	.00012	.00017	00581	.00011
SDev	.02017	.00694	.01901	.00006	.00049	.02756	.00298
%RSD	290.05	230.90	245.43	52.148	282.84	474.47	2633.9
#1	00099	.00905	.02589	.00018	.00017	.02315	.00284
#2	02089	.00898	01332	.00007	00017	02312	00232

Analysis	Report			05/19,	/00 08:56:	30 AM	657 861 page 2
#3 #4	02486 .01893	00297 00304	00318 .02160	.00007 .00018	00017 .00087	.01152 03479	.00255 00261
Errors High Low	LC Pass .25000 25000	LC Pass .50000 50000	LC Pass .10000 10000	LC Pass .05000 05000	LC Pass .05000 05000	LC Pass .30000 30000	LC Pass .05000 05000
Elem Units Avge SDev %RSD	ZN ppm .00043 .00129 298.35						
#1 #2 #3 #4	.00232 00041 00037 .00018						
Errors High Low	LC Pass .02000 02000						

Analysis Report QC Standard 05/19/00 08:59:38 AM

page 1

Method: QUANMET Sample Name: ICSA 0014-088-12 Operator: MTW

Run Time: 05/19/00 08:56:33

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

ī,	ioue: coi	NC COLL.	raccor. I					
	Elem	AG	AL	AS	B_	BA	BE	CA
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	00032	484.33	11064	06151	.00193	00008	475.65
	SDev	.00191	.44	.07303	.00257	.00024	.00007	.89
	%RSD	590.92	.09133	66.001	4.1840	12.614	90.534	.18783
	#1	00220	484.50	13200	06088	.00226	00011	474.79
	#2	00073	484.58	10068	06118	.00196	00011	475.02
	#3	.00234	483.67	19263	06507	.00175	00011	476.13
	#4	00071	484.57	01726	05893	.00175	.00003	476.66
	Errors Value Range	NOCHECK	QC Pass 500.00 20.000	NOCHECK	NOCHECK	NOCHECK	NOCHECK	QC Pass 500.00 20.000
	Elem Units Avge SDev %RSD	CD ppm 00252 .00510 202.72	CO ppm .00874 .00264 30.202	CR ppm 00641 .00544 84.839	CU ppm 00676 .00088 13.043	FE ppm 180.90 .08 .04426	K_ppm23893 .12049 50.430	LI ppm 00189 .00000
	#1	00176	.00584	01054	00791	180.95	38800	00189
	#2	00571	.01020	00427	00607	180.80	22463	00189
	#3	00698	.00730	.00029	00698	180.97	09394	00189
	#4	.00437	.01162	01111	00606	180.87	24914	00189
	Errors Value Range	NOCHECK	NOCHECK	NOCHECK	NOCHECK	QC Pass 200.00 20.000	NOCHECK	NOCHECK
	Elem Units Avge SDev %RSD	MG ppm 479.64 .35	MN ppm .00264 .00053 19.975	MO ppm .01587 .01073 67.612	NA ppm .00732 .00173 23.667	NI ppm 02147 .00916 42.670	PB ppm .12813 .06087 47.505	SB ppm .00531 .08259 1556.5
	#1	480.01	.00290	.00595	.00860	02935	.05862	.02624
	#2	479.60	.00286	.01586	.00902	01413	.09713	07895
	#3	479.76	.00293	.03079	.00563	02944	.19002	03694
	#4	479.18	.00185	.01090	.00605	01297	.16676	.11087
	Errors Value Range	QC Pass 500.00 20.000	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK
	Elem	SE	SI	SN	SR	TI	TL	V_
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	.25834	.08112	08293	.01326	02358	.47109	.00287
	SDev	.09659	.00000	.04883	.00016	.00020	.13384	.00022
	%RSD	37.386	.00342	58.882	1.1758	.84904	28.411	7.5724
	#1	.36997	.08112	02024	.01313	02341	.47057	.00267
	#2	.25400	.08112	07248	.01340	02376	.53003	.00288

Analysis	Report QC Standard		05/19/00 08:59:38 AM			657 863 page 2	
#3 #4	.27453 .13488	.08113 .08112	13432 10469	.01340 .01313	02376 02341	.59772 .28607	.00317 .00277
Errors Value Range	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK
Elem Units Avge SDev %RSD	ZN ppm .00172 .00123 71.477						
#1 #2 #3 #4	.00158 .00106 .00348 .00074		·				
Errors Value Range	NOCHECK						

Analysis Report QC Standard 05/19/00 09:02:47 AM page 1

Method: QUANMET Sample Name: ICSAB 0014-043-1 Operator: MTW

Run Time: 05/19/00 08:59:41

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.94438	483.97	.87143	.86952	.45609	.43777	475.53
SDev	.00373	2.15	.02787	.02645	.00232	.00165	1.79
%RSD	.39501	.44493	3.1985	3.0417	.50912	.37669	.37651
#1	.93923	481.91	.88884	.83185	.45364	.43558	473.13
#2	.94711	485.66	.83992	.89377	.45797	.43897	475.32
#3	.94715	485.99	.85685	.87552	.45818	.43910	477.28
#4	.94401	482.31	.90009	.87696	.45458	.43742	476.39
Errors	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass
Value	1.0000	500.00	1.0000	1.0000	.50000	.50000	500.00
Range	20.000	20.000	20.000	20.000	20.000	20.000	20.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.94991	.44757	.43477	.47189	180.61	9.5407	.94304
SDev	.00509	.00138	.00569	.00319	.63	.1726	.00675
%RSD	.53631	.30777	1.3081	.67594	.34888	1.8088	.71610
#1	.94295	.44651	.42708	.46790	179.69	9.3161	.94254
#2	.95090	.44648	.43392	.47446	180.98	9.4958	.94635
#3	.95059	.44793	.43962	.47447	181.06	9.6592	.94943
#4	.95520	.44937	.43848	.47075	180.71	9.6918	.93383
Errors	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass
Value	1.0000	.50000	.50000	.50000	200.00	10.000	1.0000
Range	20.000	20.000	20.000	20.000	20.000	20.000	20.000
Elem Units Avge SDev %RSD	MG ppm 477.79 1.77	MN ppm .45142 .00294 .65162	MO ppm .90618 .00473 .52163	NA ppm 9.6803 .0806 .83287	NI ppm .86395 .01555 1.7994	PB ppm 1.0398 .0136 1.3089	SB ppm .95797 .00604 .63080
#1	475.59	.44706	.90977	9.6216	.86098	1.0294	.95277
#2	479.00	.45252	.90500	9.7367	.86127	1.0426	.96327
#3	479.44	.45254	.90998	9.7617	.88538	1.0579	.95270
#4	477.13	.45354	.89999	9.6012	.84815	1.0291	.96313
Errors	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass
Value	500.00	.50000	1.0000	10.000	1.0000	1.0000	1.0000
Range	20.000	20.000	20.000	20.000	20.000	20.000	20.000
Elem Units Avge SDev %RSD	SE ppm 1.1776 .1268 10.765	SI ppm 1.0480 .0000	SN ppm .80581 .04225 5.2432	SR ppm .91091 .00394 .43243	TI ppm .86987 .00368 .42365	TL ppm 9.2040 .2947 3.2015	V_ ppm .45564 .00249 .54756
#1	1.0180	1.0480	Q.78132	.90686	.86545	8.8284	.45409
#2	1.1859	1.0480	Q.77952	.91415	.87308	9.5099	.45530

Analysis	Report	QC Sta	ndard	05/19	/00 09:02:	47 AM	657 865 page 2
inaliana nopolo		20 000		00, 20,	00 03,02,		page 2
#3 #4	1.1782 Q1.3282	1.0481 1.0480	.86845 Q.79397	.91442 .90821	.87273 .86822	9.3469 9.1306	.45927 .45392
Errors	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass	QC Pass
Value Range	1.0000 20.000	1.0000	1.0000 20.000	1.0000 20.000	1.0000 20.000	10.000	.50000 20.000
Elem Units Avge SDev %RSD	ZN ppm .92490 .00270 .29242		ι				
#1 #2 #3 #4	.92392 .92757 .92156 .92654						
Errors Value Range	QC Pass 1.0000 20.000						

05/19/00 09:08:10 AM

Operator: MTW

page 1

Method: QUANMET Sample Name: DDC39BT Operate Run Time: 05/19/00 09:05:04 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00039	.00273	00404	00417	.00026	00007	.01925
SDev	.00228	.00739	.04554	.01512	.00034	.00008	.00164
%RSD	590.99	270.70	1128.2	362.54	131.84	106.90	8.5290
#1	00303	00779	02576	00169	00010	00010	.01733
#2	.00153	.00679	.02966	.01524	.00020	00011	.01911
#3	.00151	.00877	05769	00952	.00072	.00004	.02134
#4	.00153	.00314	.03764	02072	.00020	00011	.01921
Errors High Low	LC Pass .50000 50000	NOCHECK	LC Pass .50000 50000	NOCHECK	LC Pass 10.000 -10.000	NOCHECK	NOCHECK
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00108	00145	00128	.00000	.03056	12048	00001
SDev	.00271	.00205	.00150	.00138	.00167	.19789	.00127
%RSD	251.48	141.45	116.81	123380.	5.4788	164.25	19668.
#1	.00450	00291	00199	00115	.02809	18379	00189
#2	.00161	00145	00256	00115	.03161	36350	.00036
#3	00196	.00145	00142	.00070	.03161	02859	.00088
#4	.00017	00290	.00085	.00162	.03091	.09394	.00062
Errors High Low	LC Pass .10000 10000	NOCHECK	LC Pass .50000 50000	NOCHECK	NOCHECK	NOCHECK	NOCHECK
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00691	00025	.00000	.01242	00349	.00190	.01317
SDev	.00913	.00084	.00248	.00400	.00261	.01826	.01582
%RSD	132.16	338.32	54356.	32.192	74.884	963.40	120.19
#1	01865	00025	00124	.00902	00459	00005	.02635
#2	00760	.00078	.00373	.00945	00667	.01537	.00525
#3	00484	00128	00124	.01751	00124	.01547	.02635
#4	.00345	00025	00124	.01369	00146	02321	00529
Errors High Low	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK	LC Pass .50000 50000	NOCHECK
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	01481	.00451	01028	.00003	00009	.00838	00000
SDev	.03640	.00903	.01172	.00016	.00059	.03046	.00301
%RSD	245.75	200.49	114.06	583.68	683.13	363.48	122760.
#1	02079	00904	02366	00020	00087	.01133	00261
#2	02078	.00904	01549	.00007	.00017	.03441	.00264

page 2

05/19/00 09:08:10 AM

Analysis Report

-.03507 -.00260 .00898 -.00505 .00007 .00052 #3 -.05262 .00309 .02284 .00256 #4 .03494 .00904 .00018 -.00017

Errors LC Pass NOCHECK NOCHECK NOCHECK NOCHECK NOCHECK

High .25000 Low -.25000

Elem ZN Units ppm Avge .00260 SDev .00097 %RSD 37.139

#1 .00341 #2 .00210 #3 .00149 #4 .00341

Errors NOCHECK

High Low

Operator: MTW

Analysis Report

Method: QUANMET Sample Name: DDA9GBT Run Time: 05/19/00 09:08:12

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode. co.	iic corr.	200000					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00264	.02150	.17866	.02727	.03692	00017	.38198
SDev	.00146	.00713	.05125	.01099	.00047	.00007	.00722
%RSD	55.286	33.173	28.688	40.299	1.2740	43.521	1.8892
#1	.00148	.01975	.22844	.03678	.03642	00023	.37203
#2	.00150	.01238	.11715	.01775	.03715	00024	.38137
#3	.00304	.02505	.15669	.03679	.03745	00011	.38655
#4	.00453	.02883	.21237	.01776	.03664	00010	.38797
Errors High Low	LC Pass .50000 50000	NOCHECK	LC Pass .50000 50000	NOCHECK	LC Pass 10.000 -10.000	NOCHECK	NOCHECK
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00177	00255	00213	.00115	.01229	.24301	00125
SDev	.00274	.00073	.00134	.00053	.00240	.27768	.00057
%RSD	154.25	28.640	63.016	46.217	19.517	114.27	45.796
#1	00518	00145	00313	.00069	.00984	06126	00145
#2	00166	00291	00199	.00069	.01264	.08577	00197
#3	.00152	00292	00028	.00161	.01124	.53503	00083
#4	00178	00291	00313	.00161	.01546	.41251	00075
Errors High Low	LC Pass .10000 10000	NOCHECK	LC Pass .50000 50000	NOCHECK	NOCHECK	NOCHECK	NOCHECK
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.05319	.00026	.00373	53373.6	.00023	01164	.00261
SDev	.01296	.00060	.00406	.2	.00782	.02037	.02337
%RSD	24.366	227.14	108.81	.00736	3353.6	174.95	896.34
#1	.05871	.00078	.00373	\$3374.0	00325	02318	00535
#2	.03385	00026	00124	\$3373.5	00668	.01529	02635
#3	.05871	00025	.00869	\$3373.4	00045	00777	.02630
#4	.06148	.00078	.00373	\$3373.5	.01131	03090	.01584
Errors High Low	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK	LC Pass .50000 50000	NOCHECK
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	03378	.09758	.00208	.00250	.00026	02614	00251
SDev	.04012	.00300	.01829	.00031	.00033	.04948	.00425
%RSD	118.76	3.0793	878.15	12.483	127.66	189.27	169.38
#1	06861	.09902	.01135	.00223	.00052	.01155	00767
#2	01687	.09307	01978		00017	09273	00261

#3 #4	06463 .01498	.09915 .09908	00505 .02180	.00277 .00277	.00017 .00052	03485 .01147	.00275 00251
Errors High Low	LC Pass .25000 25000	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK	NOCHECK
Elem Units Avge SDev %RSD	ZN ppm .01293 .00090 6.9569						
#1 #2 #3 #4	.01170 .01284 .01363 .01357						
Errors High Low	NOCHECK						

Operator: MTW

Method: QUANMET Sample Name: DDC39CT

Run Time: 05/19/00 09:11:20

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04714	1.8721	2.2012	.95340	1.8741	.04610	L.40736
SDev	.00151	.0073	.0697	.01106	.0081	.00016	.00439
%RSD	3.2098	.38756	3.1662	1.1601	.43239	.33800	1.0775
#1	.04636	1.8661	2.1665	.95790	1.8634	.04597	L.40275
#2	.04941	1.8808	2.2776	.93843	1.8777	.04624	L.41326
#3	.04641	1.8753	2.1226	.96446	1.8823	.04623	L.40605
#4	.04639	1.8661	2.2380	.95282	1.8728	.04596	L.40737
Errors	LC Pass	LC Pass	LC Pass	NOCHECK	LC Pass	LC Pass	LC Low
High	.06000	2.4000	2.4000		2.4000	.06000	60.000
Low	.04000	1.6000	1.6000		1.6000	.04000	40.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04566	.47131	.18934	.23526	1.0068	L.42680	.88747
SDev	.00108	.00426	.00333	.00177	.0040	.24654	.00410
%RSD	2.3562	.90381	1.7600	.75320	.39627	57.763	.46161
#1	.04496	.46842	.18578	.23387	1.0038	L.34716	.89067
#2	.04549	.47420	.19091	.23573	1.0073	L.78825	.88556
#3	.04723	.47565	.19318	.23757	1.0122	L.23280	.89106
#4	.04496	.46696	.18749	.23387	1.0038	L.33899	.88261
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Low	NOCHECK
High	.06000	.60000	.24000	.30000	1.2000	60.000	
Low	.04000	.40000	.16000	.20000	.80000	40.000	
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	L.06355	.47309	.00139	S3373.6	.45987	.47228	L.01147
SDev	.00761	.00223	.00287	.3	.00593	.01156	.01926
%RSD	11.973	.47174	205.95	.00832	1.2889	2.4473	167.89
#1	L.06700	.46999	.00388	S3373.8	.45240	.48181	L00970
#2	L.07253	.47412	.00388	S3373.6	.46690	.45888	L.02206
#3	L.05595	.47516	00109	S3373.3	.46034	.48206	L.03253
#4	L.05871	.47309	00109	S3373.9	.45983	.46637	L.00100
Errors	LC Low	LC Pass	NOCHECK	LC High	LC Pass	LC Pass	LC Low
High	60.000	.60000		60.000	.60000	.60000	.60000
Low	40.000	.40000		40.000	.40000	.40000	.40000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	2.1189	.08855	.01628	.92346	.00026	1.8177	.46700
SDev	.0693	.00296	.01771	.00468	.00044	.1087	.00418
%RSD	3.2692	3.3427	108.77	.50719	167.77	5.9826	.89438
#1	2.0164	.09299	.02394	.91806	.00017	1.8149	.46189
#2	2.1517	.08705	.01374	.92481	.00087	1.9538	.46705

.40000

Low

Method: QUANMET Sample Name: DD6A4T Operator: MTW

Run Time: 05/19/00 09:14:32

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00265	.05268	.14950	.14776	1.2123	00004	H733.71
SDev	.00189	.00543	.05011	.00478	.0167	.00008	3.96
%RSD	71.361	10.316	33.519	3.2359	1.3806	202.56	.53929
#1	00453	.04706	.09686	.15194	1.2305	.00003	H737.12
#2	00302	.05636	.11666	.14749	1.2210	.00003	H735.14
#3	00002	.05821	.19620	.14114	1.1929	00010	H728.00
#4	00303	.04908	.18827	.15045	1.2047	00011	H734.57
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR ppm00427 .00192 44.924	CU	FE	K_	LI
Units	ppm	ppm		ppm	ppm	ppm	ppm
Avge	.00242	00054		.00577	.03848	1.4356	00141
SDev	.00380	.00291		.00092	.00100	.1932	.00039
%RSD	157.27	534.35		15.978	2.5885	13.460	27.645
#1 #2 #3 #4	.00674 .00198 .00338	00491 .00090 .00091 .00092	00484 00371 00199 00656	.00623 .00439 .00624 .00624	.03847 .03777 .03989 .03777	1.3764 1.3600 1.7195 1.2865	00101 00153 00189 00119
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	8.9443	.09143	.00001	S3374.8	00341	.02117	.02637
SDev	.0851	.00099	.00476	.2	.01080	.02466	.02582
%RSD	.95144	1.0825	82656.	.00504	316.50	116.49	97.898
#1	8.9968	.09272	.00373	\$3374.7	01150	.03842	.04734
#2	8.9857	.09168	00124	\$3374.6	01324	01544	00529
#3	8.8172	.09065	.00373	\$3374.9	.00926	.03087	.04754
#4	8.9774	.09065	00620	\$3374.9	.00183	.03083	.01590
Errors	LC Pass	LC Pass	LC Pass	LC High	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	03966	2.7121	.02099	2.1798	04552	.00012	.00096
SDev	.07434	.0416	.01436	.0284	.00091	.07852	.00241
%RSD	187.45	1.5354	68.429	1.3011	2.0038	68133.	251.43
#1	10831	2.7482	.02212	2.2117	04630	09249	.00264
#2	03667	2.7482	.01360	2.1928	04595	.04642	.00255

Analysis	Report			05/19,	/00 09:17:	38 AM	657 873 page 2
#3 #4	07646 .06282	2.6761 2.6761	.00763 .04059	2.1467 2.1683	04422 04560	03463 .08116	00251 .00116
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .04088 .00204 4.9964						
#1 #2 #3 #4	.04169 .04218 .04183 .03783						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW

Method: QUANMET Sample Name: DD6A4TP5

Run Time: 05/19/00 09:17:45

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode. Co.	no corr.	10000					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00039	.01010	.03472	.00208	.27022	.00004	155.98
SDev	.00145	.00879	.02494	.00170	.00159	.00000	.42
%RSD	368.74	87.006	71.832	81.655	.58933	.31646	.26992
#1	.00150	.02158	.05349	.00229	.26803	.00004	156.17
#2	00001	.00321	.02186	.00442	.27163	.00004	156.39
#3	00153	.01237	.05754	.00082	.27009	.00004	155.93
#4	00153	.00323	.00597	.00081	.27112	.00004	155.41
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00050	00121	~.00384	.00115	.01015	.18379	00134
SDev	.00285	.00275	.00157	.00053	.00070	.38110	.00050
%RSD	574.68	227.84	40.827	46.302	6.9288	207.35	37.515
#1	00321	.00279	00199	.00162	.01051	.56771	00083
#2	.00169	00302	00541	.00161	.01050	32265	00163
#3	.00223	00303	00313	.00069	.01050	.35533	00101
#4	00270	00156	00484	.00069	.00910	.13478	00189
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.9610	.01989	.00000	244.95	00392	.02502	.01318
SDev	.0058	.00060	.00248	2.18	.00905	.00382	.00520
%RSD	.29329	2.9966	163590.	.89030	231.02	15.278	39.475
#1	1.9624	.01937	.00373	242.27	00474	.02320	.01577
#2	1.9541	.02040	00124	245.09	00182	.02303	.00538
#3	1.9679	.01937	00124	244.82	.00639	.03075	.01587
#4	1.9596	.02040	00124	247.61	01551	.02309	.01569
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00601	.56161	00404	.48577	00945	.02898	-:00259
SDev	.04657	.00000	.02033	.00187	.00052	.06585	.00005
%RSD	775.28	.00000	503.18	.38586	5.5046	227.19	1.9516
#1	.07466	.56161	.00114	.48315	00884	00000	00251
#2	01290	.56161	03408	.48720	00988	01153	00261

Analysis	Report			657 875 05/19/00 09:20:50 AM page 2				
#3 #4	00892 02882	.56161 .56161	.00942	.48569 .48704	00919 00988	.12742 .00005	00261 00261	
Errors H <b>igh</b> Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000	
Elem Units Avge SDev %RSD	ZN ppm .00978 .00039 3.9996							
#1 #2 #3 #4	.00958 .01036 .00955 .00962							
Errors High Low	LC Pass 100.00 02000							

657 876 Analysis Report

05/19/00 09:23:58 AM page 1

Operator: MTW Sample Name: DD6A4ST Method: QUANMET

Run Time: 05/19/00 09:20:52

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.90356	.05542	4.8882	.13720	43.256	00004	H690.35
SDev	.00584	.00608	.1085	.00174	.579	.00008	5.31
%RSD	.64665	10.970	2.2188	1.2699	1.3392	218.73	.76959
#1	.89713	.05638	4.7302	.13896	42.674	00011	H686.28
#2	.90016	.05635	4.9726	.13479	43.038	00011	H685.42
#3	.90923	.04716	4.9409	.13748	44.044	.00003	H696.17
#4	.90772	.06181	4.9090	.13754	43.267	.00003	H693.51
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.94444	00110	4.2631	.00693	.03234	1.0251	00176
SDev	.00897	.00206	.0372	.00332	.00239	.1274	.00027
%RSD	.94993	186.65	.87221	47.912	7.3993	12.430	15.498
#1	.94984	.00062	4.2306	.00531	.03147	1.0333	00189
#2	.93705	00098	4.2346	.00439	.03568	.84543	00135
#3	.93661	00003	4.3081	.01177	.03005	1.0823	00189
#4	.95426	00403	4.2790	.00623	.03217	1.1395	00189
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	8.5347	.08626	.00000	S3375.4	.00465	4.4358	.03360
SDev	.0915	.00099	.00248	.2	.00918	.0611	.02606
%RSD	1.0722	1.1463	51347.	.00530	197.34	1.3770	77.558
#1	8.4580	.08548	00124	\$3375.6	.00627	4.3569	.05498
#2	8.4884	.08549	00124	\$3375.5	00065	4.4185	.05483
#3	8.6652	.08652	.00373	\$3375.3	.01687	4.4802	.00178
#4	8.5271	.08755	00124	\$3375.2	00388	4.4878	.02281
Errors	LC Pass	LC Pass	LC Pass	LC High	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.95027	2.5724	00693	2.0821	04188	02300	.00070
SDev	.04199	.0114	.01968	.0265	.00100	.05004	.00012
%RSD	4.4183	.44204	284.17	1.2732	2.3787	217.51	16.513
#1	.99007	2.5739	01304	2.0553	04110	06933	.00061
#2	.92243	2.5558	.00558	2.0714	04110	02304	.00062

Analysis	Report			05/19	/00 09:23:	58 AM 65	7 877 page 2
#3	.98210	2.5799	03190	2.1178	04318	.04649	.00086
#4	.90650	2.5799	.01164	2.0838	04214	04613	.00071
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	100.00	50.000	50.000	100.00	100.00
Low	25000	50000	10000	05000	05000	-2.0000	05000
Elem Units Avge SDev %RSD	ZN ppm .03883 .00087 2.2294						

#1 .03782 #2 .03943 #3 .03966 #4 .03841

Errors LC Pass High 100.00 Low -.02000

Operator: MTW Sample Name: DD6A4DT Method: QUANMET

Run Time: 05/19/00 09:24:05

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode. co.	.,,	•••					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.91226	.04577	4.9588	.13467	43.834	.00003	H696.82
SDev	.00710	.00461	.1008	.00556	.547	.00000	3.74
%RSD	.77787	10.078	2.0323	4.1296	1.2483	.29469	.53722
#1	.91831	.05086	5.0958	.13960	44.274	.00003	H700.87
#2	.91831	.04713	4.9687	.13182	44.336	.00003	H699.11
#3	.90469	.04529	4.9051	.12823	43.295	.00003	H693.28
#4	.90772	.03979	4.8654	.13902	43.431	.00003	H694.02
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.95485	00178	4.3155	.00877	.03199	1.4662	00172
SDev	.01219	.00115	.0223	.00409	.00067	.1773	.00012
%RSD	1.2767	64.353	.51601	46.586	2.1021	12.095	7.1201
#1	.94990	00018	4.3383	.01177	.03216	1.5806	00171
#2	.97280	00172	4.3303	.01270	.03146	1.5071	00163
#3	.95109	00259	4.2921	.00439	.03287	1.5724	00189
#4	.94562	00264	4.3012	.00623	.03146	1.2048	00163
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	8.6196	.08729	.00249	S3375.3	.00086	4.4898	.01726
SDev	.0817	.00052	.00248	.2	.00872	.0115	.02197
%RSD	.94744	.59115	99.807	.00567	1009.2	.25655	127.30
#1	8.6735	.08755	00124	S3375.2	01058	4.4802	.02226
#2	8.7039	.08755	.00373	S3375.2	.00331	4.4955	00920
#3	8.5354	.08652	.00373	S3375.6	.01039	4.5032	.01228
#4	8.5658	.08755	.00373	S3375.4	.00034	4.4801	.04369
Errors	LC Pass	LC Pass	LC Pass	LC High	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.98310	2.6069	.01834	2.1093	04257	.01754	.00085
SDev	.04460	.0312	.00746	.0240	.00017	.04486	.00004
%RSD	4.5363	1.1973	40.671	1.1362	.40733	255.75	4.3205
#1	1.0299	2.6339	.01992	2.1286	04283	.01174	.00083
#2	.99007	2.6339	.02184	2.1312	04248	.00018	.00091

Analysis	Report			05/19,	/00 09:27:	10 AM	657 879 page 2
#3 # <b>4</b>	.92242 .99007	2.5799 2.5799	.00746 .02416	2.0848 2.0927	04248 04248	02300 .08124	.00083 .00085
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	
Elem Units Avge SDev %RSD	ZN ppm .04004 .00136 3.3871						
#1 #2 #3 #4	.04189 .03994 .03972 .03863						
Errors High Low	LC Pass 100.00 02000						

Method: QUANMET Sample Name: DD6A5T Run Time: 05/19/00 09:27:18 Operator: MTW

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Mode. Co.							
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00302	.02789	.21030	00980	1.2211	00001	H758.93
SDev	.00002	.00620	.03456	.00509	.0090	.00006	3.20
%RSD	.72364	22.226	16.432	51.962	.73589	893.73	.42103
#1	00300	.03420	.23996	00637	1.2334	.00001	H763.47
#2	00302	.02700	.21626	00637	1.2164	.00003	H758.03
#3	00305	.01970	.22437	00932	1.2218	00010	H758.25
#4	00302	.03063	.16061	01715	1.2128	.00003	H755.97
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR ppm00584 .00126 21.615	CU	FE	K_	LI
Units	ppm	ppm		ppm	ppm	ppm	ppm
Avge	00303	00089		.00439	.01845	1.1926	00130
SDev	.00192	.00299		.00131	.00091	.3631	.00029
%RSD	63.320	334.15		29.819	4.9121	30.449	22.242
#1	00074	.00091	00428	.00531	.01880	1.4662	00101
#2	00462	.00237	00541	.00531	.01950	1.3110	00145
#3	00460	00343	00712	.00254	.01740	.65756	00163
#4	00215	00343	00655	.00438	.01810	1.3355	00111
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	9.0279	.08006	.00249	S3374.0	.00030	.01154	.02384
SDev	.0608	.00052	.00476	.2	.00651	.01936	.01997
%RSD	.67394	.64627	191.28	.00529	2199.0	167.72	83.782
#1	9.1128	.08032	.00870	S3373.8	00787	.03087	.03691
#2	9.0299	.08032	.00373	S3374.1	.00625	01533	.02654
#3	8.9913	.08032	00124	S3374.2	.00476	.01535	.03709
#4	8.9774	.07928	00124	S3374.0	00196	.01529	00519
Errors	LC Pass	LC Pass	LC Pass	LC High	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	01685	2.1956	00437	2.3645	04769	.01762	.00261
SDev	.05965	.0001	.01822	.0159	.00028	.06303	.00429
%RSD	354.09	.00245	417.16	.67137	.59382	357.77	164.12
#1	05266	2.1956	00275	2.3865	04769	.02335	.00789
#2	.01897	2.1956	01108	2.3563	04803		.00264

657	881	
	page	2

05/19/00 09:30:23 AM

Analysis Report

#3	.04683	2.1955	02344	2.3651	04769	02284	00262
#4	08052	2.1956	.01981	2.3502	04734	03447	.00254
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	100.00	50.000	50.000	100.00	100.00
Low	25000	50000	10000	05000	05000	-2.0000	05000

Elem ZNUnits ppm Avge .05556 SDev .00117 %RSD 2.1035 #1 .05541 #2 .05399 .05668 #3 #4 .05616 Errors LC Pass 100.00 High

-.02000

Low

Sample Name: CCV2-1 0014-087-7 Operator: MTW Method: QUANMET

Run Time: 05/19/00 09:30:30

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode. co.	.,						
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.97794	48.665	4.9967	4.8486	4.7640	4.8426	50.197
SDev	.00419	.392	.0537	.0422	.0445	.0419	.182
%RSD	.42847	.80609	1.0752	.87074	.93447	.86464	.36162
#1	.98159	49.095	5.0666	4.9004	4.8132	4.8945	50.380
#2	.98147	48.890	4.9774	4.8569	4.7887	4.8557	50.323
#3	.97515	48.413	4.9388	4.7988	4.7365	4.8217	50.020
#4	.97357	48.261	5.0039	4.8381	4.7175	4.7983	50.063
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	55.000	5.5000	5.5000	5.5000	5.5000	55.000
Low	.90000	45.000	4.5000	4.5000	4.5000	4.5000	45.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	4.9809	4.8666	4.8566	4.7790	50.563	48.083	4.7086
SDev	.0106	.0225	.0231	.0466	.299	1.107	.0646
%RSD	.21265	.46230	.47640	.97552	.59121	2.3020	1.3720
#1	4.9823	4.8872	4.8851	4.8317	50.902	48.949	4.7720
#2	4.9881	4.8843	4.8657	4.8030	50.718	49.072	4.7561
#3	4.9655	4.8526	4.8395	4.7513	50.376	46.817	4.6597
#4	4.9876	4.8424	4.8361	4.7300	50.256	47.495	4.6465
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.5000	5.5000	5.5000	5.5000	55.000	55.000	5.5000
Low	4.5000	4.5000	4.5000	4.5000	45.000	45.000	4.5000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	48.845	4.8577	4.8554	47.680	4.8395	4.8897	4.8539
SDev	.421	.0262	.0131	.653	.0393	.0323	.0292
%RSD	.86235	.54003	.26988	1.3695	.81147	.66101	.60091
#1	49.302	4.8864	4.8642	48.332	4.8975	4.8996	4.8885
#2	49.090	4.8730	4.8691	48.151	4.8211	4.8920	4.8668
#3	48.587	4.8398	4.8442	47.149	4.8279	4.8451	4.8356
#4	48.402	4.8315	4.8442	47.089	4.8115	4.9220	4.8247
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	55.000	5.5000	5.5000	55.000	5.5000	5.5000	5.5000
Low	45.000	4.5000	4.5000	45.000	4.5000	4.5000	4.5000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	5.0441	4.8677	4.8904	4.7900	4.8101	9.5297	4.8671
SDev	.0824	.0107	.0431	.0444	.0342	.1780	.0297
%RSD	1.6331	.21986	.88231	.92672	.71008	1.8677	.61098
#1	4.9607	4.8831	4.9070	4.8398	4.8485	9.7692	4.8998
#2	5.0317	4.8649	4.8987	4.8130	4.8284	9.5395	4.8844

Analysis	Report			05/19,	/00 09:33:	35 AM	657 <sub>page</sub> 3 <sub>2</sub>
#3 #4	5.1578 5.0260	4.8584 4.8643	4.8283 4.9276	4.7643 4.7427	4.7881 4.7753	9.4624 9.3479	4.8452 4.8388
Errors High Low	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 11.000 9.0000	LC Pass 5.5000 4.5000
Elem Units Avge SDev %RSD	ZN ppm 4.8772 .0165 .33817						
#1 #2 #3 #4	4.8866 4.8872 4.8824 4.8527						
Errors High Low	LC Pass 5.5000 4.5000						

05/19/00 09:36:43 AM page 1

Operator: MTW

Method: QUANMET Sample Name: CCB1 Operate Run Time: 05/19/00 09:33:38 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: Cor	WC COLL.	raccor. I					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00076	.01173	.03267	00792	.00082	.00053	.01346
SDev	.00195	.00526	.05330	.00674	.00044	.00042	.01035
%RSD	256.99	44.834	163.15	85.160	53.053	79.847	76.875
#1	.00302	.01043	02191	00998	.00072	.00018	.00758
#2	.00153	.01951	.09713	00999	.00042	.00017	.00270
#3	00152	.00851	.05360	01357	.00072	.00075	.01773
#4	.00002	.00849	.00185	.00187	.00144	.00102	.02581
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00066	00181	.00000	.00046	.00720	10006	00096
SDev	.00348	.00183	.00257	.00088	.00210	.18172	.00074
%RSD	526.38	100.96	52681.	192.03	29.137	181.61	76.920
#1	00010	00145	.00200	.00069	.00632	02042	00101
#2	00371	.00001	00370	00023	.00492	23280	00117
#3	.00454	00436	.00143	00023	.00773	26547	.00005
#4	.00191	00146	.00029	.00162	.00984	.11844	00171
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00069	.00027	.01739	.02696	00301	02125	00009
SDev	.00597	.00060	.00248	.00906	.00321	.00386	.02015
%RSD	864.10	224.31	14.285	33.589	106.48	18.160	22101.
#1	00207	.00078	.01366	.02176	00211	02317	00535
#2	.00069	00025	.01863	.02006	00468	01546	02640
#3	.00898	00025	.01863	.02600	.00104	02322	.01573
#4	00484	.00078	.01863	.04001	00631	02314	.01565
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	01490	01501	01966	.00061	.00078	.01731	.00034
SDev	.03125	.00004	.01003	.00062	.00066	.07472	.00300
%RSD	209.76	.25508	51.014	102.64	84.132	431.71	893.43
#1	.00301	01505	02175	.00007	00017	.04629	00231
#2	.00698	01498	02392	.00007	.00087	01165	.00293

Analysis	Report			05/19,	657 885 page 2		
#3 #4	06066 00892	01505 01498	02786 00510	.00115 .00115	.00121 .00121	06951 .10411	00222 .00294
Errors High Low	LC Pass .25000 25000	LC Pass .50000 50000	LC Pass .10000 10000	LC Pass .05000 05000	LC Pass .05000 05000	LC Pass .30000 30000	LC Pass .05000 05000
Elem Units Avge SDev %RSD	ZN ppm 00017 .00053 312.08						
#1 #2 #3 #4	~.00064 00060 .00038 .00018						
Errors High Low	LC Pass .02000 02000						

Analysis Report

05/19/00 09:39:51 AM page 1

Operator: MTW Method: QUANMET Sample Na Run Time: 05/19/00 09:36:46 Sample Name: DD6A6T

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00266	.03297	.19638	00008	1.2934	.00010	H837.50
SDev	.00144	.00969	.02943	.01890	.0228	.00008	7.82
%RSD	54.130	29.381	14.989	23901.	1.7647	75.927	.93428
#1	00155	.03259	.24016	01568	1.2660	.00004	H828.14
#2	00152	.03250	.18444	.01732	1.2848	.00017	H833.96
#3	00304	.04523	.17638	01716	1.3182	.00003	H843.72
#4	00455	.02154	.18453	.01521	1.3046	.00017	H844.20
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00189	00021	00484	.00808	.01369	1.5132	00079
SDev	.00108	.00322	.00242	.00377	.00120	.3503	.00191
%RSD	57.202	1528.9	49.917	46.631	8.7795	23.153	241.92
#1	.00289	.00089	00256	.00439	.01527	1.5234	.00207
#2	.00217	.00233	00427	.00531	.01246	1.5234	00163
#3	.00035	00493	00826	.01085	.01316	1.0741	00171
#4	.00217	.00087	00428	.01177	.01387	1.9318	00189
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	13.029	.13041	.00124	5843.06	00140	01348	.02908
SDev	.185	.00133	.00287	1686.4	.00409	.03414	.01006
%RSD	1.4225	1.0227	230.55	200.03	293.18	253.35	34.612
#1	12.782	.12886	00124	S11792	00323	.00776	.03695
#2	13.020	.12990	.00373	S11792	.00362	03844	.03700
#3	13.224	.13196	.00373	S11792	00587	04637	.02640
#4	13.091	.13093	00124	S3372.6	00010	.02315	.01596
Errors	LC Pass	LC Pass	LC Pass	LC High	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00194	3.0981	.00343	2.2265	L05263	.04690	.00130
SDev	.03429	.0492	.00778	.0389	.00077	.06820	.00261
%RSD	1767.9	1.5867	226.68	1.7486	1.4612	145.40	200.03
#1	.02692	3.0545	.00347	2.1794	L05150	.07006	00261
#2	.02691	3.0606	.01175	2.2116	L05289	.04688	.00264

Analysis	Analysis Report				05/19/00 09:39:51 AM			
#3 #4	04074 02084	3.1567 3.1206	.00549 00697	2.2671 2.2478	L05323 L05289	.11643 04575	.00263 .00255	
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Low 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000	
Elem Units Avge SDev %RSD	ZN ppm .04895 .00121 2.4764							
#1 #2 #3 #4	.04862 .04838 .04807 .05074							
Errors High Low	LC Pass 100.00 02000							

05/19/00 09:43:04 AM page 1 Analysis Report Operator: MTW

Method: QUANMET Sample Name: DD6A7T Operate Run Time: 05/19/00 09:39:58 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CO	NC COII.	ractor: 1					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00077	.03662	.15265	.04098	1.2356	.00007	H779.96
SDev	.00261	.00829	.03094	.00180	.0069	.00007	3.29
%RSD	338.25	22.647	20.268	4.3867	.55987	94.038	.42131
#1	00303	.02697	.13680	.04188	1.2445	.00017	H782.66
#2	.00149	.04354	.17648	.04188	1.2341	.00004	H781.55
#3	00303	.03244	.11688	.04188	1.2362	.00003	H780.40
#4	.00149	.04354	.18046	.03828	1.2277	.00004	H775.23
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00078	00308	00556	.00508	.01792	1.3560	00104
SDev	.00298	.00073	.00164	.00116	.00120	.1450	.00057
%RSD	381.52	23.727	29.439	22.897	6.7048	10.692	55.264
#1	.00041	00345	00371	.00346	.01669	1.2212	00083
#2	.00278	00345	00541	.00531	.01950	1.3600	00135
#3	00393	00198	00541	.00623	.01739	1.5561	00034
#4	00237	00343	00769	.00531	.01810	1.2865	00163
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	12.426	.08961	.00000	\$3374.0	.00874	.00759	.00020
SDev	.050	.00084	.00248	.2	.00823	.02086	.01049
%RSD	.40312	.94134	92681.	.00533	94.133	274.67	5323.2
#1	12.467	.08961	00124	S3374.0	.01730	.03841	00503
#2	12.431	.09065	00124	S3373.8	.01380	00013	00498
#3	12.450	.08961	.00373	S3374.2	.00431	00009	00513
#4	12.354	.08858	00124	S3373.8	00046	00781	.01593
Errors	LC Pass	LC Pass	LC Pass	LC High	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	03575	2.6100	00085	1.9074	04847	.00324	00001
SDev	.02352	.0351	.01502	.0105	.00059	.04669	.00301
%RSD	65.784	1.3434	1757.3	.55133	1.2221	1441.6	27485.
#1	03277	2.6341	.01360	1.9211	04873	02282	.00255
#2	01286	2.5799	02157	1.9055	04769	.05825	00261

Analysis	Report			05/19/00 09:43:04 AM			657 889 page 2
#3	06859	2.6461	.00533	1.9076	04907	04600	.00264
#4	02879	2.5799	00078	1.8955	04838	.02352	00262
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	100.00	50.000	50.000	100.00	100.00
Low	25000	50000	10000	05000	05000	-2.0000	05000
Elem Units Avge SDev %RSD	ZN ppm .03088 .00061 1.9872						

6095

Operator: MTW

Method: QUANMET Sample Name: CCV2-2 Run Time: 05/19/00 09:43:11

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: co	NC COLL.	140001. 1					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.98469	48.905	5.0211	4.8656	4.8021	4.8903	50.775
SDev	.01026	.826	.0477	.0841	.0932	.0818	.441
%RSD	1.0420	1.6883	.94924	1.7283	1.9410	1.6736	.86936
#1	.98951	49.280	5.0596	4.8905	4.8448	4.9391	51.065
#2	.99410	49.538	5.0651	4.9159	4.8709	4.9425	51.128
#3	.97041	47.695	4.9775	4.7407	4.6648	4.7695	50.163
#4	.98475	49.106	4.9824	4.9152	4.8278	4.9103	50.744
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	55.000	5.5000	5.5000	5.5000	5.5000	55.000
Low	.90000	45.000	4.5000	4.5000	4.5000	4.5000	45.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	5.0231	4.9183	4.9023	4.8040	51.039	48.179	4.6879
SDev	.0385	.0557	.0504	.0912	.663	1.290	.1285
%RSD	.76579	1.1317	1.0288	1.8978	1.2989	2.6771	2.7407
#1	5.0403	4.9507	4.9375	4.8475	51.411	48.557	4.7038
#2	5.0484	4.9608	4.9323	4.8678	51.521	49.382	4.8071
#3	4.9658	4.8382	4.8287	4.6690	50.072	46.352	4.5065
#4	5.0380	4.9233	4.9107	4.8317	51.150	48.427	4.7343
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.5000	5.5000	5.5000	5.5000	55.000	55.000	5.5000
Low	4.5000	4.5000	4.5000	4.5000	45.000	45.000	4.5000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	49.093	4.9097	4.9027	47.688	4.8828	4.9411	4.9117
SDev	.805	.0587	.0579	1.178	.0544	.0773	.0760
%RSD	1.6392	1.1954	1.1817	2.4708	1.1136	1.5650	1.5466
#1	49.521	4.9476	4.9189	47.723	4.9290	5.0245	4.8884
#2	49.617	4.9538	4.9537	48.795	4.9208	4.9554	4.9722
#3	47.899	4.8263	4.8194	46.048	4.8112	4.8374	4.8144
#4	49.336	4.9113	4.9189	48.188	4.8702	4.9469	4.9721
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	55.000	5.5000	5.5000	55.000	5.5000	5.5000	5.5000
Low	45.000	4.5000	4.5000	45.000	4.5000	4.5000	4.5000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	5.0975	4.8982	4.9620	4.8247	4.8475	9.7618	4.9100
SDev	.1207	.0616	.0652	.0893	.0735	.1742	.0621
%RSD	2.3674	1.2574	1.3140	1.8504	1.5170	1.7844	1.2652
#1	5.1178	4.8777	4.9277	4.8657	4.8870	9.7634	4.9487
#2	5.1620	4.9619	5.0501	4.8916	4.9012	9.9591	4.9545

Analysis	Report			05/19/00 09:46:17 AM			657 891 page 2
#3 #4	4.9219 5.1885	4.8220 4.9314	4.9007 4.9695	4.6935 4.8478	4.7399 4.8620	9.5352 9.7895	4.8202 4.9164
Errors High Low	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 11.000 9.0000	LC Pass 5.5000 4.5000
Elem Units Avge SDev %RSD	ZN ppm 4.9054 .0394 .80375						
#1 #2 #3 #4	4.9294 4.9356 4.8491 4.9075						
Errors High Low	LC Pass 5.5000 4.5000						

Analysis Report

Method: QUANMET Sample Name: CCB2 Operator: MTW

Method: QUANMET Sample Na Run Time: 05/19/00 09:46:20

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00188	.01460	.02176	.01032	.00059	.00050	.01390
SDev	.00227	.01008	.04648	.00992	.00026	.00029	.01287
%RSD	120.71	69.047	213.54	96.076	43.590	58.163	92.560
#1	00152	.00133	.05367	.00397	.00020	.00018	.00011
#2	.00302	.01226	02193	.01879	.00072	.00047	.00819
#3	.00302	.02151	.06939	.01878	.00072	.00047	.01718
#4	.00302	.02328	01408	00024	.00072	.00089	.03013
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR ppm00242 .00126 52.305	CU	FE	K_	LI
Units	ppm	ppm		ppm	ppm	ppm	ppm
Avge	.00261	.00144		.00023	.00632	.10415	00080
SDev	.00343	.00205		.00119	.00385	.32065	.00037
%RSD	131.34	142.10		513.00	60.866	307.88	46.553
#1	00228	.00001	00313	00115	.00211	33082	00119
#2	.00353	.00144	00199	00023	.00703	.25731	00075
#3	.00346	.00434	00370	.00162	.00492	.41251	00093
#4	H.00572	00002	00085	.00069	.01124	.07760	00031
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge '	.00138	.00026	.01366	.02611	00984	00381	.00775
SDev	.00727	.00060	.00406	.01501	.00248	.01939	.00999
%RSD	525.99	225.82	29.691	57.494	25.198	508.35	128.86
#1	00207	00025	.00869	.01879	00831	02313	00531
#2	00484	00025	.01863	.01666	01216	00766	.01552
#3	.00069	.00078	.01366	.02048	01173	00760	.00519
#4	.01174	.00078	.01366	.04850	00716	.02312	.01562
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem Units Avge SDev %RSD	SE ppm 02386 .03072 128.77	SI ppm 01054 .00300 28.483	SN ppm 00255 .03307 1296.3	SR ppm .00034 .00038 113.14	TI ppm .00069 .00060 86.603	TL ppm .02889 .04867 168.45	V_ppm 00232 .00008 3.4202
#1	03680	00904	02796	.00007	.00017	.01157	00242
#2	.01893	01505	.03628	.00034	.00052	02323	00222

Analysis	Report			05/19/00 09:49:25 AM			657893	
#3 #4	02486 05269	00904 00904	03205 .01352	.00007	.00052 .00156	.09257 .03464	00232 00232	
Errors High Low	LC Pass .25000 25000	LC Pass .50000 50000	LC Pass .10000 10000	LC Pass .05000 05000	LC Pass .05000 05000	LC Pass .30000 30000	LC Pass .05000 05000	
Elem Units Avge SDev %RSD	ZN ppm .00077 .00112 145.47							
#1 #2 #3 #4	00037 .00043 .00073 .00231							
Errors High Low	LC Pass .02000 02000							

Method: QUANMET Sample Name: DDDFQBF Operator: MTW

Run Time: 05/19/00 09:49:27

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

11000.							
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00112	.00413	00198	00153	.00026	00006	.09511
SDev	.00076	.00528	.02404	.01623	.00034	.00007	.00046
%RSD	67.361	127.75	1212.6	1063.8	131.84	110.42	.48617
#1	.00150	.00321	.02981	00025	.00072	00010	.09563
#2	00001	.01052	.00194	00385	.00020	00010	.09507
#3	.00150	00227	02577	02078	00010	.00004	.09523
#4	.00150	.00506	01391	.01878	.00020	00010	.09451
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR ppm00285 .00151 52.934	CU	FE	K_	LI
Units	ppm	ppm		ppm	ppm	ppm	ppm
Avge	00055	00036		00023	.00386	12865	00135
SDev	.00291	.00139		.00075	.00146	.12031	.00156
%RSD	527.31	384.18		326.85	37.869	93.512	115.58
#1	.00067	00000	00370	.00069	.00562	.01225	00145
#2	.00191	00145	00427	00023	.00421	28181	00259
#3	00004	.00145	00256	00023	.00351	12661	.00088
#4	00475	00144	00085	00116	.00211	11844	00225
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem Units Avge SDev %RSD	MG ppm 00760 .00813 107.05	MN ppm 00026 .00000	MO ppm .00248 .00248 99.984	NA ppm .01337 .00400 29.935	NI ppm 00833 .00483 57.961	PB ppm 01157 .03182 275.12	SB ppm .00522 .02591 496.36
#1	00760	00026	.00373	.01497	00902	01541	.00525
#2	01589	00026	.00373	.01199	01346	04632	02651
#3	.00345	00026	.00373	.01794	00902	.03087	.00518
#4	01036	00026	00124	.00860	00180	01541	.03697
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem Units Avge SDev %RSD	SE ppm 01491 .01360 91.199	SI ppm 01205 .00347 28.793	SN ppm 01342 .03053 227.47	SR ppm .00007 .00000	TI ppm 00017 .00028 163.30	TL ppm .02024 .02738 135.22	V_ ppm 00254 .00005 1.7895
#1	.00301	00904	02791	.00007	.00017	.02312	00251
#2	01292	00904	.01953	.00007	00017	00002	00251

STL Pittsburgh

6100

Analysis	Report			05/19	/00 09:52:	33 AM	657 895 page 2
#3 #4	02884 02089	01505 01505	.00313 04845	.00007	00052 00017	00003 .05791	00251 00260
Errors High Low	LC Pass .25000 25000	LC Pass .50000 50000	LC Pass .10000 10000	LC Pass .05000 05000	LC Pass .05000 05000	LC Pass .30000 30000	LC Pass .05000 05000
Elem Units Avge SDev %RSD	ZN ppm .00590 .00153 25.880					·	
#1 #2 #3 #4	.00611 .00368 .00691 .00689						
Errors High Low	LC Pass .02000 02000						

Operator: MTW

Method: QUANMET Sample Name: DDDFQCF Run Time: 05/19/00 09:52:36

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	L00038	L.00598	L.00888	02077	L.00036	L00007	50.021
SDev	.00077	.00486	.03303	.00001	.00011	.00008	.115
%RSD	203.15	81.258	371.84	.03214	29.167	107.25	.23067
#1	L00153	L.00884	L02586	02078	L.00042	L.00004	49.902
#2	L.00002	L.00495	L.00189	02077	L.00042	L00011	49.944
#3	L00001	L00041	L.05368	02077	L.00020	L00010	50.132
#4	L.00001	L.01054	L.00582	02077	L.00042	L00011	50.107
Errors	LC Low	LC Low	LC Low	NOCHECK	LC Low	LC Low	LC Pass
High	.06000	2.4000	2.4000		2.4000	.06000	60.000
Low	.04000	1.6000	1.6000		1.6000	.04000	40.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	L00091	L00072	L00171	L00023	L.00702	46.770	00084
SDev	.00357	.00250	.00136	.00075	.00099	.457	.00038
%RSD	392.71	348.16	79.269	326.23	14.168	.97662	45.316
#1	L.00399	L.00289	L00085	L00023	L.00561	47.356	00057
#2	L00171	L00144	L00370	L00115	L.00772	46.630	00083
#3	L00457	L00288	L00142	L.00069	L.00701	46.834	00137
#4	L00135	L00144	L00086	L00023	L.00772	46.262	00057
Errors	LC Low	LC Low	LC Low	LC Low	LC Low	LC Pass	NOCHECK
High	.06000	.60000	.24000	.30000	1.2000	60.000	
Low	.04000	.40000	.16000	.20000	.80000	40.000	
Elem Units Avge SDev %RSD	MG ppm 48.837 .067	MN ppm L00026 .00000 .70206	MO ppm 00372 .00497 133.37	NA ppm 46.816 .160 .34261	NI ppm L00635 .01028 162.01	PB ppm L.03081 .01991 64.632	SB ppm L.00526 .01491 283.35
#1	48.822	L00026	00621	47.042	L.00117	L.03860	L00526
#2	48.802	L00025	.00373	46.808	L00853	L.02306	L00524
#3	48.788	L00026	00621	46.675	L.00197	L.05390	L.02637
#4	48.935	L00026	00621	46.737	L02001	L.00767	L.00517
Errors	LC Pass	LC Low	NOCHECK	LC Pass	LC Low	LC Low	LC Low
High	60.000	.60000		60.000	.60000	.60000	.60000
Low	40.000	.40000		40.000	.40000	.40000	.40000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	L.00301	.01351	00617	.00185	00251	L.03467	L00008
SDev	.03747	.00579	.02633	.00018	.00033	.02502	.00303
%RSD	1244.7	42.811	426.56	9.9329	13.206	72.165	4024.7
#1	L.03882	.00898	.02171	.00201	00260	L00006	L00270
#2	L03678	.01505	03829	.00201	00225	L.03465	L.00264

6	5.7	ge ge	ð	7
	рa	qe	Z	

05/19/00	09:55:41	ΑM
----------	----------	----

Analysis Report

#3 # <b>4</b>	L.03087 L02087	.00898 .02106	.00738 01548	.00169 .00169	00295 00225	L.05787 L.04623	L00270 L.00246
Errors High Low	LC Low 2.4000 1.6000	NOCHECK	NOCHECK	NOCHECK	NOCHECK	LC Low 2.4000 1.6000	LC Low .60000 .40000
Elem Units Avge SDev %RSD	ZN ppm L.00355 .00157 44.163						
#1 #2 #3 #4	L.00529 L.00369 L.00148 L.00372						
Errors High Low	LC Low .60000 .40000						

Operator: MTW

Sample Name: DC7M7F Method: QUANMET

Run Time: 05/19/00 09:55:48 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00056	.18227	00482	.04343	.09306	.00022	20.800
SDev	.00124	.00239	.02458	.00073	.00034	.00007	.104
%RSD	219.56	1.3111	509.51	1.6874	.36438	32.268	.49809
#1	00056	.18129	.02798	.04453	.09301	.00018	20.690
#2	00057	.18323	.00012	.04307	.09301	.00018	20.925
#3	.00095	.17952	02366	.04303	.09271	.00033	20.747
#4	00208	.18504	02374	.04308	.09352	.00018	20.841
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00008	.02313	00199	.00029	3.3017	.14499	.00703
SDev	.00304	.00204	.00081	.00116	.0120	.04686	.00217
%RSD	3929.8	8.8357	40.413	406.55	.36495	32.323	30.863
#1	00137	.02024	00313	.00190	3.2924	.10211	.00516
#2	.00312	.02457	00142	00087	3.3106	.20830	.00516
#3	.00204	.02458	00199	.00005	3.2902	.11844	.00868
#4	00347	.02314	00142	.00006	3.3134	.15112	.00912
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	14.890	6.6480	.00174	24.905	.02060	.00365	.00292
SDev	.065	.0262	.00287	.216	.00752	.02630	.01010
%RSD	.43385	.39360	165.14	.86590	36.515	720.60	345.92
#1	14.840	6.6250	.00422	24.924	.02729	00028	.00561
#2	14.879	6.6746	00075	24.807	.02286	.03065	00497
#3	14.857	6.6261	.00422	24.693	.02244	03097	.01608
#4	14.984	6.6664	00075	25.196	.00981	.01520	00504
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00982	11.725	00609	.09330	.00173	01996	00255
SDev	.05846	.063	.01075	.00027	.00020	.05978	.00005
%RSD	595.34	.53649	176.53	.28952	11.547	299.46	2.1218
#1	.00308	11.716	01073	.09343	.00156	.04951	00250
#2	.06284	11.747	.00370	.09343	.00191	08939	00260

Analysis	Report			05/19/00 09:58:53 AM			657 899 page 2
#3 #4	07652 02868	11.644 11.795	.00173 01905	.09289 .09343	.00191 .00156	.00318 04314	00250 00260
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	
Elem Units Avge SDev %RSD	ZN ppm .01685 .00168 9.9830						
#1 #2 #3 #4	.01779 .01615 .01861 .01484						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW Method: QUANMET Sample Name: DC7MAF

Run Time: 05/19/00 09:58:56

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00177	.01746	00779	.01255	2.7981	.00011	181.90
SDev	.00245	.01169	.02467	.00129	.0122	.00008	.42
%RSD	138.04	66.921	316.64	10.302	.43661	75.712	.22968
#1	00177	.00322	04439	.01320	2.7906	.00003	181.68
#2	00180	.02154	00074	.01315	2.7954	.00018	181.51
#3	00476	.01422	.00689	.01061	2.8162	.00017	182.46
#4	.00124	.03087	.00707	.01323	2.7904	.00004	181.96
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00012	.00549	00328	.00143	32.289	2.5465	.01397
SDev	.00246	.00247	.00157	.00001	.101	.1463	.00203
%RSD	2006.8	45.092	47.887	.62676	.31249	5.7463	14.568
#1	.00182	.00440	00256	.00143	32.241	2.6425	.01220
#2	00055	.00295	00484	.00143	32.213	2.4056	.01220
#3	00344	.00585	00427	.00145	32.437	2.4383	.01573
#4	.00167	.00874	00142	.00143	32.264	2.6997	.01573
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	36.139	9.4895	.00359	43.943	00095	.00253	.01851
SDev	.150	.0256	.00405	.236	.00581	.02546	.03032
%RSD	.41376	.27003	112.99	.53605	608.37	1005.9	163.83
#1	36.038	9.4781	.00358	43.791	00618	01488	02637
#2	36.065	9.4688	.00854	44.062	.00697	.03910	.03702
#3	36.361	9.5267	.00361	44.213	00038	01482	.02650
#4	36.093	9.4843	00138	43.706	00423	.00072	.03689
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01257	19.171	01546	.54233	01058	.02398	00026
SDev	.02729	.055	.02439	.00195	.00049	.06761	.00266
%RSD	217.11	.28595	157.75	.36010	4.6368	281.94	1024.5
#1	.03727	19.105	.00708	.5 <b>41</b> 61	01058	04797	.00135
#2		19.159	02988	.54199	01023	00137	00242

Analysis Report				05/19,	657 901 page 02		
#3 #4	02569 .01348	19.183 19.237	04229 .00324	.54512 .54059	01127 01023	.11238 .03288	.00264 00261
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000		LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .01863 .00150 8.0691						
#1 #2 #3 #4	.01644 .01987 .01910 .01911						
Errors High Low	LC Pass 100.00 02000						

657 902

Analysis Report

Method: QUANMET Sample Name: DC7MAFP5 Operator: MTW

Run Time: 05/19/00 10:02:03

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00008	.00462	00499	00498	.57709	.00001	37.613
SDev	.00190	.00526	.01537	.00075	.00257	.00007	.076
%RSD	2421.4	113.83	308.00	15.033	.44495	1040.5	.20219
#1	00030	.00140	00696	00535	.57722	.00004	37.538
#2	00181	.00147	01094	00532	.57979	.00004	37.587
#3	00030	.00319	01889	00386	.57774	.00004	37.718
#4	.00272	.01240	.01683	00539	.57362	00010	37.611
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR ppm00413 .00118 28.465	CU	FE	K_	LI
Units	ppm	ppm		ppm	ppm	ppm	ppm
Avge	.00110	.00080		.00055	6.8272	.45539	.00303
SDev	.00185	.00138		.00088	.0152	.13544	.00156
%RSD	168.52	173.56		161.50	.22279	29.741	51.524
#1	.00361	.00260	00484	.00124	6.8312	.56771	.00472
#2	.00097	.00116	00541	00061	6.8404	.45335	.00264
#3	00084	00029	00313	.00032	6.8319	.26547	.00106
#4	.00065	00029	00313	.00124	6.8052	.53503	.00371
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	7.4440	2.0007	.00102	8.9498	00183	03233	.01314
SDev	.0232	.0023	.00475	.0403	.00949	.01257	.01005
%RSD	.31201	.11464	465.79	.45066	518.15	38.883	76.442
#1	7.4191	1.9994	.00475	8.9266	.00097	03226	.02632
#2	7.4578	2.0015	00519	9.0035	.00896	04774	.00540
#3	7.4689	2.0036	.00475	8.9568	01380	03235	.01564
#4	7.4302	1.9984	00022	8.9122	00346	01695	.00520
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00406	4.0021	.01689	.11193	00277	~.03895	00257
SDev	.02659	.0179	.00656	.00049	.00020	.07104	.00009
%RSD	654.20	.44820	38.822	.43836	7.2169	182.39	3.6226
#1	.02000	3.9796	.02109	.11224	00295	03612	00250
#2	.01605	4.0096	.00858	.11224	00295	12883	00269

Analysis	Report			05/19	657 903 page 2		
#3 #4	03572 .01592	4.0216 3.9976	.01483	.11202 .11121	00260 00260	.04498 03583	00249 00259
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .00487 .00110 22.564						
#1 #2 #3 #4	.00407 .00642 .00412 .00488						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW

Method: QUANMET Sample Name: DC7MASF

Run Time: 05/19/00 10:05:11

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00074	.01243	00602	.01190	2.7628	.00015	229.38
SDev	.00190	.00155	.01635	.00967	.0065	.00007	.13
%RSD	255.35	12.461	271.62	81.241	.23489	47.756	.05710
#1	00038	.01246	00795	.02637	2.7588	.00018	229.33
#2	00036	.01246	.01579	.00662	2.7671	.00004	229.28
#3	00339	.01429	00804	.00659	2.7693	.00018	229.57
#4	.00115	.01050	02387	.00803	2.7558	.00018	229.35
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00155	.00370	00399	.00139	31.758	49.370	.01386
SDev	.00232	.00278	.00235	.00001	.033	.508	.00024
%RSD	150.35	75.003	58.914	.33504	.10303	1.0289	1.7199
#1 #2 #3 ·	00049 00034 L00503 00032	.00297 .00587 .00588 .00008	00370 00142 00370 00712	.00139 .00139 .00139 .00139	31.717 31.792 31.775 31.749	48.941 49.693 49.913 48.933	.01371 .01397 .01415 .01363
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	84.191	9.3375	.00351	90.461	00403	.00259	.00534
SDev	.163	.0061	.00406	.640	.01050	.02912	.00012
%RSD	.19305	.06522	115.69	.70767	260.57	1125.9	2.3308
#1	83.983	9.3334	00147	90.123	00959	.02377	.00529
#2	84.306	9.3417	.00351	90.946	01266	.03154	.00522
#3	84.334	9.3437	.00351	91.046	.01095	02240	.00551
#4	84.143	9.3313	.00847	89.727	00481	02256	.00532
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.08426	19.132	02958	.53626	01275	.00547	00252
SDev	.01314	.056	.01211	.00089	.00017	.11379	.00008
%RSD	15.592	.29509	40.923	.16540	1.3605	2080.3	3.0071
#1	.07515	19.075	01977	.53578	01266	.15352	00261
#2	.07543	19.189	01975	.53659	01301	04409	00251

Analysis	Report			05/19	657 905 page 2		
#3 #4	.08332 .10313	19.171 19.093	04458 03424	.53735 .53535	01266 01266	.02558 11313	00252 00243
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .01442 .00131 9.0715						
#1 #2 #3 #4	.01268 .01459 .01586 .01457						
Errors High Low	LC Pass 100.00 02000						

657 906

Analysis Report

Operator: MTW

Method: QUANMET Sample Name: DC7MADF Run Time: 05/19/00 10:08:18

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00152	.00789	04162	.00566	2.7513	.00011	229.41
SDev	.00074	.00441	.02441	.00430	.0083	.00008	.93
%RSD	48.815	55.865	58.650	76.073	.30349	71.275	.40566
#1	00189	.00515	05155	00079	2.7478	.00004	229.82
#2	00041	.00511	01958	.00763	2.7563	.00004	228.01
#3	00190	.01438	07144	.00782	2.7414	.00018	229.85
#4	00186	.00694	02391	.00799	2.7599	.00017	229.95
Error	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00124	.00442	00313	.00138	31.649	50.030	.01375
SDev	.00263	.00264	.00147	.00001	.081	.677	.00038
%RSD	212.51	59.650	46.938	.62106	.25638	1.3530	2.7406
#1	.00459	.00731	00256	.00139	31.679	49.864	.01345
#2	00131	.00153	00484	.00137	31.540	50.191	.01363
#3	00032	.00298	00371	.00138	31.643	49.219	.01363
#4	.00199	.00587	00143	.00139	31.733	50.844	.01430
Error	rs LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	84.240	9.3057	00272	91.360	.00064	.02383	.00535
SDev	.191	.0235	.00625	1.210	.01355	.03565	.00008
%RSD	.22674	.25213	230.02	1.3247	2102.8	149.62	1.5080
#1	84.162	9.3137	.00350	90.719	.01894	.05472	.00546
#2	84.375	9.2724	00149	92.262	00139	00705	.00535
#3	84.007	9.3096	01141	89.979	01380	00704	.00526
#4	84.416	9.3272	00146	92.483	00117	.05467	.00534
Erro:		LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High		100.00	50.000	400.00	100.00	100.00	100.00
Low		01500	04000	-5.0000	04000	10000	06000
Elem Unit Avge SDev %RSD	s ppm .00028 .04209	SI ppm 19.147 .078 .40912	SN ppm 00064 .00947 1484.2	SR ppm .53445 .00135 .25215	TI ppm 01318 .00020 1.5193	TL ppm .07878 .05999 76.150	V_ ppm 00135 .00260 192.86
#1	03244	19.231	.01126	.53447	01301	.16535	00252
#2	.05859	19.069	01152	.53481	01335	.06238	00262

Analysis	Report			05/19,	657 907 page 2		
#3 #4	02859 .00358	19.093 19.195	.00091 00321	.53264 .53589	01301 01335	.02678 .06063	00281 .00255
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .01347 .00199 14.786						
#1 #2 #3 #4	.01204 .01616 .01381 .01189						
Errors High Low	LC Pass 100.00 02000						

Method: QUANMET Sample Name: DC7MCF Operator: MTW

Run Time: 05/19/00 10:11:26

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00023	.01069	00804	.08469	.06016	.00001	91.355
SDev	.00227	.00790	.03065	.00201	.00045	.00007	.128
%RSD	1006.0	73.856	381.07	2.3680	.74203	1041.2	.13996
#1	.00166	00030	02680	.08413	.05957	00010	91.257
#2	00287	.01434	.02865	.08209	.06060	.00004	91.463
#3	00136	.01076	.00489	.08626	.06009	.00004	91.232
#4	.00167	.01796	03890	.08629	.06039	.00004	91.468
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00030	.00904	00327	.00056	6.8439	.22872	.01540
SDev	.00228	.00139	.00108	.00088	.0177	.27096	.00214
%RSD	764.31	15.409	32.911	159.08	.25912	118.47	13.893
#1	.00118	.00867	00256	00060	6.8284	.28181	.01617
#2	00077	.01012	00484	.00033	6.8678	15928	.01661
#3	00333	.01013	00256	.00125	6.8326	.32265	.01661
#4	.00171	.00722	00313	.00125	6.8467	.46968	.01220
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	33.123	7.2890	.00102	27.168	.01511	00900	.00548
SDev	.090	.0081	.00476	.344	.00945	.01406	.00871
%RSD	.27235	.11069	464.88	1.2673	62.567	156.20	158.88
#1	33.045	7.2872	00022	26.788	.01352	01671	.00551
#2	33.253	7.3006	.00475	27.556	.02879	02438	.01616
#3	33.087	7.2820	00519	26.990	.00746	.00645	.00543
#4	33.106	7.2861	.00475	27.338	.01066	00136	00517
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI ppm00590 .00020 3.3962	TL	V_
Units	ppm	ppm	ppm	ppm		ppm	ppm
Avge	.02104	12.114	00848	.20530		04413	00256
SDev	.02211	.036	.01748	.00055		.07657	.00009
%RSD	105.09	.30127	206.08	.26674		173.50	3.5268
#1	.03591	12.083	02659	.20484	00572	02950	00259
#2	.00421	12.167	01825	.20592	00572	14564	00250

05/19/	00	10:	14:	31	AM
--------	----	-----	-----	----	----

Analysis Re	port
-------------	------

page 2

#3	.00010	12.107	00176	.20484	00607	.03989	00268
#4	.04393	12.101	.01268	.20561	00607	04127	00249
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	100.00	50.000	50.000	100.00	100.00
Low	25000	50000	10000	05000	05000	-2.0000	05000
Elem	ZN						

Elem ZN
Units ppm
Avge .02648
SDev .00307
%RSD 11.586
#1 .02447
#2 .02687
#3 .03066
#4 .02392

Errors LC Pass High 100.00 Low -.02000

Method: QUANMET Sample Name: DC7MDF Operator: MTW

Run Time: 05/19/00 10:14:34

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00010	.02236	.01867	01837	.68327	.00004	120.66
SDev	.00145	.00382	.07317	.00242	.00270	.00016	.30
%RSD	1388.9	17.089	391.96	13.192	.39541	391.90	.24898
#1	.00124	.01778	.08129	01921	.68700	00010	120.30
#2	00179	.02144	.07331	02070	.68186	.00018	121.02
#3	.00124	.02697	00621	01858	.68340	00010	120.75
#4	00027	.02324	07372	01499	.68083	.00018	120.58
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR ppm00470 .00215 45.803	CU	FE	K_	LI
Units	ppm	ppm		ppm	ppm	ppm	ppm
Avge	00227	00214		.00093	.04581	1.5255	.01720
SDev	.00246	.00138		.00116	.00202	.3438	.00143
%RSD	108.41	64.706		125.73	4.4043	22.538	8.2881
#1	L00583	00322	00655	.00069	.04634	1.3192	.01648
#2	00116	00033	00256	.00254	.04704	1.8828	.01925
#3	00027	00178	00313	.00069	.04704	1.1558	.01604
#4	00181	00323	00655	00023	.04282	1.7440	.01705
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem Units Avge SDev %RSD	MG ppm 21.275 .057	MN ppm 1.3558 .0021 .15709	MO ppm .01118 .00287 25.643	NA ppm 17.742 .099 .55528	NI ppm 00316 .00739 233.68	PB ppm 01739 .00737 42.414	SB ppm .02626 .01489 56.710
#1	21.267	1.3540	.00870	17.862	01380	02321	.01566
#2	21.237	1.3571	.01367	17.734	00087	01538	.04732
#3	21.358	1.3581	.00870	17.751	00131	02316	.02629
#4	21.239	1.3540	.01367	17.621	.00332	00780	.01579
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem Units Avge SDev %RSD	SE ppm 01277 .04396 344.33	SI ppm 10.413 .008	SN ppm .00740 .00453 61.245	SR ppm .45117 .00162 .35922	TI ppm 00711 .00028 3.9829	TL ppm .05641 .05297 93.904	V_ ppm 00237 .00006 2.3568
#1	04858	10.419	.00735	.45360	00711	.06509	00242
#2	.05091	10.413	.01371	.45036	00746	.04192	00232

05/19/00 10:17:39 AM

page 2

#3 #4	03266 02074	10.401 10.419	.00533	.45036 .45036	00676 00711	00438 .12303	00242 00233
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .00688 .00190 27.566						
#1 #2 #3 #4	.00557 .00552 .00957 .00687						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW Method: QUANMET Sample Name: CCV2-3 Run Time: 05/19/00 10:17:41

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Mode: Co.	NC COII.	raccor, r					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.97412	48.027	4.9237	4.8116	4.6919	4.8330	50.803
SDev	.00150	.106	.0986	.0253	.0095	.0059	.118
%RSD	.15398	.22066	2.0019	.52572	.20255	.12294	.23234
#1	.97529	47.895	4.9346	4.7991	4.6812	4.8376	50.883
#2	.97217	48.113	4.7822	4.7989	4.6981	4.8340	50.628
#3	.97372	47.988	4.9738	4.7989	4.6868	4.8244	50.855
#4	.97531	48.112	5.0042	4.8495	4.7014	4.8359	50.847
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	55.000	5.5000	5.5000	5.5000	5.5000	55.000
Low	.90000	45.000	4.5000	4.5000	4.5000	4.5000	45.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	5.0384	4.8916	4.8757	4.7017	50.538	46.664	4.5347
SDev	.0212	.0094	.0066	.0123	.067	.545	.0502
%RSD	.42116	.19297	.13509	.26060	.13285	1.1675	1.1077
#1	5.0436	4.9003	4.8811	4.6885	50.569	45.862	L4.4654
#2	5.0182	4.8960	4.8703	4.7134	50.468	47.038	4.5813
#3	5.0259	4.8916	4.8697	4.6941	50.498	46.964	4.5331
#4	5.0660	4.8785	4.8817	4.7107	50.616	46.793	4.5591
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC. Pass
High	5.5000	5.5000	5.5000	5.5000	55.000	55.000	5.5000
Low	4.5000	4.5000	4.5000	4.5000	45.000	45.000	4.5000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	48.449	4.8608	4.8679	46.769	4.8576	4.9427	4.8538
SDev	.074	.0072	.0117	.439	.0189	.0420	.0233
%RSD	.15305	.14908	.24130	.93797	.38826	.84936	.48066
#1	48.360	4.8647	4.8691	46.150	4.8837	4.9546	4.8460
#2	48.487	4.8512	4.8592	47.146	4.8466	4.8850	4.8247
#3	48.421	4.8595	4.8592	46.785	4.8414	4.9852	4.8670
#4	48.529	4.8678	4.8840	46.996	4.8586	4.9462	4.8773
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	55.000	5.5000	5.5000	55.000	5.5000	5.5000	5.5000
Low	45.000	4.5000	4.5000	45.000	4.5000	4.5000	4.5000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	4.9863	5.0087	4.9903	4.7356	4.7738	9.5994	4.8609
SDev	.0949	.0383	.0400	.0091	.0052	.1424	.0076
%RSD	1.9040	.76448	.80077	.19260	.10798	1.4835	.15640
#1	4.9038	4.9547	5.0333	4.7276	4.7715	9.5874	4.8664
#2	5.1223	5.0146	5.0083	4.7412	4.7715	9.4379	4.8520
							~ ~ ~

05/19/00	10:20:47	AM	657	913 page	2
,,	,			2-25	_

9.7851

9.5871

LC Pass

11.000

9.0000

4.8572

4.8680

LC Pass

5.5000

4.5000

4.7708

4.7816

LC Pass

5.5000

4.5000

Analysis	Report
----------	--------

#3 #4	4.9473 4.9716	5.0207 5.0449	4.9402 4.9796	4.7282 4.7456
	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000
Elem Units Avge SDev %RSD	ZN ppm 4.8641 .0137 .28137			
#1 #2 #3 #4	4.8629 4.8628 4.8487 4.8820			
Errors High Low	LC Pass 5.5000 4.5000			

Method: QUANMET Sample Name: CCB3 Operator: MTW

Run Time: 05/19/00 10:20:49

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: Co.	NC COII.	140001.					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00151	.01181	00407	00369	.00088	.00050	.00722
SDev	.00001	.00759	.02805	.00832	.00043	.00032	.00329
%RSD	.86833	64.294	689.71	225.88	49.362	63.545	45.494
#1 #2 #3 #4	.00150 .00153 .00150	.01414 .00667 .00493 .02149	02194 01800 .03773 01406	.00758 00999 00996 00236	.00072 .00042 .00144 .00093	.00018 .00031 .00089 .00061	.00372 .00646 .01164 .00707
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00051	00000	00156	.00069	.00913	03880	00137
SDev	.00111	.00205	.00205	.00000	.00470	.19288	.00062
%RSD	218.10	82180.	131.09	.24037	51.410	497.12	45.111
#1	00010	00145	00256	.00069	.00632	.02042	00137
#2	.00216	.00289	.00143	.00070	.00492	.20830	00145
#3	.00016	00145	00313	.00069	.01546	18379	00057
#4	00018	00000	00199	.00069	.00984	20013	00207
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem Units Avge SDev %RSD	MG ppm 00553 .00913 165.20	MN ppm 00025 .00000	MO ppm .01366 .00702 51.419	NA ppm .01804 .00504 27.931	NI ppm 00810 .00436 53.860	PB ppm 01928 .00440 22.795	SB ppm 00538 .00861 159.91
#1	01036	00025	.00869	.02091	01053	01547	00544
#2	.00345	00025	.02359	.01539	00631	02304	.00515
#3	.00069	00025	.01366	.02346	01267	01548	01593
#4	01589	00025	.00869	.01242	00289	02314	00531
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00494	.08108	01347	.00063	.00043	.02018	00103
SDev	.02003	.00003	.02055	.00052	.00044	.02193	.00271
%RSD	405.27	.04266	152.50	81.893	100.66	108.67	263.60
#1	02485	.08106	.01549	.00050	.00052	.03470	00242
#2	.02290	.08113	02997	00004	00017	.03464	.00304

Analysis	Report			05/19,	/00 10:23:	54 AM	657 915 page 2
#3 #4	00890 00892	.08106 .08106	02594 01347	.00104 .00104	.00052	.02303 01167	00232 00241
Errors High Low	LC Pass .25000 25000	LC Pass .50000 50000	LC Pass .10000 10000	LC Pass .05000 05000	LC Pass .05000 05000	LC Pass .30000 30000	LC Pass .05000 05000
Elem Units Avge SDev %RSD	ZN ppm .00192 .00161 84.005						
#1 #2 #3 #4	00038 .00233 .00232 .00340						
Errors High Low	LC Pass .02000 02000						

Operator: MTW Method: QUANMET Sample Name: DDCAWB Run Time: 05/19/00 10:23:57

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: co.	NC COII.	raccor. z					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00113	.00415	.01586	00151	.00013	00003	.02879
SDev	.00259	.00953	.02717	.01648	.00052	.00008	.00352
%RSD	228.57	229.39	171.28	1088.5	405.50	241.34	12.230
#1	00152	00958	00585	02077	00031	00010	.02510
#2	.00153	.00683	.01777	.00188	.00072	.00003	.03358
#3	00001	.00690	00201	00596	.00042	.00004	.02840
#4	.00453	.01247	.05355	.01879	00031	00010	.02809
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR ppm00128 .00256 199.88	CU	FE	K_	LI
Units	ppm	ppm		ppm	ppm	ppm	ppm
Avge	.00076	00036		.00023	.01018	.06331	00148
SDev	.00161	.00139		.00092	.00070	.30733	.00035
%RSD	212.58	381.57		398.90	6.8992	485.47	23.407
#1	.00295	00001	00142	00023	.00983	09394	00119
#2	.00039	00145	.00199	.00162	.01124	.50236	00189
#3	.00062	00145	00427	00023	.00983	.03676	00119
#4	00093	.00145	00143	00023	.00983	19196	00163
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00345	00077	00248	.01507	00533	00772	00265
SDev	.01696	.00060	.00248	.00717	.00265	.01661	.02172
%RSD	490.99	77.120	100.06	47.548	49.687	214.97	819.51
#1	01589	00129	00124	.00520	00297	01545	02631
#2	.02003	00026	00124	.01751	00875	.01535	01589
#3	01589	00129	00124	.02218	00609	00773	.01574
#4	00207	00026	00621	.01539	00352	02307	.01585
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	02384	.07057	00518	.00008	00026	.00569	00102
SDev	.01820	.00902	.01562	.00011	.00044	.03600	.00245
%RSD	76.339	12.788	301.51	139.88	167.77	632.61	240.77
#1	03280	.05703	02184	00004	00087	00009	00260
#2	03677	.07512	.00096	.00007	00017	01170	.00256

Analysis	Report			05/19,	/00 10:27:	02 AM	657 917 page 2
#3 #4	02882 .00302	.07505 .07507	.01352 01337	.00023	00017 .00017	02324 .05779	00261 00141
Errors High Low	LC Pass .25000 25000	LC Pass .50000 50000	LC Pass .10000 10000	LC Pass .05000 05000	LC Pass .05000 05000	LC Pass .30000 30000	.05000
Elem Units Avge SDev %RSD	ZN ppm .01176 .00193 16.426						
#1 #2 #3 #4	.01069 .01366 .01308 .00959						
Errors High Low	LC Pass .02000 02000						

Method: QUANMET Sample Name: DDCAWC Operator: MTW

Run Time: 05/19/00 10:27:05

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: Co	NC COII.	raccor. r					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04873	1.8843	1.9425	.92837	1.8926	.04861	L.04054
SDev	.00151	.0070	.0422	.01858	.0134	.00012	.00178
%RSD	3.1060	.37398	2.1738	2.0015	.70864	.23724	4.3828
#1	.04646	1.8816	1.8919	.90101	1.8941	.04861	L.04186
#2	.04949	1.8925	1.9752	.94207	1.8930	.04861	L.04186
#3	.04949	1.8870	1.9236	.93341	1.9079	.04875	L.04033
#4	.04949	1.8761	1.9793	.93698	1.8752	.04847	L.03810
Errors	LC Pass	LC Pass	LC Pass	NOCHECK	LC Pass	LC Pass	LC Low
High	.06000	2.4000	2.4000		2.4000	.06000	60.000
Low	.04000	1.6000	1.6000		1.6000	.04000	40.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04497	.48895	.19361	.23713	1.0432	L.22463	.92576
SDev	.00333	.00342	.00296	.00220	.0078	.05698	.01997
%RSD	7.4005	.70016	1.5267	.92800	.74381	25.368	2.1571
#1	.04075	.49150	.19574	.23944	1.0401	L.24914	.93731
#2	.04529	.48424	.18947	.23574	1.0394	L.27364	.91005
#3	.04495	.49147	.19347	.23851	1.0549	L.14295	.94793
#4	.04889	.48859	.19574	.23482	1.0387	L.23280	.90774
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Low	NOCHECK
High	.06000	.60000	.24000	.30000	1.2000	60.000	
Low	.04000	.40000	.16000	.20000	.80000	40.000	
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	L02418	.48917	.97867	L.02102	.49083	.47144	.47163
SDev	.00983	.00155	.00625	.00212	.01238	.00762	.03132
%RSD	40.675	.31718	.63874	10.084	2.5220	1.6171	6.6417
#1	L01865	.48995	.97743	L.02133	.50206	.47536	.46396
#2	L03246	.48995	.97743	L.01921	.49243	.47515	.46375
#3	L03246	.48996	.98737	L.01964	.47326	.46001	.51617
#4	L01312	.48685	.97246	L.02388	.49556	.47525	.44266
Errors	LC Low	LC Pass	NOCHECK	LC Low	LC Pass	LC Pass	LC Pass
High	60.000	.60000		60.000	.60000	.60000	.60000
Low	40.000	.40000		40.000	.40000	.40000	.40000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.8783	9.3409	1.9640	.95687	.94513	1.9533	.48073
SDev	.0099	.0388	.0370	.00574	.00365	.0768	.00012
%RSD	.52958	.41512	1.8857	.59993	.38645	3.9318	.02488
#1	1.8812	9.3139	1.9096	.9572 <b>7</b>	.94383	2.0343	.48072
#2	1.8653	9.3319	1.9738	.95673	.94521	1.9649	.48070

Analysis	Report			05/19/	657 919 page 2		
#3 #4	1.8773 1.8892	9.3980 9.3199	1.9802 1.9923	.96376 .94971	.95007 .94140	1.9647 1.8491	.48090 .48062
Errors High Low	LC Pass 2.4000 1.6000	NOCHECK	NOCHECK	NOCHECK	NOCHECK	LC Pass 2.4000 1.6000	LC Pass .60000 .40000
Elem Units Avge SDev %RSD	ZN ppm .48566 .00319 .65594						
#1 #2 #3 #4	.48486 .48332 .49034 .48411						
Errors High Low	LC Pass .60000 .40000						

Operator: MTW Sample Name: DD9PJ Method: QUANMET

Run Time: 05/19/00 10:30:18

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode. Co.							
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00113	.03068	.01065	.17232	.00489	00003	1.0576
SDev	.00259	.00750	.04729	.00107	.00034	.00008	.0108
%RSD	228.12	24.452	444.13	.62239	6.9389	284.83	1.0200
#1	.00454	.03430	.01757	.17354	.00483	00010	1.0551
#2	.00151	.01964	.00975	.17143	.00483	00010	1.0449
#3	00000	.03618	05000	.17291	.00453	.00004	1.0594
#4	00152	.03258	.06527	.17142	.00535	.00004	1.0709
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_ppm .47377 .24673 52.077	LI
Units	ppm	ppm	ppm	ppm	ppm		ppm
Avge	00059	00181	00099	.71372	.04952		.01288
SDev	.00300	.00182	.00349	.00641	.00186		.00058
%RSD	506.66	100.22	351.22	.89842	3.7545		4.4901
#1	.00081	00145	.00257	.70703	.05128	.69840	.01213
#2	L00502	00434	00370	.72087	.04917	.20013	.01275
#3	.00156	00146	00427	.71718	.05058	.33082	.01345
#4	.00028	00000	.00143	.70980	.04706	.66573	.01319
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.10292	.00415	.00373	.80271	.00808	.15791	.01052
SDev	.00874	.00052	.00406	.03060	.00632	.03238	.01053
%RSD	8.4893	12.457	108.66	3.8120	78.200	20.509	100.07
#1	.10568	.00389	.00870	.75792	.00911	.20033	.02632
#2	.09463	.00492	.00373	.80844	.01481	.16169	.00530
#3	.11397	.00389	.00373	.82415	00046	.14635	.00520
#4	.09739	.00389	00123	.82033	.00888	.12328	.00527
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01610	13.295	.02489	.00669	00009	.01402	00251
SDev	.03222	.180	.01231	.00016	.00044	.03459	.00008
%RSD	200.16	1.3562	49.469	2.3327	503.32	246.76	3.1729
#1	00877	13.038	.00737	.00682	.00052	00048	00240
#2	.05888	13.326	.03624	.00655	00052	.02273	00251

Analysis	Report			05/19,	/00 10:33::	23 AM	657 921 page 2
#3 #4	00877 .02306	13.458 13.356	.02797 .02797	.00682 .00655	00017 00017	02363 .05745	00251 00260
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .10978 .00350 3.1919						
#1 #2 #3 #4	.10591 .10772 .11262 .11287						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW

Method: QUANMET Sample Name: DD9PJP5 Operat Run Time: 05/19/00 10:33:26 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

noue. ou							
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00001	.00594	00897	.00754	.00121	.00001	.21010
SDev	.00214	.00683	.02906	.00974	.00047	.00007	.00223
%RSD	24410.	114.98	324.11	129.23	38.857	1202.1	1.0602
#1	00303	.00130	.01391	.02198	.00072	.00004	.20725
#2	.00150	00050	00990	.00293	.00093	00010	.21268
#3	.00153	.00871	.00980	.00441	.00175	.00003	.21039
#4	00004	.01428	04967	.00082	.00144	.00006	.21009
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00176	00217	00228	.14288	.01071	09598	00060
SDev	.00177	.00251	.00248	.00075	.00225	.22690	.00020
%RSD	100.73	115.76	109.07	.52716	21.011	236.40	32.729
#1	00280	00434	00370	.14380	.00843	03676	00083
#2	.00016	00290	.00143	.14288	.00913	06126	00067
#3	00366	.00146	00314	.14288	.01265	.12661	00049
#4	00074	00290	00370	.14196	.01264	41251	00039
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00760	.00000	.00124	.16494	00758	.04424	00533
SDev	.01286	.00052	.00287	.00331	.01052	.00742	.02280
%RSD	169.26	15291.	230.62	2.0090	138.68	16.771	428.20
#1	.02003	00025	.00373	.16695	00781	.03840	00529
#2	00484	00025	.00373	.16525	01596	.03840	03703
#3	.01727	00026	00124	.16016	.00731	.04627	.00535
#4	00207	.00078	00124	.16738	01387	.05387	.01567
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	02085	2.8533	00726	.00119	00009	03482	00224
SDev	.03484	.0783	.01939	.00016	.00052	.05007	.00426
%RSD	167.05	2.7431	266.99	13.665	600.00	143.80	190.69
#1	.01097	2.9524	02795	.00104	00052	.03469	00122
#2	.00700	2.8683	01776	.00115	00052	05795	00250

STL Pittsburgh

page 1

Analysis	Report			05/19,	657 923 page 2		
#3 #4	04473 05666	2.8263 2.7661	.01554 .00111	.00115 .00142	.00052 .00017	08122 03479	.00255 00777
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .02442 .00111 4.5358						
#1 #2 #3 #4	.02296 .02457 .02451 .02565						-
Errors High Low	LC Pass 100.00 02000						

Method: QUANMET Sample Name: DD9PJS Operate Run Time: 05/19/00 10:36:34 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E Operator: MTW

Mode.		<del></del>					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04912	1.9136	1.9412	1.1263	1.9181	.04885	.93042
SDev	.00191	.0100	.0652	.0055	.0122	.00021	.00735
%RSD	3.8811	.52381	3.3592	.49033	.63703	.42764	.79006
#1	.05101	1.9274	2.0107	1.1277	1.9197	.04875	.93213
#2	.04647	1.9128	1.9830	1.1277	1.9278	.04904	.92010
#3	.04949	1.9035	1.8837	1.1184	1.9005	.04861	.93751
#4	.04952	1.9107	1.8875	1.1313	1.9246	.04902	.93193
Error High Low	2.0000 01000	LC Pass 600.00 20000	LC Pass 100.00 30000	LC Pass 100.00 20000	LC Pass 100.00 20000	LC Pass 15.000 00500	LC Pass 600.00 -5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04716	.48566	.19389	.92799	1.0714	.37779	.99055
SDev	.00142	.00290	.00313	.00546	.0037	.31477	.02394
%RSD	3.0005	.59665	1.6147	.58789	.34911	83.318	2.4164
#1	.04720	.49001	.19175	.92730	1.0717	.39617	.99667
#2	.04521	.48421	.19119	.93376	1.0668	.04493	1.0105
#3	.04766	.48422	.19460	.92083	1.0710	.79642	.95578
#4	.04856	.48420	.19802	.93007	1.0759	.27364	.99929
Erroi	rs LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem Unit: Avge SDev %RSD	s ppm .06838 .01505	MN ppm .49383 .00259 .52351	MO ppm .97371 .00745 .76523	NA ppm .77108 .01068 1.3856	NI ppm .50546 .01316 2.6045	PB ppm .62547 .02311 3.6952	SB ppm .48236 .03036 6.2930
#1	.06700	.49409	.97744	.77575	.48751	.60630	.47415
#2	.05319	.49098	.96253	.77406	.51913	.65243	.50636
#3	.08910	.49306	.97744	.75537	.50664	.63702	.50605
#4	.06424	.49719	.97744	.77915	.50854	.60611	.44287
Erro		LC Pass					
High		100.00	50.000	400.00	100.00	100.00	100.00
Low		01500	04000	-5.0000	04000	10000	06000
Elem Unit Avge SDev %RSD	s ppm 1.9450 .0444	SI ppm 13.166 .147 1.1195	SN ppm 1.9423 .0476 2.4529	SR ppm .97430 .00461 .47299	TI ppm .95241 .00199 .20892	TL ppm 1.9327 .1136 5.8769	V_ ppm .48193 .00264 .54675
#1	1.9251	13.170	2.0090	.97537	.95215	1.8835	.48071
#2	1.9569	13.339	1.8974	.97769	.95354	1.9531	.48042
							712

Analysis	Report			05/19/00 10:39:39 AM 657 page 52				
#3 #4	1.8973 2.0007	13.176 12.978	1.9388 1.9240	.96754 .97661	.94972 .95423	1.8141 2.0803	.48071 .48588	
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	100.00	
Elem Units Avge SDev %RSD	ZN ppm .57734 .00307 .53161							
#1 #2 #3 #4	.58001 .57314 .57918 .57702							
Errors High Low	LC Pass 100.00 02000							

Analysis Report

Operator: MTW

Method: QUANMET Sample Name: DD9PJD Operat Run Time: 05/19/00 10:39:42 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CO	MC Corr.	Factor: 1					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04880	1.9556	2.0281	1.1493	1.9491	.04964	1.0304
SDev	.00087	.0096	.0705	.0208	.0052	.00008	.0049
%RSD	1.7901	.49333	3.4743	1.8071	.26607	.16471	.47152
#1	.04955	1.9456	1.9268	1.1313	1.9498	.04957	1.0352
#2	.04955	1.9491	2.0420	1.1673	1.9416	.04971	1.0242
#3	.04804	1.9639	2.0538	1.1313	1.9527	.04957	1.0332
#4	.04804	1.9639	2.0896	1.1673	1.9524	.04971	1.0288
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04762	.49902	.19831	.96009	1.0979	.34920	1.0013
SDev	.00469	.00477	.00177	.00350	.0034	.17575	.0095
%RSD	9.8450	.95676	.89292	.36444	.30643	50.329	.94879
#1	.05275	.50010	.20087	.95778	1.1027	.60855	.99905
#2	.04324	.49288	.19688	.95778	1.0949	.24914	.98936
#3	.05041	.50445	.19745	.96517	1.0978	.23280	1.0117
#4	.04407	.49867	.19802	.95963	1.0964	.30632	1.0051
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.07529	.50262	.99731	.80112	.51762	.65663	.47697
SDev	.00504	.00176	.00405	.01121	.00781	.02053	.02769
%RSD	6.7000	.35101	.40663	1.3990	1.5083	3.1262	5.8060
#1	.08082	.50443	.99234	.79486	.52874	.66819	.45323
#2	.06976	.50030	.99731	.78934	.51119	.62946	.45334
#3	.07253	.50237	1.0023	.81439	.51342	.67606	.49527
#4	.07805	.50340	.99731	.80590	.51713	.65280	.50602
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.9481	11.622	1.9960	.99017	.96689	1.9324	.49142
SDev	.0248	.020	.0619	.00280	.00131	.0555	.00008
%RSD	1.2739	.17078	3.1004	.28284	.13579	2.8700	.01539
#1	1.9213	11.633	2.0585	.98985	.96706	1.8599	.49132
#2	1.9809	11.603	1.9302	.98633	.96 <b>4</b> 98	1.9758	.49141

Analysis	Report			05/19,	657 927 page 2		
#3 #4	1.9491 1.9411	11.645 11.609	2.0380 1.9574	.99239 .99212	.96776 .96776	1.9757 1.9179	.49151 .49142
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000		LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	
Elem Units Avge SDev %RSD	ZN ppm .59545 .00254 .42724						
#1 #2 #3 #4	.59594 .59172 .59706 .59708						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW

Sample Name: DD5V9 Method: QUANMET

Run Time: 05/19/00 10:42:49
Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CO	MC COLL.	ractor: I					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00266	.06316	.01828	.05284	.02413	00000	.69764
SDev	.00077	.00479	.01755	.00148	.00011	.00007	.00399
%RSD	28.776	7.5797	95.988	2.8023	.43887	1579.4	.57240
#1	.00305	.06544	.03712	.05269	.02408	.00003	.69407
#2	.00305	.06365	.02918	.05121	.02408	00011	.69661
#3	.00151	.05632	.00150	.05268	.02429	.00004	.69651
#4	.00305	.06723	.00531	.05480	.02408	.00003	.70337
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00003	00001	.00143	1.1271	.02863	.93937	00280
SDev	.00151	.00167	.00140	.0070	.00176	.26058	.00264
%RSD	5229.7	12588.	97.963	.61673	6.1386	27.740	94.514
#1	00170	00146	.00142	1.1299	.02810	.89444	00189
#2	.00028	.00143	00029	1.1290	.03091	1.0006	00111
#3	.00193	.00143	.00143	1.1327	.02669	.61672	00674
#4	00040	00146	.00313	1.1170	.02880	1.2457	00145
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04559	.00182	.00621	2.2368	00531	.04430	.02090
SDev	.00993	.00084	.00287	.0274	.00602	.00387	.01365
%RSD	21.782	46.434	46.154	1.2244	113.46	8.7272	65.294
#1	.03937	.00182	.00373	2.2519	.00355	.03850	.03682
#2	.04766	.00182	.00373	2.2629	00780	.04626	.01569
#3	.03661	.00078	.00870	2.2319	00988	.04625	.00504
#4	.05871	.00285	.00870	2.2005	00709	.04619	.02605
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem Units Avge SDev %RSD	SE ppm .02199 .03123 142.01	SI ppm H73.297 2.311 3.1523	SN ppm .03476 .02173 62.521	SR ppm .00412 .00000	TI ppm .00061 .00052 85.714	TL ppm .01707 .06307 369.41	V_ ppm .00173 .00190 109.69
#1	02477	H76.033	.02397	.00412	.00017	03503	.00265
#2	.03892	H74.273	.01559		.00087	.02282	.00265

Analysis	Report			05/19,	/00 10:45:	55 AM	657 929 page 2
#3 #4	.03492 .03891	H72.026 H70.855	.03417 .06530	.00412 .00412	.00017 .00121	02343 .10393	00112 .00275
Errors High Low	LC Pass 100.00 25000	LC High 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .17787 .00227 1.2752						
#1 #2 #3 #4	.17480 .17968 .17751 .17949						
Errors High Low	LC Pass 100.00 02000						

657 930 05/19/00 10:49:07 AM page 1 Analysis Report

Operator: MTW

Sample Name: DD9NN Method: QUANMET

Run Time: 05/19/00 10:46:02

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: co.							
Elem Units Avge SDev %RSD	AG ppm .00114 .00076 66.569	AL ppm .07925 .00435 5.4920	AS ppm .03109 .02481 79.792	B_ ppm .22842 .00179 .78560	BA ppm .00329 .00000	BE ppm 00003 .00008 284.72	CA ppm .50051 .00281 .56045
#1	.00151	.07845	.05693	.23111	.00329	.00004	.49893
#2	.00152	.08563	.03699	.22752	.00329	00010	.49928
#3	.00151	.07646	00267	.22751	.00329	.00004	.49913
#4	.00000	.07647	.03310	.22752	.00329	00010	.50471
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00207	00182	.00128	1.9964	.06234	4.2741	00170
SDev	.00127	.00182	.00126	.0106	.00120	.1844	.00073
%RSD	61.061	100.11	98.356	.53063	1.9240	4.3142	43.132
#1	.00360	00001	.00085	1.9978	.06111	4.3742	00145
#2	.00249	00146	.00200	1.9830	.06392	4.2762	00233
#3	.00064	00145	00028	1.9960	.06182	4.0148	00078
#4	.00157	00435	.00257	2.0089	.06252	4.4314	00225
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.06009	.00312	.00373	2.1482	.02787	.72811	00792
SDev	.01070	.00052	.00702	.0344	.01044	.01833	.01008
%RSD	17.807	16.615	188.09	1.6026	37.443	2.5169	127.28
#1	.05871	.00285	00620	2.1470	.01744	.71659	01582
#2	.07529	.00389	.00870	2.1118	.04070	.70885	00523
#3	.05042	.00286	.00870	2.1394	.02165	.74740	.00521
#4	.05595	.00286	.00373	2.1946	.03170	.73959	01583
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	02266	H27.160	.09670	.00277	.00026	.07178	00250
SDev	.02630	.408	.01494	.00000	.00052	.05933	.00014
%RSD	116.10	1.5020	15.454	.00000	200.00	82.651	5.4221
#1	05649	H27.695	.08373	.00277	.00052	01218	00270
#2	.00720	H27.262	.11068	.00277	.00052	.12677	00241

Analysis	Report			05/19,	/00 10:49:0		57 931 page 2
#3 #4	02465 01669		.08383 .10856	.00277 .00277	.00052 00052	.08046 .09206	00241 00250
Errors High Low	LC Pass 100.00 25000	LC High 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .21352 .00110 .51588						
#1 #2 #3 #4	.21355 .21219 .21489 .21345						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW

Analysis Report

Method: QUANMET Sample Name: CCV2-4 Run Time: 05/19/00 10:49:15

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.98043	48.201	5.0265	4.8616	4.7074	4.8848	51.380
SDev	.00388	.226	.0635	.0371	.0284	.0199	.123
%RSD	.39582	.46797	1.2626	.76268	.60248	.40750	.23919
#1	.97995	48.115	5.1111	4.8318	4.6950	4.8702	51.291
#2	.98473	48.521	5.0391	4.9150	4.7490	4.9102	51.449
#3	.98161	48.171	4.9828	4.8572	4.6998	4.8912	51.517
#4	.97542	47.997	4.9730	4.8424	4.6858	4.8678	51.263
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	55.000	5.5000	5.5000	5.5000	5.5000	55.000
Low	.90000	45.000	4.5000	4.5000	4.5000	4.5000	45.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	5.0791	4.9432	4.9195	4.7170	50.880	47.695	4.5631
SDev	.0169	.0213	.0110	.0291	.162	.646	.0630
%RSD	.33231	.43151	.22395	.61595	.31808	1.3548	1.3809
#1	5.0730	4.9277	4.9096	4.7024	50.783	47.659	4.5648
#2	5.1042	4.9681	4.9267	4.7597	51.071	48.533	4.6515
#3	5.0716	4.9538	4.9312	4.7098	50.951	47.634	4.5232
#4	5.0676	4.9234	4.9107	4.6959	50.714	46.956	4.5131
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.5000	5.5000	5.5000	5.5000	55.000	55.000	5.5000
Low	4.5000	4.5000	4.5000	4.5000	45.000	45.000	4.5000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	48.732	4.9025	4.9089	47.453	4.9289	4.9365	4.9752
SDev	.247	.0167	.0091	.576	.0590	.0383	.0437
%RSD	.50754	.34041	.18493	1.2131	1.1970	.77595	.87907
#1	48.659	4.8916	4.8989	47.453	4.8605	4.9862	4.9616
#2	49.087	4.9206	4.9188	48.266	4.9825	4.9408	5.0048
#3	48.670	4.9123	4.9039	47.096	4.9736	4.8944	5.0150
#4	48.512	4.8854	4.9138	46.999	4.8991	4.9244	4.9196
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	55.000	5.5000	5.5000	55.000	5.5000	5.5000	5.5000
Low	45.000	4.5000	4.5000	45.000	4.5000	4.5000	4.5000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	5.0233	H20.283	5.0106	4.7698	4.7940	9.6736	4.8956
SDev	.0700	.862	.0414	.0254	.0170	.0955	.0161
%RSD	1.3928	4.2508	.82584	.53242	.35545	.98688	.32789
#1	5.1036	H20.831	4.9717	4.7602	4.7829	9.5503	4.8799
#2	5.0091	H21.084	4.9863	4.8064	4.8152	9.7206	4.9151

Errors LC Pass High 5.5000 Low 4.5000

#4

4.8855

Operator: MTW Method: QUANMET Sample Name: CCB4

Run Time: 05/19/00 10:52:27

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CO	NC Corr.	Factor: 1					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00150	.01683	.02174	00384	.00070	.00054	.00574
SDev	.00247	.00462	.01807	.00835	.00034	.00027	.00497
%RSD	164.34	27.414	83.119	217.38	49.238	50.468	86.682
#1	.00150	.01591	.04559	01209	.00072	.00032	.00215
#2	00152	.01226	.00589	.00334	.00020	.00032	.00093
#3	.00453	.01590	.02572	00998	.00093	.00061	.01113
#4	.00151	.02326	.00976	.00336	.00093	.00089	.00875
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR ppm00242 .00126 52.338	CU	FE	K_	LI
Units	ppm	ppm		ppm	ppm	ppm	ppm
Avge	00069	.00000		.00046	.01019	.14907	00141
SDev	.00175	.00265		.00046	.00475	.09465	.00067
%RSD	254.62	355130.		100.23	46.613	63.488	47.864
#1	.00007	00290	00199	00023	.00773	.04493	00091
#2	.00093	00145	00313	.00069	.00492	.09394	00207
#3	00062	.00290	00085	.00070	.01265	.24097	00075
#4	00313	.00146	00370	.00070	.01546	.21646	00189
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01036	00025	.01739	.02080	00630	.01925	00540
SDev	.00731	.00000	.00476	.00541	.00541	.01597	.01485
%RSD	70.553	1.0757	27.356	25.995	85.894	83.007	274.81
#1	.00898	00025	.01366	.02473	00796	.03843	00538
#2	.00069	00025	.01366	.01284	01323	.01537	.00506
#3	.01727	00025	.02360	.02218	00182	.00007	.00512
#4	.01451	00025	.01863	.02346	00219	.02310	02642
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.02193	H3.2152	00468	.00073	.00043	02614	00225
SDev	.04227	.3363	.01226	.00023	.00059	.05781	.00009
%RSD	192.80	10.458	262.05	32.248	136.63	221.10	4.1473
#1	.05873	Н3.6372	01968	.00050	.00017	01162	00232
#2	.05474	Н3.3068	.00727	.00061	00017	02318	00232

Analysis Report			05/19/00 10:55:33 AM			page 2	
#3 #4	.00303 02880	H3.0605 H2.8563	.00313 00943	.00077 .00104	.00121	10436 .03458	00212 00223
Errors High Low	LC Pass .25000 25000	LC High .50000 50000	LC Pass .10000 10000	LC Pass .05000 05000	LC Pass .05000 05000	LC Pass .30000 30000	LC Pass .05000 05000
Elem Units Avge SDev %RSD	ZN ppm 00098 .00082 83.913						
#1 #2 #3 #4	00117 .00019 00175 00120						
Errors High Low	LC Pass .02000 02000						

Operator: MTW Method: QUANMET Sample Name: DDCAEB

Run Time: 05/19/00 10:59:12

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: co	ac corr.	140001.					
Elem	AG	AL	AS	B_ /	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00151	.00823	00304	01411	.00023	.00007	.04425
SDev	.00125	.00642	.01361	.00385	.00024	.00007	.00218
%RSD	82.602	78.005	448.19	27.287	103.76	99.305	4.9366
#1 #2 #3 #4	00001 .00304 .00150	.00504 .01597 .01058 .00131	00199 02204 .00193 .00996	01358 00998 01929 01358	.00042 00010 .00042 .00020	.00004 .00003 .00018 .00004	.04516 .04688 .04206 .04292
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm /	ppm
Avge	00077	00036	00171	.00023	.00702	.17971	00156
SDev	.00076	.00072	.00118	.00053	.00057	.25567	.00031
%RSD	97.989	200.81	69.382	230.54	8.1650	142.27	19.562
#1	.00009	.00000	00199	00023	.00702	04493	00171
#2	00143	.00000	00142	.00069	.00773	.46152	00119
#3	00036	.00000	00028	00023	.00703	.33082	00145
#4	00140	00145	00313	.00069	.00632	02859	00189
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA /	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00138	00051	.00373	.02420	00381	01157	.00787
SDev	.00794	.00052	.00406	.00568	.00677	.01602	.01007
%RSD	574.46	100.82	108.83	23.467	177.65	138.55	128.00
#1	00760	00026	.00373	.01666	00188	00769	.01580
#2	00760	00026	.00373	.02600	.00446	01543	00528
#3	.00069	00129	00124	.02388	01151	03081	.01573
#4	.00898	00025	.00869	.03025	00631	.00767	.00523
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
Low	-5.0000	01500	04000	-5.0000	~.04000	10000	06000
Elem Units Avge SDev %RSD	SE ppm .00600 .03304 551.05	SI ppm H1.1383 .0000	SN ppm 00720 .02031 281.96	SR ppm .00032 .00034 104.60	TI ppm .00000 .00035 414e6	TL ppm .01152 .03407 295.89	V_ ppm 00122 .00258 211.37
#1 #2	03679 .00301	H1.1383 H1.1383	01130 .02170	.00012 00004	00052 .00017	02322 .03464	00251 .00265

Analysis	Report			05/19,	/00 11:02:		57 937 page 2
#3 #4	.01495 .04280	H1.1383 H1.1383	01337 02584	.00061 .00061	.00017 .00017	.04626 01162	00260 00242
Errors High Low	LC Pass .25000 25000	LC High .50000 50000	LC Pass .10000 10000	LC Pass .05000 ~.05000	LC Pass .05000 05000	LC Pass .30000 30000	LC Pass .05000 05000
Elem Units Avge SDev %RSD	ZN ppm .00965 .00146 15.125						
#1 #2 #3 #4	.00879 .00958 .00850 .01173						
Errors High Low	LC Pass .02000 02000						

Operator: MTW

Analysis Report

Method: QUANMET Sample Name: DDCAEC

Run Time: 05/19/00 11:02:25

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: Co	NC COLL.	raccor. x					
Elem Units Avge SDev	AG ppm .04987 .00145	AL ppm 1.9099 .0106	AS ppm 1.9582 .0372 1.9016	B_ / ppm .90244 .00100 .11071	BA ppm 1.8438 .0042 .22612	BE ppm .04907 .00013 .26955	CA ppm / 52.333 .146 .27905
%RSD #1 #2 #3 #4	2.9172 .05100 .04798 .05103 .04949	1.9090 1.9217 1.9127 1.8961	1.9354 2.0068 1.9670 1.9235	.90248 .90101 .90313 .90312	1.8393 1.8430 1.8494 1.8434	.04903 .04918 .04916 .04889	52.551 52.242 52.283 52.256
Errors	LC Pass	LC Pass	LC Pass	NOCHECK	LC Pass	LC Pass	LC Pass
High	.06000	2.4000	2.4000		2.4000	.06000	60.000
Low	.04000	1.6000	1.6000		1.6000	.04000	40.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm /	ppm
Avge	.04594	.49150	.19389	.23528	1.0401	47.299	.88571
SDev	.00460	.00117	.00320	.00053	.0043	.584	.00901
%RSD	10.019	.23769	1.6498	.22666	.40964	1.2339	1.0175
#1	.04939	.49149	.19460	.23482	1.0373	47.185	.87976
#2	L.03945	.49007	.19233	.23574	1.0387	48.075	.89639
#3	.04904	.49293	.19802	.23575	1.0464	47.275	.88979
#4	.04589	.49150	.19062	.23482	1.0380	46.662	.87691
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	NOCHECK
High	.06000	.60000	.24000	.30000	1.2000	60.000	
Low	.04000	.40000	.16000	.20000	.80000	40.000	
Elem	MG /	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	49.757	.49228	.98985	46.828	.48785	.50239	.52405
SDev	.154	.00099	.00860	.384	.00286	.02226	.01003
%RSD	.30852	.20049	.86912	.81948	.58536	4.4302	1.9147
#1	49.700	.49305	.98240	46.535	.49003	.52165	.51616
#2	49.692	.49099	.99730	47.335	.49061	.48310	.53718
#3	49.985	.49305	.98240	46.916	.48525	.52169	.51626
#4	49.650	.49202	.99730	46.526	.48554	.48314	.52662
Errors	LC Pass	LC Pass	NOCHECK	LC Pass	LC Pass	LC Pass	LC Pass
High	60.000	.60000		60.000	.60000	.60000	.60000
Low	40.000	.40000		40.000	.40000	.40000	.40000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.9837	11.191	2.0377	.94122	.95302	1.9678	.48224
SDev	.0444	.046	.0316	.00161	.00229	.0884	.00249
%RSD	2.2365	.41483	1.5531	.17063	.24047	4.4919	.51654
#1	1.9887	11.158	2.0588	.94008	.95284	2.0228	.48081
#2	2.0245	11.146	2.0507	.94073	.95146	1.8491	.48109

Analysis	Report			05/19/	MA OE	657 939 page 2	
#3 #4	2.0006 1.9210	11.242 11.218	1.9905 2.0506	.94359 .94046	.95631 .95146	1.9532 2.0459	.48597 .48109
Errors High Low	LC Pass 2.4000 1.6000	NOCHECK	NOCHECK	NOCHECK	NOCHECK	LC Pass 2.4000 1.6000	LC Pass .60000 .40000
Elem Units Avge SDev %RSD	ZN ppm .51680 .00516 .99835						
#1 #2 #3 #4	.52104 .52128 .51379 .51108						
Errors High Low	LC Pass .60000 .40000						

05/19/00 11:08:38 AM page 1

Operator: MTW Method: QUANMET Sample Name: DD7WE Run Time: 05/19/00 11:05:33

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00616	2.5839	.11348	.17005	.08185	00009	46.383
SDev	.00088	.0142	.06044	.01503	.00043	.00008	.104
%RSD	14.321	.54878	53.267	8.8353	.52921	87.438	.22338
#1	.00691	2.6036	.06663	.19249	.08190	00002	46.497
#2	.00694	2.5705	.05888	.16456	.08139	00003	46.401
#3	.00540	2.5834	.14628	.16159	.08169	00016	46.388
#4	.00539	2.5780	.18211	.16157	.08242	00016	46.245
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00043	00009	.06297	.04091	3.5237	H1260.8	.00055
SDev	.00330	.00119	.00271	.00046	.0073	13.9	.00029
%RSD	771.29	1325.9	4.3110	1.1272	.20807	1.1006	53.293
#1 #2 #3 #4	00055 .00502 00279 .00003	.00136 00155 00008 00009	.06582 .06011 .06468 .06126	.04068 .04068 .04068 .04160	3.5313 3.5243 3.5257 3.5137	H1260.2 H1250.8 H1251.7 H1280.7	.00057 .00013 .00075
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	3.9779	.39169	.04399	72.875	.00265	.03259	.02295
SDev	.0129	.00060	.00406	.902	.00391	.02555	.01583
%RSD	.32330	.15187	9.2213	1.2378	147.71	78.381	68.982
#1	3.9655	.39221	.04399	72.594	.00726	.03651	.04669
#2	3.9876	.39118	.04399	72.305	00224	.05952	.01503
#3	3.9904	.39118	.04896	72.384	.00218	.03644	.01491
#4	3.9683	.39221	.03902	74.215	.00340	00209	.01516
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00602	17.722	.01671	.21185	.02878	.03805	.02031
SDev	.02068	.052	.02766	.00053	.00035	.05304	.00258
%RSD	343.47	.29588	165.56	.25147	1.2048	139.40	12.681
#1	.00793	17.784	.00802	.21209	.02930	.00034	.01903
#2	03587	17.682	.01822	.21106	.02861	.01197	

Analysis	Report			05/19/00 11:08:38 AM					
#3 #4	.00791 00407	17.748 17.676	.05339 01281	.21214 .21214	.02861 .02861	.02356 .11633	.01912 .01892		
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000		
Elem Units Avge SDev %RSD	ZN ppm .27431 .00475 1.7318								
#1 #2 #3 #4	.27364 .26910 .28062 .27388								
Errors High Low	LC Pass 100.00 02000								

Operator: MTW Sample Name: DD7WEP5 Method: QUANMET

Run Time: 05/19/00 11:08:46

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Mode: co.	NC COII.	raccox.					
Elem Units Avge SDev %RSD	AG ppm .00167 .00214 128.02	AL ppm .51043 .00530 1.0377	AS ppm .03950 .04441 112.42	B_ ppm .00161 .00106 65.706	BA ppm .01636 .00000	BE ppm 00000 .00007 2552.1	CA ppm 8.9320 .0132 .14806
#1	.00319	.51075	01120	.00002	.01636	.00003	8.9238
#2	00136	.50357	.02863	.00213	.01636	00011	8.9263
#3	.00316	.51649	.04448	.00214	.01636	.00004	8.9261
#4	.00168	.51091	.09609	.00214	.01636	.00003	8.9518
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00014	00110	.00969	.00790	.69789	257.27	00125
SDev	.00074	.00139	.00057	.00046	.00337	3.19	.00056
%RSD	509.55	125.81	5.8778	5.8320	.48279	1.2415	44.752
#1	.00114	00292	.00997	.00813	.69754	253.78	00117
#2	00063	00001	.00997	.00813	.69333	260.49	00163
#3	.00013	00002	.00883	.00721	.69965	259.42	00171
#4	00005	00147	.00997	.00813	.70106	255.41	00049
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.77638	.07663	.00383	14.814	00562	.00147	00541
SDev	.01416	.00052	.00702	.176	.00225	.01820	.02276
%RSD	1.8233	.67290	183.42	1.1913	40.078	1238.7	420.92
#1	.76050	.07638	.01376	14.610	00575	.01488	02655
#2	.79089	.07637	.00383	14.974	00259	00038	.02618
#3	.76878	.07740	00114	14.949	00803	.01498	00539
#4	.78536	.07637	00114	14.724	00611	02360	01588
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01151	3.7994	00974	.04283	.00520	02993	.00106
SDev	.03639	.0350	.01614	.00027	.00020	.04226	.00251
%RSD	316.28	.92154	165.68	.63062	3.8490	141.18	236.95
#1	.00156	3.8294	00725	.04243	.00503	07625	.00286
#2	.06521	3.7634	.01162	.04297	.00503	05306	.00138

Analysis	Report			05/19/	657 943 page 2		
#3 #4	00640 01435	3.8294 3.7754	01750 02582	.04297 .04297	.00538 .00538	00677 .01634	00258 .00258
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .05434 .00127 2.3416						
#1 #2 #3 #4	.05451 .05338 .05608 .05340						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW

Analysis Report

Method: QUANMET Sample Name: DD7WES

Run Time: 05/19/00 11:11:54

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Elem Units Avge SDev %RSD	AG ppm .05937 .00213 3.5804	AL ppm 5.5857 .0281 .50224	AS ppm 2.0936 .0382 1.8232	B_ ppm 1.0772 .0197 1.8281	BA ppm 1.8400 .0146 .79224	BE ppm .04632 .00030 .65015	CA ppm 94.613 .067
#1	.06237	5.5638	2.1306	1.0592	1.8254	.04605	94.710
#2	.05787	5.6239	2.0585	1.0967	1.8588	.04674	94.589
#3	.05937	5.5894	2.1224	1.0915	1.8434	.04632	94.557
#4	.05786	5.5656	2.0630	1.0613	1.8324	.04618	94.599
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04515	.47341	.24988	.27132	4.5008	H1234.4	.85768
SDev	.00291	.00495	.00369	.00158	.0122	11.4	.01255
%RSD	6.4442	1.0455	1.4786	.58112	.27044	.91957	1.4634
#1	.04819	.47124	.24532	.27109	4.4926	H1227.8	.84773
#2	.04119	.46833	.25387	.27293	4.5186	H1250.3	.87463
#3	.04550	.47992	.25159	.27202	4.4989	H1234.4	.85973
#4	.04573	.47413	.24874	.26924	4.4933	H1224.9	.84861
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	50.980	.84492	.98167	114.39	.47807	.51465	.53955
SDev	.280	.00230	.01174	1.21	.00608	.02396	.01812
%RSD	.54934	.27177	1.1958	1.0594	1.2721	4.6559	3.3583
#1	50.734	.84207	.97298	113.50	.48547	.54929	.54760
#2	51.341	.84725	.99782	116.05	.47212	.49520	.51563
#3	51.060	.84621	.98292	114.53	.48048	.50327	.53696
#4	50.786	.84414	.97298	113.47	.47421	.51086	.55802
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	2.0083	H27.329	1.9407	1.1035	.94261	1.9196	.48993
SDev	.0304	.097	.0054	.0077	.00429	.0433	.00270
%RSD	1.5145	.35589	.27921	.69748	.45509	2.2544	.55180
#1	2.0013	H27.209	1.9376	1.0955	.93758	1.9197	.48588
#2	2.0531	H27.443	1.9478	1.1129	.94799	1.9311	.49154

Analysis	Report			05/19	657945		
#3 #4	1.9894 1.9894	H27.353 H27.311	1.9355 1.9418	1.1063 1.0993	.94313 .94175	1.9659 1.8618	.49124 .49105
Errors High Low	LC Pass 100.00 25000	LC High 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .74289 .00215 .28878						
#1 #2 #3 #4	.74335 .74554 .74227 .74040						
Errors High Low	LC Pass 100.00 02000						

Method: QUANMET Sample Name: DD7WED Run Time: 05/19/00 11:15:07 Operator: MTW

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

ľ	iode: COr	COPP.	ractor: 1					
	Elem	AG	AL	AS	B_	BA	BE	CA
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	.05817	5.4929	2.0937	1.0824	1.8105	.04598	93.804
	SDev	.00191	.0195	.0630	.0157	.0117	.00029	.354
	%RSD	3.2814	.35450	3.0076	1.4520	.64462	.63125	.37767
	#1	.05629	5.5072	2.0796	1.0949	1.8159	.04591	94.026
	#2	.05778	5.5052	2.0876	1.0801	1.8175	.04606	93.582
	#3	.05777	5.4650	2.1794	1.0610	1.7931	.04564	93.431
	#4	.06083	5.4943	2.0280	1.0935	1.8157	.04633	94.177
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
	Low	01000	20000	30000	20000	20000	00500	-5.0000
	Elem	CD	CO	CR	CU	FE	K_	LI
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	.04334	.47490	.24618	.26762	4.3817	H1214.1	.84760
	SDev	.00168	.00597	.00109	.00158	.0164	13.6	.01218
	%RSD	3.8852	1.2562	.44297	.59116	.37455	1.1223	1.4365
	#1	.04329	.47997	.24646	.26924	4.3907	H1232.4	.86328
	#2	.04220	.46838	.24532	.26738	4.3808	H1213.3	.84784
	#3	.04214	.47130	.24760	.26554	4.3590	H1211.4	.84563
	#4	.04574	.47996	.24532	.26832	4.3963	H1199.5	.83364
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC High	LC Pass
	High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
	Low	00500	05000	01000	02500	10000	-5.0000	05000
	Elem	MG	MN	MO	NA	NI	PB	SB
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	50.507	.83584	.96179	112.61	.47153	.51863	.54759
	SDev	.196	.00368	.00476	1.23	.00586	.03153	.01485
	%RSD	.38885	.44041	.49461	1.0928	1.2430	6.0796	2.7120
	#1	50.521	.83894	.95806	114.19	.46800	.52647	.54757
	#2	50.651	.83377	.96303	112.65	.46757	.53387	.55812
	#3	50.225	.83170	.95806	112.40	.48011	.47229	.52660
	#4	50.631	.83895	.96800	111.20	.47043	.54190	.55808
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
	Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
	Elem	SE	SI	SN	SR	TI	TL	V_
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	1.9720	H26.812	1.9222	1.0863	.92380	1.7761	.48309
	SDev	.0571	.101	.0200	.0056	.00423	.0880	.00300
	%RSD	2.8968	.37811	1.0412	.51723	.45812	4.9525	.62205
	#1	2.0248	H26.896	1.9377	1.0893	.92614	1.8049	.48560
	#2	1.9691	H26.818	1.8943	1.0896	.92545	1.8862	.48053

Analysis	Report			05/19,	657 947 page 2		
#3 #4	2.0008 1.8935	H26.668 H26.866	1.9210 1.9356	1.0779 1.0885	.91747 .92614	1.7242 1.6890	.48044 .48579
Errors High Low	LC Pass 100.00 25000	20.000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000		LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .73362 .00288 .39299						
#1 #2 #3 #4	.73075 .73232 .73395 .73748						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW Method: QUANMET Sample Name: DD7NT Run Time: 05/19/00 11:18:19

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

140ac. 00.	.,,						
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00120	6.7612	.02758	.16649	.00239	.00012	1.9606
SDev	.00076	.0389	.04947	.00000	.00048	.00014	.0164
%RSD	62.973	.57590	179.38	.00091	20.104	113.01	.83634
#1	.00158	6.7851	02726	.16649	.00196	.00030	1.9405
#2	.00158	6.8016	.02823	.16649	.00226	.00001	1.9666
#3	.00158	6.7410	.01681	.16649	.00308	.00015	1.9792
#4	.00007	6.7172	.09252	.16649	.00226	.00001	1.9560
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem .	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00031	00038	.06824	.02125	.12754	9.2936	.01220
SDev	.00142	.00072	.00164	.00046	.00135	.3035	.00000
%RSD	460.13	187.73	2.3982	2.1742	1.0555	3.2654	.00000
#1	00097	00002	.06923	.02102	.12719	8.9485	.01220
#2	00065	00002	.06866	.02102	.12860	9.4386	.01220
#3	.00177	00003	.06923	.02194	.12859	9.6347	.01220
#4	00138	00147	.06581	.02101	.12578	9.1527	.01220
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	2.5757	.02521	.01616	S11792	.00088	.00032	.02034
SDev	.0264	.00052	.00641	.00000	.01203	.03417	.01056
%RSD	1.0251	2.0521	39.676	.00000	1362.4	10525.	51.931
#1	2.5702	.02444	.01368	S11792	00923	.04270	.01500
#2	2.5730	.02547	.00871	S11792	.00455	03434	.03619
#3	2.6117	.02548	.02361	S11792	.01631	.01190	.01519
#4	2.5481	.02547	.01865	S11792	00809	01896	.01499
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	01446	1.4883	.03938	.00689	.01344	.10878	.00818
SDev	.02805	.0406	.00997	.00026	.00044	.03599	.00012
%RSD	194.02	2.7250	25.306	3.7546	3.2472	33.087	1.5094
#1	04033	1.5229	.03418	.00655	.01335	.09140	.00813
#2	.00345	1.4628	.05290	.00709	.01405	.16087	.00804

Analysis	Report			05/19/	/00 11:21:2	25 AM	657 949 page 2
#3 #4	.01539 03635	1.5229 1.4448	.03005 .04039	.00709 .00682	.01335 .0130 <b>1</b>	.10297 .07986	.00833
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .02696 .00064 2.3680						
#1 #2 #3 #4	.02705 .02621 .02776 .02681						
Errors High Low	LC Pass 100.00 02000						

Method: QUANMET Sample Name: CCV2-5 Operate Run Time: 05/19/00 11:21:27 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Operator: MTW

Mode: COI	NC Corr.	Factor: 1					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.97243	48.126	4.9489	4.8270	4.7232	4.9037	51.209
SDev	.00130	.184	.0721	.0192	.0202	.0194	.161
%RSD	.13365	.38251	1.4564	.39837	.42835	.39553	.31462
#1	.97411	48.329	4.9853	4.8352	4.7484	4.9269	51.431
#2	.97094	48.186	4.8997	4.8424	4.7283	4.9112	51.183
#3	.97234	48.098	5.0320	4.7990	4.7152	4.8938	51.045
#4	.97235	47.889	4.8787	4.8315	4.7008	4.8828	51.179
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	55.000	5.5000	5.5000	5.5000	5.5000	55.000
Low	.90000	45.000	4.5000	4.5000	4.5000	4.5000	45.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	5.0463	4.9303	4.8989	4.7333	50.692	47.040	4.5214
SDev	.0304	.0210	.0144	.0218	.171	.324	.0350
%RSD	.60330	.42642	.29297	.46003	.33659	.68884	.77384
#1	5.0471	4.9509	4.9192	4.7569	50.923	47.332	4.5310
#2	5.0723	4.9436	4.8976	4.7412	50.706	47.201	4.5210
#3	5.0033	4.9047	4.8925	4.7301	50.530	47.038	4.5589
#4	5.0622	4.9219	4.8862	4.7052	50.606	46.589	L4.4748
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.5000	5.5000	5.5000	5.5000	55.000	55.000	5.5000
Low	4.5000	4.5000	4.5000	4.5000	45.000	45.000	4.5000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	48.792	4.8831	4.8865	48.459	4.9332	4.9188	4.9124
SDev	.199	.0174	.0125	.272	.0375	.0343	.0945
%RSD	.40749	.35600	.25632	.56127	.76058	.69740	1.9236
#1	49.048	4.9061	4.9039	48.583	4.9838	4.9559	4.9941
#2	48.844	4.8843	4.8840	48.532	4.9145	4.8864	4.9942
#3	48.675	4.8647	4.8741	48.662	4.9374	4.8929	4.8254
#4	48.601	4.8771	4.8840	48.059	4.8972	4.9397	4.8359
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	55.000	5.5000	5.5000	55.000	5.5000	5.5000	5.5000
Low	45.000	4.5000	4.5000	45.000	4.5000	4.5000	4.5000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	5.0276	5.3185	4.9689	4.7957	4.7877	9.6379	4.8942
SDev	.0759	.0159	.1056	.0213	.0158	.2150	.0161
%RSD	1.5090	.29881	2.1249	.44345	.33059	2.2303	.32937
#1	4.9568	5.3098	5.0835	4.8218	4.8083	9.4558	4.9161
#2	4.9680	5.3396	4.8704	4.8017	4.7920	9.8171	4.8964
~~~ ~ '	7						C1 E /

pag	ge	2
-----	----	---

4.8820 4.8822

LC Pass 5.5000 4.5000

#3 #4	5.1026 5.0830	5.3214 5.3033	5.0331 4.8884	4.7875 4.7718	4.7767 4.7739	9.8308 9.4478
	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 5.5000 4.5000	LC Pass 11.000 9.0000
Elem Units Avge SDev %RSD	ZN ppm 4.8530 .0051 .10564					
#1 #2 #3 #4	4.8485 4.8592 4.8491 4.8553					
Errors High Low	LC Pass 5.5000 4.5000					

05/19/00 11:27:41 AM page 1

Operator: MTW Method: QUANMET Sample Na Run Time: 05/19/00 11:24:35 Sample Name: CCB5

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: co	WC COII.	100001 -					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00188	.01411	00505	.00441	.00077	.00082	.00843
SDev	.00190	.00578	.02209	.01491	.00038	.00018	.00229
%RSD	101.21	40.980	436.98	338.29	49.332	22.295	27.215
#1	.00453	.01955	00609	.00124	.00093	.00061	.00788
#2	.00150	.01233	.02576	.00758	.00102	.00103	.01149
#3	00001	.00676	02585	01358	.00020	.00089	.00595
#4	.00150	.01778	01403	.02239	.00093	.00075	.00839
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR ppm00071 .00261 368.43	CU	FE	K_	LI
Units	ppm	ppm		ppm	ppm	ppm	ppm
Avge	.00085	00145		.00069	.00738	.22463	00197
SDev	.00206	.00205		.00075	.00091	.17108	.00071
%RSD	243.09	140.90		109.06	12.303	76.161	36.226
#1	00030	00435	.00257	.00069	.00773	.03676	00223
#2	.00393	00001	00028	.00069	.00843	.44518	00285
#3	00006	00145	00142	00023	.00632	.16745	00153
#4	00019	00000	00370	.00162	.00703	.24914	00127
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem Units Avge SDev %RSD	MG ppm .01243 .01180 94.933	MN ppm 00025 .00000	MO ppm .01242 .00248 19.997	NA ppm .06453 .01499 23.238	NI ppm 00374 .00629 168.25	PB ppm 01352 .03468 256.48	SB ppm .01042 .01050 100.80
#1	.01451	00025	.01366	.04978	00845	01553	.00507
#2	.02003	00025	.00869	.08544	.00475	.00770	.00520
#3	00484	00025	.01366	.06082	00858	.01540	.02616
#4	.02003	00025	.01366	.06209	00267	06166	.00523
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	03878	.28680	.00212	.00081	.00043	00005	00234
SDev	.01471	.02703	.01894	.00026	.00044	.04123	.00005
%RSD	37.949	9.4249	894.23	31.914	100.66	83357.	2.0385
#1	02484	.27329	00101	.00061	.00052	.05787	00231
#2	02882	.32735	.02382	.00088		03481	00241

Analysis	alysis Report				05/19/00 11:27:41 AM			
#3 #4	04475 05669	.27329 .27329	.00737 02170	.00061 .00115	00017 .00087	02320 00006	00232 00232	
Errors High Low	LC Pass .25000 25000	LC Pass .50000 50000	LC Pass .10000 10000	LC Pass .05000 05000	LC Pass .05000 05000	LC Pass .30000 30000	LC Pass .05000 05000	
Elem Units Avge SDev %RSD	ZN ppm .00117 .00156 133.45							
#1 #2 #3 #4	.00287 .00013 .00208 00041							
Errors High Low	LC Pass .02000 02000							

Operator: MTW

Method: QUANMET Sample Name: DD7WE/2 K

Run Time: 05/19/00 11:32:26

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CO	MC Corr.	Factor: 1					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00231	1.2760	.02760	.06501	.04115	.00005	23.118
SDev	.00145	.0068	.03496	.01798	.00025	.00007	.071
%RSD	62.507	.53259	126.67	27.653	.61042	143.96	.30523
#1	.00043	1.2695	.02068	.05605	.04136	.00001	23.182
#2	.00194	1.2787	.05241	.05602	.04105	.00001	23.139
#3	.00345	1.2842	.05635	.05600	.04084	.00016	23.017
#4	.00345	1.2714	01905	.09198	.04136	.00001	23.135
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00102	.00032	.03049	.01676	1.7651	633.64	.00009
SDev	.00212	.00217	.00298	.00692	.0086	6.07	.00229
%RSD	207.03	689.24	9.7719	41.312	.48558	.95773	2523.9
#1	.00175	00295	.03447	.02022	1.7772	642.12	00189
#2	00180	.00141	.02821	.02022	1.7646	633.94	.00207
#3	.00326	.00140	.03106	.02022	1.7576	629.15	00189
#4	.00088	.00140	.02821	.00637	1.7611	629.35	.00207
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.9928	.19546	.01889	36.050	00522	.01054	.01807
SDev	.0108	.00084	.00000	.350	.00885	.02228	.03379
%RSD	.54291	.43195	.00679	.96999	169.54	211.28	187.02
#1	1.9983	.19546	.01889	36.529	01830	.01429	.00474
#2	1.9900	.19649	.01889	36.091	00296	01642	.00490
#3	2.0038	.19442	.01889	35.773	.00032	.00669	00559
#4	1.9789	.19546	.01889	35.807	.00007	.03761	.06821
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.02038	8.4548	01256	.10680	.01422	.00887	.00816
SDev	.02983	.0591	.01188	.00079	.00035	.05781	.00001
%RSD	146.38	.69853	94.603	.74134	2.4390	651.87	.07246
#1	01042	8.5389	01521	.10790	.01370	02887	.00817
#2	.06117	8.4428	.00341	.10682	.01439	.06387	
-							616

Analysis	Report			05/19/00 11:35:32 AM			657 955 page 2
#3 #4	.01737 .01340	8.4008 8.4368	01319 02526	.10628 .10617	.01439 .01439	05186 .05233	.00816 .00815
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000		LC Pass 50.000 05000	LC Pass 100.00 -2.0000	
Elem Units Avge SDev %RSD	ZN ppm .13686 .00138 1.0100						
#1 #2 #3 #4	.13791 .13596 .13540 .13816						
Errors High Low	LC Pass 100.00 02000						

Method: QUANMET Sample Name: DD7WEP10 K . Run Time: 05/19/00 11:35:35 Operator: MTW

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CO	NC COII.	raccor. I					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00159	.27400	01567	.00610	.00874	.00007	4.7229
SDev	.00175	.00489	.03104	.00843	.00025	.00007	.0139
%RSD	110.55	1.7848	198.10	138.13	2.8725	99.498	.29441
#1	.00309	.27080	02059	.00361	.00843	.00004	4.7384
#2	.00007	.27822	05244	00061	.00895	.00018	4.7307
#3	.00312	.27814	01279	.00298	.00895	.00003	4.7126
#4	.00007	.26884	.02314	.01844	.00865	.00004	4.7098
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00008	00037	.00613	.00372	.36791	137.72	00194
SDev	.00334	.00274	.00303	.00088	.00265	1.27	.00082
%RSD	4014.0	738.66	49.390	23.754	.72025	.92322	42.451
#1	00277	.00145	.00485	.00257	.36598	136.80	00277
#2	.00483	.00143	.00370	.00442	.37160	138.32	00189
#3	00018	00001	.01054	.00441	.36598	136.55	00083
#4	00154	00435	.00542	.00349	.36809	139.23	00225
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.42480	.04065	.00502	7.7977	00509	.00168	01330
SDev	.01934	.00060	.00745	.0976	.00519	.02307	.00525
%RSD	4.5528	1.4668	148.35	1.2513	102.04	1376.9	39.481
#1	.41789	.04116	.00875	7.7281	00011	.03064	00543
#2	.43447	.04116	00119	7.8067	01174	00792	01592
#3	.44552	.04013	00119	7.7235	00197	.00746	01587
#4	.40132	.04013	.01371	7.9323	00653	02348	01601
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	01557	1.8441	00052	.02292	.00243	03830	00118
SDev	.03133	.0384	.01817	.00030	.00066	.04013	.00251
%RSD	201.22	2.0847	3505.9	1.3170	27.355	104.79	212.46
#1	.02024	1.8831	.02590	.02271	.00260	02670	00240
#2	05535	1.8711	01139	.02330	.00191	09624	00259

Errors LC Pass High 100.00 Low -.02000

Operator: MTW Method: QUANMET Sample Name: DD7WES/2 K Run Time: 05/19/00 11:38:42

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: COI	MC Corr.	ractor: 1					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.02968	2.7582	1.0049	.53666	.92913	.02343	46.948
SDev	.00195	.0198	.0545	.00242	.00556	.00014	.092
%RSD	6.5769	.71936	5.4196	.45058	.59843	.57829	.19666
#1	.02892	2.7559	1.0287	.53381	.92869	.02360	46.905
#2	.03043	2.7376	1.0607	.53590	.92148	.02332	46.838
#3	.02741	2.7542	.99695	.53953	.93229	.02346	47.027
#4	.03195	2.7853	.93312	.53741	.93404	.02332	47.021
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.02276	.23670	.12580	.13266	2.2570	624.03	.42936
SDev	.00180	.00321	.00205	.00119	.0116	9.22	.00563
%RSD	7.9071	1.3553	1.6280	.89953	.51249	1.4779	1.3105
#1	.02473	.23488	.12622	.13404	2.2551	630.34	.43041
#2	.02168	.23345	.12850	.13127	2.2417	610.50	.42170
#3	.02377	.24068	.12451	.13313	2.2684	625.84	.43007
#4	.02085	.23779	.12394	.13220	2.2628	629.43	.43525
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	25.596	.42298	.49208	57.032	.23971	.24564	.16304
SDev	.139	.00169	.00248	.831	.01196	.04643	.10156
%RSD	.54197	.39929	.50466	1.4562	4.9907	18.901	62.288
#1	25.583	.42298	.49580	57.568	.25113	.26692	.26583
#2	25.406	.42091	.49084	55.803	.23086	.28997	.23408
#3	25.677	.42298	.49084	57.257	.22800	.18204	.06555
#4	25.718	.42504	.49084	57.502	.24884	.24363	.08672
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.95343	13.257	.99105	.55829	.47053	.92501	.24434
SDev	.06136	.080	.01636	.00306	.00156	.08067	.00005
%RSD	6.4352	.59996	1.6506	.54792	.33167	8.7213	.01993
#1	.99422	13.233	.99619	.55808	.47044	.95978	.24442
#2	.92253	13.155	.97948	.55403	.46836	.83254	

Analysis	Report			05/19,	/00 11:41:4	48 AM	657 959 page 2
#3 #4	1.0142 .88281	13.317 13.323	.97662 1.0119	.56051 .56051	.47148 .47183	1.0175 .89021	.24432 .24432
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	
Elem Units Avge SDev %RSD	ZN ppm .36376 .00188 .51688						
#1 #2 #3 #4	.36354 .36251 .36251 .36648						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW Sample Name: DD7WED/2 K Method: QUANMET

Run Time: 05/19/00 11:41:50

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CO	WC Corr.	ractor: 1					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.02816	2.7419	1.0190	.53917	.92527	.02350	46.674
SDev	.00151	.0150	.0592	.01226	.00449	.00018	.218
%RSD	5.3760	.54565	5.8079	2.2733	.48563	.75909	.46811
#1	.02740	2.7323	1.1005	.52501	.92170	.02346	46.639
#2	.02740	2.7285	.97737	.55488	.92303	.02332	46.435
#3	.03043	2.7450	.97325	.53733	.92457	.02346	46.657
#4	.02740	2.7617	1.0248	.53944	.93177	.02375	46.965
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.02318	.23924	.12323	.13243	2.2124	621.07	.42835
SDev	.00351	.00334	.00220	.00158	.0089	5.82	.00338
%RSD	15.133	1.3959	1.7861	1.1908	.40296	.93655	.78827
#1	.02168	.24215	.12109	.13035	2.2003	614.96	.42372
#2	.01940	.23636	.12451	.13312	2.2122	617.88	.42795
#3	.02401	.23634	.12166	.13219	2.2157	623.40	.43077
#4	.02762	.24212	.12565	.13405	2.2213	628.04	.43095
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	25.517	.42012	.47966	56.920	.23832	.25735	.20801
SDev	.150	.00130	.00476	.507	.00737	.03115	.09592
%RSD	.58646	.30942	.99150	.89029	3.0924	12.105	46.110
#1	25.450	.41986	.47593	56.400	.24855	.21293	.06589
#2	25.406	.41987	.48586	56.643	.23550	.25924	.23430
#3	25.475	.41883	.48090	57.088	.23801	.27468	.26603
#4	25.737	.42193	.47593	57.548	.23123	.28254	.26583
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.94929	13.110	.94989	.55568	.46558	.87622	.24410
SDev	.06231	.079	.02975	.00271	.00165	.09140	.00009
%RSD	6.5635	.60032	3.1319	.48838	.35463	10.431	.03796
#1	.90646	13.059	.96629	.55376	.46454	.97186	.24402
#2	.92640	13.071	.93604	.55457	.46454	.89072	

100.00

-.02000

High

Low

Operator: MTW

Sample Name: CCV2-6 Method: QUANMET

Run Time: 05/19/00 11:44:58

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CC	onc Corr.	ractor: 1					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.97724	48.492	5.0282	4.8297	4.7509	4.9577	51.715
SDev	.00218	.099	.0905	.0207	.0176	.0192	.129
%RSD	.22333	.20465	1.7989	.42927	.37125	.38709	.24971
#1	.97414	48.450	5.0636	4.8319	4.7431	4.9427	51.610
#2	.97873	48.430	5.1073	4.7997	4.7449	4.9465	51.816
#3	.97881	48.640	4.8986	4.8441	4.7770	4.9851	51.596
#4	.97729	48.448	5.0433	4.8429	4.7384	4.9564	51.836
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	1.1000	55.000	5.5000	5.5000	5.5000	5.5000	55.000
Low	.90000	45.000	4.5000	4.5000	4.5000	4.5000	45.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	5.1080	4.9663	4.9376	4.7759	50.982	48.445	4.6146
SDev	.0117	.0128	.0053	.0251	.083	.487	.0287
%RSD	.22873	.25793	.10633	.52656	.16291	1.0057	.62132
#1	5.1009	4.9508	4.9380	4.7606	50.876	48.933	4.6355
#2	5.0973	4.9768	4.9301	4.7616	50.968	48.753	4.6111
#3	5.1106	4.9608	4.9415	4.8133	51.074	47.871	4.6365
#4	5.1234	4.9768	4.9409	4.7680	51.012	48.222	4.5754
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.5000	5.5000	5.5000	5.5000	55.000	55.000	5.5000
Low	4.5000	4.5000	4.5000	4.5000	45.000	45.000	4.5000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	49.277	4.9146	4.9077	48.229	4.9874	4.9445	4.9100
SDev	.191	.0081	.0227	.253	.0257	.0151	.0855
%RSD	.38701	.16451	.46308	.52407	.51539	.30588	1.7408
#1	49.023	4.9030	4.8791	48.478	4.9818	4.9327	4.9733
#2	49.266	4.9195	4.9039	48.258	4.9918	4.9486	4.8363
#3	49.479	4.9154	4.9139	48.302	5.0191	4.9326	4.8364
#4	49.338	4.9206	4.9337	47.878	4.9570	4.9642	4.9940
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	55.000	5.5000	5.5000	55.000	5.5000	5.5000	5.5000
Low	45.000	4.5000	4.5000	45.000	4.5000	4.5000	4.5000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	4.9660	5.0382	4.9895	4.8405	4.8142	9.7242	4.9349
SDev	.0486	.0166	.0478	.0159	.0107	.1379	.0089
%RSD	.97819	.33045	.95754	.32855	.22188	1.4182	.17950
#1	4.9248	5.0216	4.9552	4.8300	4.8038	9.6531	4.9247
#2	5.0207	5.0276	5.0518	4.8362	4.8128	9.9183	4.9303
							~1~

Low

4.5000

Operator: MTW Sample Name: CCB6 Method: QUANMET

Run Time: 05/19/00 11:48:06 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CO	NC COTF.	ractor: 1					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00340	.02099	.04356	00368	.00098	.00071	.01077
SDev	.00227	.01186	.02635	.01755	.00030	.00037	.00603
%RSD	66.877	56.509	60.490	476.50	30.434	52.382	56.009
#1 #2 #3 #4	.00454 00001 .00453	.00673 .01595 .02876 .03251	.06155 .06944 .02958 .01366	00998 01147 01568 .02240	.00072 .00072 .00123 .00123	.00032 .00047 .00103 .00103	.00458 .00666 .01515 .01667
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00083	00072	00014	.00092	.00896	.31653	00156
SDev	.00192	.00187	.00256	.00089	.00508	.10276	.00045
%RSD	229.83	258.70	1862.2	95.900	56.699	32.464	28.982
#1	.00124	00290	.00200	00023	.00492	.27364	00189
#2	00340	.00001	00370	.00069	.00422	.19196	00189
#3	00069	00145	.00143	.00162	.01335	.41251	00153
#4	00049	.00145	00028	.00162	.01335	.38800	00093
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00898	00025	.01242	.02377	00523	.00385	.02093
SDev	.01496	.00084	.00248	.00439	.00995	.02033	.01364
%RSD	166.65	334.75	19.994	18.464	190.02	528.45	65.195
#1	.00345	00129	.01366	.02218	00517	.02306	.01562
#2	00207	00025	.01366	.01836	00595	00000	.00516
#3	.00345	00025	.01366	.02813	01707	.01539	.02600
#4	.03108	.00078	.00869	.02643	.00726	02305	.03692
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00494	.13663	.00682	.00071	.00061	00297	00202
SDev	.03083	.00301	.02064	.00057	.00059	.06371	.00066
%RSD	623.90	2.2044	302.57	79.201	97.590	2146.0	32.975
#1	00894	.14115	.01352	00004	00017	04635	00102
#2	01690	.13513	00308	.00061	.00052	.08103	00232

Analysis	Report			05/19,	/00 11:51:	65 12 am	7 965 page 2
#3 #4	03278 .03885	.13513 .13513	.03219 01535	.00115 .00115	.00087 .00121	.01147 05803	00231 00241
Errors High Low	LC Pass .25000 ~.25000	LC Pass .50000 50000	LC Pass .10000 10000	LC Pass .05000 05000	LC Pass .05000 05000	LC Pass .30000 30000	LC Pass .05000 05000
Elem Units Avge SDev %RSD	ZN ppm .00123 .00162 131.50						
#1 #2 #3 #4	.00152 .00151 .00289 00099						
Errors High Low	LC Pass .02000 02000						

Method: QUANMET Sample Name: DDDCMB Run Time: 05/19/00 11:54:15 Operator: MTW

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

ľ	lode: CUI	NC COLL.	ractor: I					
	Elem	AG	AL	AS	B_	BA	BE	CA
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	.00038	.01003	04580	.00866	.00039	.00014	.30121
	SDev	.00075	.00479	.03244	.00849	.00024	.00007	.00264
	%RSD	195.58	47.775	70.827	98.081	63.071	47.075	.87463
	#1	.00002	.00506	02196	.01521	.00020	.00017	.29809
	#2	.00150	.01414	02990	.01162	.00072	.00018	.30154
	#3	00001	.01412	09348	.01163	.00020	.00004	.30073
	#4	.00002	.00679	03787	00383	.00042	.00017	.30449
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
	Low	01000	20000	30000	20000	20000	00500	-5.0000
	Elem Units Avge SDev %RSD	CD ppm .00029 .00255 868.40	CO ppm 00290 .00264 91.200	CR ppm00271 .00108 39.824	CU ppm .00092 .00089 96.119	FE ppm .01686 .00099 5.8927	K_ ppm .33899 .03056 9.0160	LI ppm 00189 .00000
	#1	.00353	00001	00257	.00162	.01616	.37166	00189
	#2	00043	00435	00427	.00069	.01616	.33899	00189
	#3	.00068	00580	00199	00023	.01826	.29815	00189
	#4	00261	00144	00199	.00162	.01686	.34716	00189
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
	Low	00500	05000	01000	02500	10000	-5.0000	05000
	Elem	MG	MN	MO	NA	NI	PB	SB
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	00069	.00001	.00124	.02770	00902	01356	.01039
	SDev	.00529	.00052	.00497	.00350	.00358	.00738	.02505
	%RSD	765.94	10206.	399.20	12.638	39.636	54.427	241.06
	#1	.00069	00026	00621	.02388	00839	02315	01583
	#2	00484	.00078	.00373	.03237	01386	00779	.03668
	#3	00484	00025	.00373	.02728	00860	01559	00551
	#4	.00622	00025	.00373	.02728	00523	00772	.02623
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
	Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
	Elem	SE	SI	SN	SR	TI	TL	V_
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	00889	.11263	.02954	.00015	00009	.00853	.00002
	SDev	.02131	.02749	.02051	.00006	.00017	.05626	.00292
	%RSD	239.63	24.411	69.445	36.842	200.00	659.30	15648.
	#1	.01100	.09915	.00717	.00007	00017	02335	.00245
	#2	03277	.10509	.02604	.00018	00017	.09253	00251
								~ ~ ~

Analysis	Report			05/19	/00 11:57:	21 AM	657 967
#3 #4	.00703 02083	.09308 .15321	.05688 .02806	.00018	00017 .00017	02329 01176	00251 .00265
Errors High Low	LC Pass .25000 25000	LC Pass .50000 50000	LC Pass .10000 10000	LC Pass .05000 05000	LC Pass .05000 05000	LC Pass .30000 30000	LC Pass .05000 05000
Elem Units Avge SDev %RSD	ZN ppm .01918 .00053 2.7725						
#1 #2 #3 #4	.01904 .01960 .01958 .01848						
Errors High Low	LC Pass .02000 02000						

05/19/00 12:00:28 PM

page 1

Operator: MTW Sample Name: DDDCMC Method: QUANMET

Run Time: 05/19/00 11:57:23 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CO	NC COII.	raccor. x					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04600	1.8616	1.8347	.93770	1.8753	.04954	L.08879
SDev	.00145	.0071	.0315	.02523	.0178	.00031	.00381
%RSD	3.1486	.38298	1.7168	2.6901	.94894	.61678	4.2906
#1	.04484	1.8543	1.8488	.94564	1.8693	.04938	L.08812
#2	.04638	1.8579	1.7891	.94563	1.8788	.04964	L.08624
#3	.04789	1.8635	1.8407	.90098	1.8554	.04922	L.09436
#4	.04489	1.8707	1.8604	.95855	1.8978	.04991	L.08644
Errors	LC Pass	LC Pass	LC Pass	NOCHECK	LC Pass	LC Pass	LC Low
High	.06000	2.4000	2.4000		2.4000	.06000	60.000
Low	.04000	1.6000	1.6000		1.6000	.04000	40.000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04243	.47232	.18734	.23388	1.0125	L.46356	.89760
SDev	.00299	.00479	.00248	.00075	.0048	.51044	.01107
%RSD	7.0390	1.0137	1.3222	.32247	.47640	110.11	1.2331
#1	.04566	.46544	.18834	.23387	1.0098	L.60038	.89453
#2	L.03857	.47269	.18378	.23388	1.0098	L.59221	.89388
#3	.04352	.47560	.18947	.23296	1.0105	L.92712	.88833
#4	.04195	.47556	.18777	.23480	1.0197	L26547	.91368
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Low	NOCHECK
High	.06000	.60000	.24000	.30000	1.2000	60.000	
Low	.04000	.40000	.16000	.20000	.80000	40.000	
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	L01934	.47005	.95011	L.10401	.46906	.45750	.45859
SDev	.01852	.00259	.00951	.00638	.00954	.03118	.03897
%RSD	95.743	.55020	1.0011	6.1304	2.0327	6.8153	8.4973
#1	L00484	.46720	.93769	L.09818	.46665	.42065	.41111
#2	L03246	.47031	.95756	L.09903	.45663	.47490	.49556
#3	L00207	.46927	.94762	L.10794	.47544	.49032	.44288
#4	L03799	.47341	.95756	L.11091	.47753	.44413	.48480
Errors	LC Low	LC Pass	NOCHECK	LC Low	LC Pass	LC Pass	LC Pass
High	60.000	.60000		60.000	.60000	.60000	.60000
Low	40.000	.40000		40.000	.40000	.40000	.40000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.8294	4.6283	1.8703	.95511	.93360	1.7742	.46470
SDev	.0402	.0510	.0755	.00797	.00432	.0698	.00437
%RSD	2.1969	1.1023	4.0349	.83417	.46248	3.9352	.93972
#1	1.8015	4.5681	1.9012	.95157	.93169	1.8380	.45931
#2	1.8135	4.6042	1.7712	.95638	.93481	1.7800	.46484

Analysis	Report			05/19,	657 969 page 2		
#3 # <b>4</b>	1.8891 1.8135	4.6643 4.6764	1.8599 1.9491	.94693 .96557	.92891 .93897	1.6758 1.8031	.46466 .47000
Errors High Low	LC Pass 2.4000 1.6000	NOCHECK	NOCHECK	NOCHECK	NOCHECK	LC Pass 2.4000 1.6000	LC Pass .60000 .40000
Elem Units Avge SDev %RSD	ZN ppm .45852 .00323 .70376						
#1 #2 #3 #4	.45472 .45882 .45797 .46257						
Errors High Low	LC Pass .60000 .40000						

05/19/00 12:03:41 PM page 1 Analysis Report

Method: QUANMET Sample Name: DD5VC Run Time: 05/19/00 12:00:36 Operator: MTW

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CO	NC Corr.	Factor: I					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00466	.74370	.09242	.00299	.10078	.00004	5.1485
SDev	.00123	.00537	.02482	.00188	.00011	.00000	.0216
%RSD	26.520	.72233	26.856	62.854	.10509	.77814	.41938
#1	.00466	.75098	.09832	.00547	.10073	.00004	5.1320
#2	.00466	.73820	.11830	.00124	.10073	.00004	5.1483
#3	.00315	.74187	.05866	.00188	.10094	.00004	5.1345
#4	.00617	.74376	.09440	.00336	.10073	.00004	5.1790
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00150	00189	.11397	.12654	.56218	11.846	.00014
SDev	.00243	.00139	.00099	.00088	.00239	.298	.00134
%RSD	162.07	73.476	.86661	.69873	.42534	2.5174	941.79
#1	00141	00007	.11482	.12631	.56130	11.987	.00194
#2	.00157	00298	.11254	.12723	.55990	11.595	00083
#3	.00130	00298	.11425	.12538	.56200	11.603	00093
#4	.00453	00154	.11425	.12723	.56552	12.200	.00039
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.75773	.01487	.00505	56.550	.01616	.21143	.00169
SDev	.01105	.00060	.00625	.327	.00323	.03422	.01576
%RSD	1.4585	4.0170	123.73	.57798	19.962	16.184	930.85
#1	.75221	.01539	.01374	56.364	.01388	.23076 '.16901 .19982 .24613	.01476
#2	.75221	.01538	.00381	56.614	.01566		00613
#3	.75221	.01435	.00381	56.240	.02086		01673
#4	.77431	.01435	00116	56.981	.01424		.01487
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem Units Avge SDev %RSD	SE ppm .02195 .03038 138.38	SI ppm H27.228 .102 .37629	SN ppm .03575 .01960 54.820	SR ppm .04610 .00000	TI ppm .01509 .00049 3.2511	TL ppm .03773 .07159 189.71	V_ ppm 00225 .00012 5.3883
#1	.01300	H27.088	.04874	.04610	.01543	.02905	00209
#2	.01697	H27.238	.00726	.04610	.01509	04039	00228

Analysis Report	A	na	1y	si	S	R	eı	þ	2	rı	t
-----------------	---	----	----	----	---	---	----	---	---	----	---

#4

High Low .41965

100.00

Errors LC Pass

05/19/00 12:03:41 PM

page 2

#3	00690	H27.250	.04860	.04610	.01439	.02906	00228
#4	.06475	H27.334	.03840	.04610	.01543	.13323	00237
Errors	LC Pass	LC High	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	100.00	50.000	50.000	100.00	100.00
Low	25000	50000	10000	05000	05000	-2.0000	05000
Elem Units Avge SDev %RSD	ZN ppm .41855 .00285 .68184						
#1 #2 #3	.41506 .41774 .42177						

Operator: MTW

Method: QUANMET Sample Name: DD5VE Operate Run Time: 05/19/00 12:03:49 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: Col	NC COII.	ractor. I					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.13957	17.779	.03917	.06517	1.5337	00048	43.521
SDev	.00000	.024	.06516	.00218	.0013	.00007	.077
%RSD	.00310	.13294	166.35	3.3480	.08461	14.645	.17726
#1	.13957	17.813	.11240	.06570	1.5356	00059	43.488
#2	.13956	17.776	.00147	.06208	1.5331	00045	43.506
#3	.13956	17.758	03017	.06568	1.5328	00045	43.457
#4	.13957	17.770	.07299	.06722	1.5333	00045	43.632
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.02092	.02062	1.5403	1.6204	13.987	5.3401	.00937
SDev	.00624	.00218	.0056	.0018	.017	.2602	.00215
%RSD	29.832	10.569	.36670	.11413	.11940	4.8731	22.899
#1	.01319	.02173	1.5416	1.6213	13.990	5.2237	.01220
#2	.02462	.01735	1.5399	1.6176	13.973	5.1420	.00780
#3	.02714	.02169	1.5331	1.6213	13.976	5.7220	.00762
#4	.01874	.02172	1.5468	1.6213	14.009	5.2727	.00987
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	5.5652	.22008	.01451	23.913	.27071	.91383	00613
SDev	.0098	.00051	.00625	.081	.00551	.02183	.00008
%RSD	.17670	.23204	43.081	.33763	2.0374	2.3887	1.2273
#1	5.5569	.22034	.00582	24.011	.26257	.89843	00611
#2	5.5708	.22034	.01575	23.881	.27370	.89833	00609
#3	5.5763	.22034	.01575	23.822	.27206	.91387	00608
#4	5.5569	.21932	.02072	23.937	.27449	.94469	00624
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01929	16.240	.27482	.19958	.39380	03350	.02137
SDev	.03100	.045	.02186	.00014	.00040	.05553	.00012
%RSD	160.69	.27515	7.9532	.06767	.10169	165.77	.57672
#1	.02229	16.174	.25258	.19965	.39414	.01858	.02121
#2	.03815	16.264	.26293	.19965	.39345	09701	.02140

Analysis	Report			05/19,	657 973 page 2		
#3 #4	.04214 02540	16.252 16.270	.28155 .30224	.19938 .19965	.39414 .39345	.00714 06269	.02138 .02151
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm 6.2305 .0026						
#1 #2 #3 #4	6.2329 6.2270 6.2304 6.2318						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW Method: QUANMET Sample Name: DD9NT

Run Time: 05/19/00 12:06:57

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: Co.	NC COLL.	raccor. x					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01888	8.6398	.08066	.05652	.21910	00033	14.453
SDev	.00124	.0257	.05739	.01619	.00151	.00008	.063
%RSD	6.5523	.29757	71.154	28.639	.68952	22.723	.43255
#1	.02041	8.6421	.07264	.04844	.21946	00027	14.408
#2	.01738	8.6677	.07641	.04846	.22028	00027	14.479
#3	.01888	8.6441	.01706	.08081	.21977	00040	14.529
#4	.01887	8.6054	.15651	.04839	.21689	00040	14.396
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01033	.00479	.18277	.41943	2.9966	14.060	.00286
SDev	.00128	.00145	.00135	.00276	.0147	.323	.00024
%RSD	12.401	30.186	.73565	.65749	.48892	2.2953	8.2447
#1	.00907	.00407	.18376	.42012	3.0025	13.833	.00295
#2	.00944	.00407	.18262	.42289	3.0082	14.528	.00303
#3	.01105	.00696	.18376	.41827	3.0004	14.021	.00295
#4	.01176	.00407	.18092	.41642	2.9751	13.858	.00251
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	2.4390	.07352	.00542	65.286	.07977	.55760	.01739
SDev	.0256	.00052	.00475	.551	.00918	.02522	.01583
%RSD	1.0485	.70641	87.807	.84463	11.508	4.5225	91.063
#1	2.4486	.07377	.00417	65.112	.08023	.52673	.00423
#2	2.4597	.07378	.00914	65.808	.09229	.55757	.02543
#3	2.4459	.07377	00079	64.589	.07567	.58850	.03577
#4	2.4017	.07274	.00914	65.637	.07088	.55758	.00412
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01593	H26.596	.06086	.13219	.08609	.01571	.00562
SDev	.07964	.047	.01607	.00068	.00052	.08080	.00236
%RSD	499.96	.17688	26.414	.51423	.60423	514.22	41.919
#1	.09753	H26.548	.06131	.13233	.08583	.11986	.00818
#2	.04980	H26.656	.04072	.13287	.08583	07705	.00698

Analysis	Report			05/19	657 975 page 2		
#3 #4	08952 .00590	H26.608 H26.572	.08008 .06131	.13233 .13125	.08687 .08583	.00409 .01595	.00422 .00311
Errors High Low	LC Pass 100.00 25000	LC High 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm 1.5865 .0100 .63201						
#1 #2 #3 #4	1.5809 1.5911 1.5981 1.5758						
Errors High Low	LC Pass 100.00 02000						

Sample Name: DD9NTP5 Operator: MTW Method: QUANMET

Run Time: 05/19/00 12:10:10

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: cor	WC COII.	100001					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00506	1.8096	.03770	.00042	.04491	00003	3.1414
SDev	.00145	.0157	.02165	.00106	.00030	.00008	.0088
%RSD	28.747	.86999	57.419	252.99	.66127	249.93	.27878
#1	.00316	1.7977	.03384	00011	.04517	00010	3.1405
#2	.00467	1.8014	.00998	00011	.04466	00010	3.1299
#3	.00619	1.8326	.06137	.00201	.04517	.00004	3.1444
#4	.00621	1.8068	.04562	00011	.04466	.00003	3.1508
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00212	00042	.03590	.08661	.63068	3.0489	00136
SDev	.00239	.00217	.00136	.00076	.00373	.2471	.00052
%RSD	112.57	518.23	3.7793	.87258	.59174	8.1038	38.111
#1	.00482	.00139	.03733	.08754	.63366	2.9447	00189
#2	.00225	00296	.03447	.08568	.62594	2.8549	00145
#3	00099	.00140	.03505	.08661	.63366	2.9856	00065
#4	.00240	00151	.03675	.08661	.62945	3.4103	00145
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.50907	.01487	.00382	13.667	.01263	.10874	.00501
SDev	.01034	.00060	.00406	.179	.00378	.01153	.00859
%RSD	2.0307	4.0211	106.18	1.3118	29.950	10.606	171.40
#1 #2 #3 #4	.50907 .49802 .50631 .52288	.01538 .01435 .01435 .01538	.00382 00115 .00382 .00879	13.462 13.887 13.712	.01566 .01081 .00817 .01588	.09915 .12216 .11455 .09909	00549 .00506 .00494 .01556
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	01560	6.0279	.00629	.02697	.01777	.01385	00147
SDev	.01048	.0427	.01562	.00034	.00052	.05934	.00200
%RSD	67.137	.70894	248.53	1.2556	2.9268	428.48	136.33
#1	01062	6.0519	.00314	.02735	.0171 <b>7</b>	00645	00243
#2	03054	6.0279	00923	.02692	.01821	.00524	00253

05/19/00 12:13:15 PM

page 2

#3	01460	6.0640	.02801	.02708	.01751	04120	00243
#4	00665	5.9679	.00323	.02654	.01821	.09781	.00153
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	20.000	100.00	50.000	50.000	100.00	100.00
Low	25000	50000	10000	05000	05000	-2.0000	05000
Elem	ZN						

Units ppm Avge .34333 SDev .00143 %RSD .41558 #1 .34464 .34443 #2 .34174 #3 #4 .34252 Errors LC Pass

100.00

-.02000

High

Low

Method: QUANMET Sample Name: DD9NTS Operator: MTW

Run Time: 05/19/00 12:13:18

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CONC Corr. Factor: 1

Mode. ce	,,,,,						
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.06486	10.403	1.9511	.96418	2.0596	.04977	14.424
SDev	.00077	.050	.0458	.00005	.0173	.00057	.079
%RSD	1.1832	.48056	2.3479	.00529	.84138	1.1435	.54535
#1	.06522	10.371	1.8889	.96412	2.0449	.04922	14.314
#2	.06371	10.353	1.9605	.96415	2.0446	.04936	14.423
#3	.06526	10.461	1.9992	.96422	2.0711	.05019	14.471
#4	.06526	10.428	1.9558	.96422	2.0778	.05033	14.489
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.04788	.46796	.36956	.64521	3.9424	14.262	.91081
SDev	.00107	.00372	.00305	.00615	.0303	.313	.01195
%RSD	2.2295	.79445	.82506	.95277	.76869	2.1922	1.3116
#1	.04856	.46363	.36614	.63943	3.9080	14.470	.92523
#2	.04643	.46653	.36785	.64036	3.9262	14.340	.91544
#3	.04774	.46939	.37184	.65052	3.9642	14.438	.90419
#4	.04879	.47228	.37240	.65053	3.9712	13.801	.89836
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	2.3948	.53737	1.0375	66.299	.54793	.96674	.46971
SDev	.0269	.00362	.0122	.705	.00843	.00780	.03367
%RSD	1.1245	.67412	1.1728	1.0631	1.5376	.80642	7.1683
#1	2.3685	.53298	1.0424	67.185	.54856	.95504	.44077
#2	2.3768	.53608	1.0275	66.461	.54203	.97063	.50403
#3	2.4265	.53919	1.0524	66.029	.54152	.97063	.49321
#4	2.4072	.54125	1.0275	65.520	.55962	.97065	.44084
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.9873	H33.917	2.0804	1.0760	1.0999	1.7473	.46901
SDev	.0782	.291	.0255	.0085	.0092	.1290	.00344
%RSD	3.9373	.85777	1.2272	.79362	.83701	7.3808	.73352
#1	1.9196	H33.588	2.0647	1.0684	1.0895	1.7505	.46555
#2	1.9793	H33.774	2.0526	1.0691	1.0950	1.9240	.46656
							~ ~ ~ ~

Analysis	Report			05/19,	657 979 page 2		
#3 #4	2.0989 1.9516	H34.068 H34.236	2.1022 2.1019	1.0816 1.0850	1.1058 1.1093	1.6920 1.6224	.47220 .47172
Errors High Low	LC Pass 100.00 25000	LC High 20.000 50000		LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm 2.0207 .0173 .85495				N.		
#1 #2 #3 #4	1.9990 2.0148 2.0367 2.0322						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW

Method: QUANMET Sample Name: DD9NTD Operate Run Time: 05/19/00 12:16:31 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode:	CON	C Corr.	Factor: 1					
Elem Unit Avge SDev %RSD	s j	AG ppm .06827 .00123 1.8070	AL ppm 10.221 .085 .82894	AS ppm 1.9407 .0701 3.6105	B_ppm .95236 .01275 1.3383	BA ppm 2.0672 .0214 1.0361	BE ppm .05015 .00058 1.1617	CA ppm 14.313 .037 .25951
#1		.06978	10.180	1.8745	.96033	2.0541	.04976	14.315
#2		.06827	10.137	1.9464	.94952	2.0482	.04962	14.261
#3		.06676	10.233	1.9058	.93561	2.0706	.05032	14.345
#4		.06828	10.334	2.0361	.96399	2.0961	.05089	14.333
Erro	. :	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High		2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low		01000	20000	30000	20000	20000	00500	-5.0000
Elem	is j	CD	CO	CR	CU	FE	K_	LI
Unit		ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge		.05226	.48294	.37353	.63575	3.8366	14.270	.89802
SDev		.00250	.00447	.00395	.00687	.0222	.252	.01357
%RSD		4.7944	.92614	1.0567	1.0811	.57915	1.7677	1.5110
#1		.05149	.48260	.37011	.63206	3.8248	14.479	.90080
#2		.05522	.47680	.37011	.62928	3.8121	14.389	.88232
#3		.05303	.48692	.37695	.63668	3.8480	13.907	.89411
#4		.04928	.48546	.37695	.64499	3.8614	14.307	.91484
Erro		LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High		100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low		00500	05000	01000	02500	10000	-5.0000	05000
Elem Unit Avge SDev %RSI	cs e v	MG ppm 2.3713 .0248 1.0465	MN ppm .54558 .00285 .52214	MO ppm .95798 .00405 .42320	NA ppm 65.737 .884 1.3450	NI ppm .56142 .00279 .49678	PB ppm .98456 .02296 2.3325	SB ppm .43652 .02500 5.7270
#1		2.3464	.54429	.96295	65.776	.55862	.97881	.44176
#2		2.3547	.54222	.95798	64.704	.55948	1.0172	.45236
#3		2.3989	.54739	.95302	65.610	.56326	.97891	.45220
#4		2.3851	.54843	.95798	66.860	.56431	.96333	.39977
Erre	h	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
Hig		600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low		-5.0000	01500	04000	-5.0000	04000	10000	06000
Ele Uni Avg SDe %RS	ts e v	SE ppm 1.9452 .0571 2.9372	sī ppm H20.164 .094 .46558	SN ppm 1.9646 .0518 2.6376	SR ppm 1.0815 .0108 1.0015	TI ppm 1.0103 .0074 .73481	TL ppm 1.7511 .0446 2.5460	V_ ppm .47554 .00007 .01536
#1		1.9710	H20.102	1.9698	1.0750	1.0059	1.7628	.47562
#2		1.8595	H20.102	1.9492	1.0712	1.0031	1.8093	.47553
		-						610

657 981

						657	981 .
Analysis	Report			05/19/00 12:19:36 PM			page 2
#3 #4	1.9751 1.9751	H20.150 H20.300	2.0319 1.9075	1.0842 1.0956	1.0125 1.0198	1.7162 1.7161	.47545 .47554
Errors High Low	LC Pass 100.00 25000	LC High 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm 2.0216 .0133 .65972						
#1 #2 #3 #4	2.0206 2.0039 2.0265 2.0355						
Errors High Low	LC Pass 100.00 02000						

Method: QUANMET Sample Name: DD9NW Operator: MTW

Run Time: 05/19/00 12:19:44

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00359	5.1141	.05675	.03770	2.1494	.00062	6.4637
SDev	.00214	.0234	.00874	.00319	.0118	.00007	.0237
%RSD	59.550	.45757	15.408	8.4677	.54944	11.473	.36674
#1	.00359	5.0966	.05987	.04145	2.1493	.00051	6.4565
#2	.00208	5.1479	.05944	.03364	2.1657	.00065	6.4474
#3	.00662	5.1114	.06372	.03786	2.1448	.00065	6.4988
#4	.00208	5.1003	.04395	.03784	2.1379	.00065	6.4521
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00029	.00837	.09913	.27021	1.6811	3.9903	.00333
SDev	.00084	.00303	.00099	.00046	.0053	.1906	.00020
%RSD	290.30	36.150	.99435	.17106	.31510	4.7775	5.9096
#1	00050	.00765	.09999	.26952	1.6790	4.0066	.00314
#2	00091	.00474	.09828	.27044	1.6868	3.7207	.00339
#3	.00095	.00909	.09999	.27044	1.6839	4.0720	.00321
#4	00070	.01201	.09828	.27044	1.6748	4.1618	.00358
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.9085	.03654	.02260	17.742	.05756	.37483	.00610
SDev	.0237	.00052	.00625	.170	.00441	.02182	.02332
%RSD	1.2426	1.4143	27.649	.95796	7.6664	5.8216	382.13
#1	1.8795	.03628	.02384	17.653	.05555	.34401	.01395
#2	1.9071	.03629	.02385	17.977	.05562	.37472	.01401
#3	1.9375	.03628	.01391	17.588	.06416	.39022	02808
#4	1.9099	.03732	.02881	17.749	.05491	.39037	.02453
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.02704	17.889	.03395	.25081	.86198	.03465	.03415
SDev	.03321	.071	.01659	.00125	.00290	.04063	.00012
%RSD	122.81	.39967	48.856	.49700	.33661	117.27	.34967
#1	.00912	17.802	.04381	.25065	.86233	00586	.03417
#2	01075	17.916	.01691	.25254	.86441	.06356	.03417

Analysis	Report			05/19	657 98; page 2		
#3 #4	.05690 .05288	17.970 17.868	.05189 .02319	.25049 .24957	.86337 .85782	.07514 .00574	.03398 .03426
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000		LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .32510 .00126 .38843						
#1 #2 #3 #4	.32328 .32518 .32594 .32598						
Errors High Low	LC Pass 100.00 02000						

3

Method: QUANMET Sample Name: DD9NWP5 Run Time: 05/19/00 12:22:52 Operator: MTW

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00123	1.0269	.00381	00464	.42977	.00024	1.2983
SDev	.00287	.0117	.05817	.00752	.00349	.00008	.0083
%RSD	232.19	1.1409	1525.9	161.95	.81181	32.849	.64015
#1	00293	1.0095	03476	01084	.43388	.00017	1.2867
#2	.00312	1.0314	05482	.00187	.42771	.00031	1.3049
#3	.00163	1.0351	.06825	.00186	.43131	.00030	1.2980
#4	.00312	1.0314	.03658	01146	.42617	.00017	1.3037
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00104	.00059	.01838	.05404	.33794	.91691	00163
SDev	.00204	.00182	.00256	.00089	.00202	.23146	.00078
%RSD	197.19	307.05	13.947	1.6375	.59753	25.243	48.186
#1	.00101	.00095	.01624	.05335	.33495	.58404	00244
#2	00199	.00096	.01852	.05520	.33917	1.0660	00057
#3	.00026	.00240	.02193	.05427	.33917	.93529	00189
#4	00343	00194	.01681	.05335	.33847	1.0823	00161
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.38128	.00705	.00502	3.5747	.00533	.07262	.00222
SDev	.00523	.00000	.00248	.0311	.00414	.03695	.02168
%RSD	1.3717	.01476	49.499	.87039	77.670	50.876	978.58
#1	.37369	.00705	.00378	3.6138	.00716	.05335	01618
#2	.38474	.00705	.00874	3.5442	.00960	.10732	.01541
#3	.38197	.00705	.00378	3.5846	.00467	.03026	01621
#4	.38474	.00705	.00378	3.5561	00010	.09955	.02584
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01118	4.0682	.00618	.05026	.17115	.03705	.00400
SDev	.04496	.0296	.01093	.00038	.00057	.05865	.00257
%RSD	401.97	.72693	176.93	.76003	.33090	158.31	64.174
#1	05549	4.0577	.00661	.05080	.17045	04976	.00268
#2	.04004	4.1118	.00055	.05026	.17115	.06599	.00278

Analysis	Report			05/19,	/00 12:25:	58 PM	657 985 page 2
#3 #4	.02412 .03606	4.0578 4.0457	00374 .02129	.04999 .04999	.17184 .17115	.05436 .07760	.00785 .00268
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .06553 .00077 1.1792						
#1 #2 #3 #4	.06475 .06505 .06644 .06588						
Errors High Low	LC Pass 100.00 02000						

Method: QUANMET Sample Name: CCV2-7 Run Time: 05/19/00 12:26:01 Operator: MTW

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

M	lode: CON	Corr.	ractor: 1					
	Elem	AG	AL	AS	B_	BA	BE	CA
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	.98884	48.980	5.1467	4.8910	4.8056	5.0980	52.658
	SDev	.00403	.194	.0529	.0218	.0287	.0234	.159
	%RSD	.40775	.39664	1.0286	.44506	.59822	.45987	.30191
	#1	.98852	49.065	5.1706	4.9014	4.8135	5.1124	52.756
	#2	.98533	48.690	5.2102	4.8584	4.7631	5.0630	52.755
	#3	.98694	49.065	5.1033	4.9013	4.8247	5.1097	52.423
	#4	.99456	49.100	5.1028	4.9029	4.8212	5.1070	52.699
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	1.1000	55.000	5.5000	5.5000	5.5000	5.5000	55.000
	Low	.90000	45.000	4.5000	4.5000	4.5000	4.5000	45.000
	Elem	CD	CO	CR	CU	FE	K_	LI
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	5.1936	5.0605	5.0340	4.8358	51.770	47.973	4.5111
	SDev	.0159	.0145	.0115	.0319	.096	.384	.0378
	%RSD	.30667	.28716	.22801	.65941	.18475	.80039	.83786
	#1	5.1887	5.0735	5.0434	4.8384	51.849	48.141	4.5023
	#2	5.2155	5.0402	5.0326	4.7903	51.640	48.059	L4.4648
	#3	5.1776	5.0605	5.0184	4.8624	51.757	47.414	4.5220
	#4	5.1928	5.0677	5.0417	4.8522	51.835	48.280	4.5552
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	5.5000	5.5000	5.5000	5.5000	55.000	55.000	5.5000
	Low	4.5000	4.5000	4.5000	4.5000	45.000	45.000	4.5000
	Elem	MG	MN	MO	NA	NI	PB	SB
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	49.952	5.0029	4.9773	47.949	5.1341	4.9952	5.0871
	SDev	.203	.0108	.0238	.323	.0154	.0289	.0623
	%RSD	.40651	.21551	.47800	.67414	.29977	.57851	1.2240
	#1	50.048	5.0148	4.9885	47.900	5.1244	4.9743	5.0790
	#2	49.653	4.9910	4.9487	47.577	5.1560	5.0198	5.0268
	#3	50.007	4.9973	4.9686	47.951	5.1224	4.9663	5.1742
	#4	50.101	5.0087	5.0034	48.365	5.1336	5.0203	5.0685
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	55.000	5.5000	5.5000	55.000	5.5000	5.5000	5.5000
	Low	45.000	4.5000	4.5000	45.000	4.5000	4.5000	4.5000
	Elem	SE	SI	SN	SR	TI	TL	V_
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	5.0375	H7.3295	5.0285	4.9306	4.8840	9.7673	5.0244
	SDev	.0448	.0919	.0245	.0252	.0159	.1374	.0105
	%RSD	.88974	1.2536	.48703	.51139	.32563	1.4070	.20902
	#1	4.9960	H7.2575	5.0486	4.9395	4.8960	9.7693	5.0327
	#2	5.0152	H7.4495	5.0402	4.8930	4.8606	9.6328	5.0100
								7101

LC Pass

5.5000

4.5000

Errors

High Low

Operator: MTW

Method: QUANMET Sample Name: CCB7 Operate Run Time: 05/19/00 12:29:14 Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

110001							
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00037	.01596	00606	01157	.00064	.00064	.00733
SDev	.00076	.00637	.02473	.00105	.00023	.00029	.00583
%RSD	205.53	39.894	407.94	9.1008	35.736	45.342	79.651
#1	.00150	.01785	04184	00999	.00042	.00033	.00514
#2	00001	.00677	.01389	01210	.00051	.00047	.00250
#3	00001	.01776	00211	01209	.00093	.00089	.00585
#4	00001	.02147	.00581	01210	.00072	.00089	.01581
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.01000	.20000	.30000	.20000	.20000	.00500	5.0000
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00167	00073	00128	.00069	.00650	.21646	00196
SDev	.00254	.00277	.00261	.00075	.00365	.15839	.00035
%RSD	152.13	379.39	204.04	108.76	56.106	73.173	17.721
#1	.00241	.00144	00256	00023	.00492	.20013	00161
#2	00186	00434	00313	.00069	.00211	.10211	00244
#3	.00195	00145	00199	.00162	.00983	.11844	00189
#4	.00416	.00144	.00257	.00069	.00913	.44518	00189
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	.00500	.05000	.01000	.02500	.10000	5.0000	.05000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem Units Avge SDev %RSD	MG ppm 00069 .00160 230.94	MN ppm 00025 .00000	MO ppm .01118 .00287 25.662	NA ppm .01889 .00651 34.443	NI ppm 00868 .00814 93.739	PB ppm 01158 .04932 425.94	SB ppm .00778 .01018 130.83
#1	.00069	00025	.00869	.01751	00973	06160	.01571
#2	00207	00025	.00869	.01284	00267	.03071	.00519
#3	.00069	00025	.01366	.01709	01980	04630	00554
#4	00207	00025	.01366	.02813	00253	.03087	.01575
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	5.0000	.01500	.04000	5.0000	.04000	.10000	.06000
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	04276	H.65923	00927	.00042	.00017	.04338	00236
SDev	.01321	.04637	.01743	.00050	.00057	.05538	.00006
%RSD	30.895	7.0336	188.01	120.27	326.60	127.68	2.4997
#1	04077	H.72380	01337	.00007	.00017	.04627	00242
#2	06068	H.65773	.00106	00004	00052	03471	00242

Analysis	Report			05/19	657 989 page 2		
#3 #4	04076 02882		.00722 03200	.00061 .00104	.00017	.09256 .06939	00232 00231
Errors High Low	LC Pass .25000 25000	LC High .50000 50000	LC Pass .10000 10000	LC Pass .05000 05000	LC Pass .05000 05000	LC Pass .30000 30000	LC Pass .05000 05000
Elem Units Avge SDev %RSD	ZN ppm 00076 .00109 143.34						
#1 #2 #3 #4	.00018 00143 00193 .00015						
Errors High Low	LC Pass .02000 02000						

Method: QUANMET Sample Name: DD9NWS Run Time: 05/19/00 12:32:27 Operator: MTW

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Ī	Mode: COI	NC Corr.	Factor: 1					
	Elem	AG	AL	AS	B_	BA	BE	CA
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	.04630	5.3364	1.7822	.84361	3.6160	.04717	5.2723
	SDev	.00146	.0384	.0174	.00284	.0289	.00024	.0220
	%RSD	3.1561	.71906	.97787	.33627	.79939	.51163	.41648
	#1	.04515	5.2957	1.7905	.84231	3.6074	.04696	5.2475
	#2	.04667	5.3121	1.7625	.84022	3.5791	.04696	5.2976
	#3	.04821	5.3743	1.7738	.84596	3.6427	.04737	5.2817
	#4	.04516	5.3635	1.8018	.84594	3.6350	.04738	5.2622
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
	Low	01000	20000	30000	20000	20000	00500	-5.0000
	Elem	CD	CO	CR	CU	FE	K_	LI
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	.04039	.45207	.27378	.47425	2.2870	4.0679	.83153
	SDev	.00242	.00314	.00203	.00380	.0095	.2978	.00978
	%RSD	5.9943	.69574	.74093	.80059	.41483	7.3204	1.1757
	#1	.04271	.45208	.27378	.47310	2.2740	4.0393	.82579
	#2	.04022	.45355	.27264	.46941	2.2860	4.4967	.82226
	#3	.04153	.44770	.27207	.47771	2.2944	3.8432	.84429
	#4	.03709	.45496	.27663	.47679	2.2937	3.8923	.83379
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
	Low	00500	05000	01000	02500	10000	-5.0000	05000
	Elem	MG	MN	MO	NA	NI	PB	SB
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	1.3393	.47547	.82736	18.640	.50417	.81276	.11307
	SDev	.0144	.00155	.00625	.196	.00940	.02029	.01802
	%RSD	1.0719	.32648	.75554	1.0516	1.8651	2.4958	15.933
	#1	1.3241	.47366	.81867	18.481	.49898	.81086	.11576
	#2	1.3435	.47470	.83357	18.481	.51603	.84170	.10517
	#3	1.3573	.47676	.82861	18.883	.49469	.80296	.09458
	#4	1.3324	.47676	.82861	18.714	.50698	.79551	.13678
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
	Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
	Elem	SE	SI	SN	SR	TI	TL	V_
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	1.8400	H35.444	1.7697	1.0915	1.4849	1.3328	.46379
	SDev	.0916	.326	.0300	.0077	.0077	.0359	.00259
	%RSD	4.9784	.91929	1.6945	.70684	.51782	2.6953	.55939
	#1	1.7982	H35.065	1.7470	1.0881	1.4803	1.2808	.46233
	#2	1.8340	H35.288	1.8132	1.0822	1.4765	1.3618	.46262
								~ ~ ~

Analysis	Report			05/19	657 991 page 2		
#3 #4	1.9694 1.7585	H35.648 H35.774	1.7531 1.7657	1.0985 1.0970	1.4921 1.4907	1.3385 1.3501	.46768 .46253
Errors High Low	LC Pass 100.00 25000	LC High 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .75616 .00169 .22360						
#1 #2 #3 #4	.75459 .75828 .75675 .75502						
Errors High Low	LC Pass 100.00 02000		•				

Method: QUANMET Sample Name: DD9NWD Run Time: 05/19/00 12:35:40 Operator: MTW

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

ľ	Mode: COI	Corr.	Factor: 1					
	Elem	AG	AL	AS	B_	BA	BE	CA
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	.04901	6.2145	1.9057	.84313	3.7783	.04808	6.7328
	SDev	.00151	.0181	.0196	.00344	.0162	.00013	.0295
	%RSD	3.0886	.29123	1.0269	.40783	.42776	.27940	.43851
	#1	.04977	6.2117	1.8978	.83863	3.7892	.04819	6.7218
	#2	.04977	6.1899	1.8821	.84579	3.7557	.04791	6.7464
	#3	.04978	6.2280	1.9254	.84225	3.7778	.04818	6.7654
	#4	.04674	6.2282	1.9175	.84584	3.7907	.04805	6.6975
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
	Low	01000	20000	30000	20000	20000	00500	-5.0000
	Elem	CD	CO	CR	CU	FE	K_	LI
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	.04177	.48445	.29032	.49990	2.2646	4.1087	.83776
	SDev	.00154	.00522	.00444	.00106	.0068	.2864	.00916
	%RSD	3.6867	1.0785	1.5291	.21275	.29905	6.9715	1.0933
	#1	.04300	.48335	.28519	.50082	2.2625	3.8514	.84107
	#2	.03959	.49208	.29374	.49898	2.2561	4.3007	.82597
	#3	.04268	.48191	.29431	.49898	2.2716	4.4069	.83628
	#4	.04183	.48046	.28804	.50082	2.2681	3.8759	.84773
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
	Low	00500	05000	01000	02500	10000	-5.0000	05000
	Elem	MG	MN	MO	NA	NI	PB	SB
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	1.9803	.50904	.90187	18.399	.54670	.85557	.10418
	SDev	.0094	.00099	.01102	.192	.00864	.01708	.01720
	%RSD	.47655	.19405	1.2213	1.0434	1.5807	1.9969	16.511
	#1	1.9679	.50775	.89814	18.466	.54549	.88058	.10429
	#2	1.9845	.50981	.89317	18.146	.53494	.85004	.12514
	#3	1.9900	.50879	.91801	18.380	.55319	.84198	.08300
	#4	1.9789	.50981	.89814	18.603	.55319	.84967	.10429
	Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
	High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
	Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
	Elem	SE	SI	SN	SR	TI	TL	V_
	Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
	Avge	1.8867	H25.500	1.9806	1.1487	1.7909	1.2171	.47976
	SDev	.0364	.103	.0269	.0042	.0042	.0609	.00086
	%RSD	1.9293	.40555	1.3569	.36654	.23691	5.0023	.17895
	#1	1.9056	H25.466	1.9506	1.1516	1.7945	1.2576	.47936
	#2	1.8379	H25.370	2.0108	1.1427	1.7848	1.1650	.47928
								C10

Analysis	Report			05/19/	/00 12:38:4	16 PM	657 993 page 2
#3 #4	1.8817 1.9215	H25.605 H25.557	1.9940 1.9672	1.1491 1.1516	1.7928 1.7914	1.2807 1.1650	.48105 .47937
Errors High Low	LC Pass 100.00 25000	LC High 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm .80040 .00829 1.0352						
#1 #2 #3 #4	.79717 .79048 .80928 .80467						
Errors High Low	LC Pass 100.00 02000						

Operator: MTW Method: QUANMET Sample Name: DD9PE

Run Time: 05/19/00 12:38:54

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: Cor	NC COII.	raccor. I					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.02216	7.4960	.02933	.04749	.17488	00050	10.762
SDev	.00145	.0285	.01045	.00002	.00114	.00007	.029
%RSD	6.5411	.38024	35.621	.03863	.65222	14.080	.26865
#1	.02330	7.5220	.02313	.04752	.17553	00054	10.797
#2	.02178	7.4654	.03158	.04747	.17450	00054	10.770
#3	.02027	7.5184	.04305	.04750	.17604	00054	10.728
#4	.02330	7.4782	.01955	.04749	.17347	00040	10.752
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00344	.00054	.07778	.59643	2.5035	4.2333	.00119
SDev	.00258	.00187	.00203	.00226	.0083	.2069	.00076
%RSD	74.877	344.06	2.6070	.37899	.33077	4.8875	64.195
#1	.00596	00019	.07778	.59735	2.5149	4.1373	.00218
#2	.00533	.00271	.07892	.59458	2.4952	4.3660	.00119
#3	.00157	00163	.07493	.59920	2.5030	3.9903	.00031
#4	.00090	.00127	.07949	.59458	2.5009	4.4395	.00109
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	1.7393	.04680	.00534	6.6279	.04299	.08419	.01273
SDev	.0091	.00000	.00625	.0740	.00572	.00999	.02165
%RSD	.52487	.00854	117.03	1.1171	13.310	11.867	170.14
#1	1.7330	.04680	.01404	6.6787	.04790	.08798	00572
#2	1.7496	.04679	00087	6.5759	.03772	.08043	.02586
#3	1.7303	.04680	.00410	6.7033	.04797	.07253	00559
#4	1.7441	.04680	.00410	6.5538	.03836	.09581	.03635
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem	SE	SI	SN	SR	TI	TL	V_
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	00478	11.861	.04018	.05546	.05107	.00044	.00283
SDev	.06688	.123	.01641	.00026	.00066	.03039	.00012
%RSD	1399.9	1.0379	40.844	.46630	1.2856	6868.1	4.2411
#1	.07187	12.035	.03854	.05566	.05150	.03217	.00300
#2	.00017	11.843	.04903	.05539	.05011	02554	.00272

Analysis	Report			05/19	/00 12:41:	59 PM	657 995 page 2
#3 #4	09134 .00019	11.819 11.747	.01785 .05529	.05566 .05512	.05115 .05150	.02073 02559	.00281 .00281
Errors High Low	LC Pass 100.00 25000	LC Pass 20.000 50000	LC Pass 100.00 10000	LC Pass 50.000 05000	LC Pass 50.000 05000	LC Pass 100.00 -2.0000	LC Pass 100.00 05000
Elem Units Avge SDev %RSD	ZN ppm 1.9912 .0065 .32495			•			
#1 #2 #3 #4	1.9872 1.9940 1.9990 1.9848						
Errors High Low	LC Pass 100.00 02000					,	

Operator: MTW Sample Name: DD9PX Method: QUANMET

Run Time: 05/19/00 12:42:02

Comment: STL PITTSBURGH ICP METALS ANALYSIS-INSTRUMENT JA61E

Mode: CC	ONC Corr.	ractor: 1					
Elem	AG	AL	AS	B_	BA	BE	CA
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.01183	9.2400	.09458	.12169	.23767	.00044	18.788
SDev	.00213	.0539	.03091	.01933	.00192	.00000	.078
%RSD	18.047	.58294	32.684	15.882	.80751	.03235	.41440
#1	.01031	9.2341	.08867	.13660	.23768	.00044	18.752
#2	.01485	9.1865	.10103	.14014	.23593	.00044	18.701
#3	.01032	9.3147	.13166	.10431	.24034	.00044	18.879
#4	.01183	9.2249	.05696	.10571	.23674	.00044	18.821
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	2.0000	600.00	100.00	100.00	100.00	15.000	600.00
Low	01000	20000	30000	20000	20000	00500	-5.0000
Elem	CD	CO	CR	CU	FE	K_	LI
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	.00378	.00336	.41297	.70998	5.1706	6.3550	.03321
SDev	.00456	.00344	.00382	.00485	.0251	.1289	.00038
%RSD	120.76	102.45	.92485	.68288	.48508	2.0282	1.1578
#1	.00714	.00552	.41682	.71021	5.1717	6.2938	.03275
#2	.00472	.00553	.41568	.70559	5.1415	6.5388	.03304
#3	00290	00170	.40998	.71667	5.2026	6.3428	.03353
#4	.00615	.00408	.40941	.70744	5.1667	6.2448	.03353
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	100.00	100.00	100.00	100.00	400.00	1000.0	20.000
Low	00500	05000	01000	02500	10000	-5.0000	05000
Elem	MG	MN	MO	NA	NI	PB	SB
Units	ppm	ppm	ppm	ppm	ppm	ppm	ppm
Avge	12.158	.20790	.01071	8.2136	.20427	.38543	.00606
SDev	.065	.00052	.00248	.1455	.00807	.02225	.00530
%RSD	.53209	.25077	23.196	1.7720	3.9519	5.7716	87.447
#1	12.152	.20764	.01443	8.1654	.19252	.37972	.00326
#2	12.119	.20763	.00946	8.0843	.20693	.40287	.00351
#3	12.252	.20868	.00947	8.4223	.20671	.35635	.00346
#4	12.111	.20764	.00947	8.1824	.21093	.40281	.01400
Errors	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass	LC Pass
High	600.00	100.00	50.000	400.00	100.00	100.00	100.00
Low	-5.0000	01500	04000	-5.0000	04000	10000	06000
Elem Units Avge SDev %RSD	SE ppm 01592 .03160 198.49	SI ppm 18.661 .080	SN ppm .06847 .02399 35.043	SR ppm .11646 .00074 .63871	TI ppm .08297 .00033 .40018	TL ppm .04322 .04956 114.66	V_ ppm .00876 .00005 .60910
#1	06168	18.625	.07415	.11693	.08271	01760	.00884
#2	00210	18.577	.03690	.11558	.08341	.06377	.00874
							C O O '

## FINAL PAGE

## **PART I**

## **ADMINISTRATIVE RECORD**

**PART I** 

FINAL PAGE