



THE MEMPHIS DEPOT TENNESSEE

ADMINISTRATIVE RECORD COVER SHEET

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UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

REGION IV

345 COURTLAND STREET, N.E.
ATLANTA, GEORGIA 30365

AUG 24 1994

4WD-FFA

Commander

Attn: DDMT-DE (Frank Novitski)
Defense Distribution Depot Memphis
2163 Airways Blvd.
Memphis, Tennessee 38114

SUBJ: Defense Distribution Depot, Memphis, Tennessee (DDMT)
Response to EPA Comments, May 3, 1994

Dear Mr. Novitski:

In a May 3, 1994 letter to Mr. Joseph Franzmathes, DDMT submitted its response to the comments EPA had on the following four documents:

- o Generic Remedial Investigation/Feasibility Study Work Plan;
- o Generic Quality Assurance Plan;
- o Generic Health and Safety Plan;
- o Operable Unit 1 Field Sampling Plan.

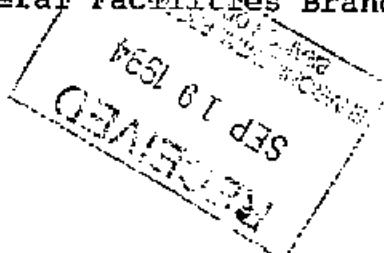
EPA's comments were enclosed in a March 28, 1994 letter from EPA to DDMT. The vast majority of the responses to EPA's comments were satisfactory. However, there were three comments that need to be addressed further. These comments are enclosed with this letter. If you have any questions or comments, please contact me at 404/347-3016, vmx. 6431.

Sincerely,

Martha Befry
Remedial Project Manager
Federal Facilities Branch

Enclosure

cc: Jordan English, TDEC
Christine Kartman, DDMT-DE



RECEIVED
DDMT-DE
AUG 29 1994
C.G. 541-4600

EPA COMMENTS

GENERIC RI/FS WORK PLAN, OPERABLE UNIT 1 FIELD SAMPLING PLAN
DDMT RESPONSE TO EPA COMMENTS, MAY 3, 1994
DEFENSE DISTRIBUTION DEPOT
MEMPHIS, TENNESSEE

GENERIC RI/FS WORK PLAN

EPA Comment #24: Apparently, the intent of this comment was misunderstood. Information regarding the condition of the wells in Allen field should be gathered regardless of whether or not additional offsite monitor wells are to be installed. This information, if available, may have a major impact on the decision regarding necessity and/or placement of additional monitor wells.

EPA Comment #25: Again, the intent of the original comment was misunderstood. EPA is concerned about chlorinated solvents in the groundwater, not in the halogenated by-products produced by treatment of water supplies by chlorine.

EPA Comment # 37: It is stated that preliminary remediation goals (PRGs) which were developed by EPA Region IX were used in the document. This guidance should not be used for PRGs in Region IV. Instead, EPA Region IV recommends that Region III's Risk-Based Concentration Table (updated quarterly) be used for determining PRBs at the site. Region III has developed a table for selecting contaminants of potential concern (COPCs). This table should be used for selecting COPCs and/or as a chemical screening guidance. The table is titled Selecting Exposure Routes and Contaminants of Concern by Risk-Based Screening. A copy of this has been attached to these comments.



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

Region III

341 Chestnut Street
Philadelphia, Pennsylvania 19107

July 11, 1994

SUBJECT: Risk-Based Concentration Table, Third Quarter 1994

FROM: Roy L. Smith, Ph.D., Senior Toxicologist
Technical Support Section (3HW13)

TO: RBC Table mailing list

Roy L. Smith

Attached is the EPA Region III risk-based concentration (RBC) table, which we have distributed quarterly to all interested parties since 1991. If you are not currently on the mailing list, but would like to be, please contact Anna Poulton (phone: 215-597-3179, fax: 215-597-9890) and give her your name, address, and phone and fax numbers.

The table contains reference doses and carcinogenic potency slopes (obtained from IRIS through July 1, 1994, HEAST through November March 1994, the Superfund Health Risk Technical Support Center, and other EPA sources) for nearly 600 chemicals. These toxicity constants have been combined with "standard" exposure scenarios to calculate RBCs - chemical concentrations corresponding to fixed levels of risk (i.e., a hazard quotient of 1, or lifetime cancer risk of 10^{-6} , whichever occurs at a lower concentration) in water, air, fish tissue, and soil.

The Region III toxicologists use the table to screen sites not yet on the NPL, respond rapidly to citizen inquiries, and spot-check formal baseline risk assessments. The background materials provide the complete basis for all the calculations, with the intent of showing users exactly how the RBCs were developed. Simply put, RBCs are risk assessments run in reverse. For a single contaminant in a single medium, under standard default exposure assumptions, the RBC corresponds to the target risk or hazard quotient.

The calculations also have several important limitations. Specifically excluded from consideration are (1) transfers from soil to air and groundwater, and (2) cumulative risk from multiple contaminants or media. Also, the toxicity information in the table has been assembled by hand, and (despite extensive checking and years of use) may contain errors. It's advisable to cross-check before relying on any RfDs or CPSs in the table. If you find any errors, please send me a note.

Lately, many callers have asked whether the risk-based concentrations can be used as valid no-action levels or cleanup levels, especially for soils. The answer is a bit complex. First, it is important to realize that the RBC table does not constitute regulation or guidance, and should not be viewed as a substitute for a site-specific risk assessment. For sites where:

1. A single medium is contaminated;
2. A single contaminant contributes nearly all of the health risk;

EPA Region III Risk-Based Concentrations: R.L. Smith (7/11/94)

3. Volatilization or leaching of that contaminant from soil is expected not to be significant;

The exposure scenarios used in the RBC table are appropriate for the site;

5. The fixed risk levels used in the RBC table are appropriate for the site; and

6. Risk to ecological receptors is expected not to be significant;

the risk-based concentrations would probably be protective as no-action levels or cleanup goals. However, to the extent that a site deviates from this description, as most do, the RBCs would not necessarily be appropriate.

To summarize, the table should generally not be used to (1) set cleanup or no-action levels at CERCLA or RCRA Corrective Action sites, (2) substitute for EPA guidance for preparing baseline risk assessments, or (3) determine if a waste is hazardous under RCRA.

This issue of the RBC table includes new toxicity constants and media concentrations, which are marked on the table in underlined boldface print. On pg. 4, the source of the toxicological constants labeled "ECAO-Cincinnati" has been clarified.

I get many telephone calls about the RBC table, but am often unable to answer the phone. Many of you have the same problem, so we play a lot of "phone tag". Last quarter, I suggested that you fax me (at 215-597-9890) your technical questions and concerns, so I could respond by return fax. This has worked very well for me, and I hope you have been satisfied with my responses. I would like to continue this method. Of course, if you don't have access to a fax machine, I will also continue to respond to voice mail messages.

Attachment

Risk-Based Concentration Table Background Information

General: Separate carcinogenic and non-carcinogenic risk-based concentrations were calculated for each compound for each pathway. The concentration in the table is the lower of the two, rounded to two significant figures. The following terms and values were used in the calculations:

Exposure variables	Value	Name
1-General:		
Carcinogenic potency slope oral (risk per mg/kg/d):	*	CPS _O
Carcinogenic potency slope inhaled (risk per mg/kg/d):	*	CPS _I
Reference dose oral (mg/kg/d):	*	RfDo
Reference dose inhaled (mg/kg/d):	*	RfDi
Target cancer risk:	1e-06	TR
Target hazard quotient:	1	THQ
Body weight, adult (kg):	70	BWa
Body weight, age 1-6 (kg):	15	BWc
Averaging time carcinogens (d):	25550	ATc
Averaging time non-carcinogens (d):	ED*365	ATn
Inhalation, adult (m ³ /d):	20	IRAb
Inhalation, child (m ³ /d):	12	IRAc
Inhalation factor, age-adjusted (m ³ -y/kg-d):	11.66	IFAdj
Tap water ingestion, adult (L/d):	2	IRWa
Tap water ingestion, age 1-6 (L/d):	1	IRWc
Tap water ingestion factor, age-adjusted (L-y/kg-d):	1.09	IFWadj
Fish ingestion (g/d):	54	IRF
Soil ingestion, adult (mg/d):	100	IRS _a
Soil ingestion, age 1-6 (mg/d):	200	IRS _c
Soil ingestion factor, age adjusted (mg-y/kg-d):	114.29	IFSadj
2-Residential:		
Exposure frequency (d/y):	350	EFr
Exposure duration, total (y):	30	EDtot

Exposure variables	Value	Name
Exposure duration, age 1-6 (y)	6	EDc
Volatilization factor (L/m ³)	0.5	VP
3-Occupational:		
Exposure frequency (d/y)	250	EPo
Exposure duration (y)	25	EDo
* - Contaminant-specific toxicity parameters		

The priority among sources of toxicological constants was as follows: (1) IRIS, (2) HEAST, (3) HEAST alternative method, (4) EPA Superfund Health Risk Technical Support Center, (5) withdrawn from IRIS or HEAST, and (6) other EPA documents. Each source was used only if numbers from higher-priority sources were unavailable. The EPA Superfund Health Risk Technical Support Center, part of the Chemical Mixtures Branch of ECAO-Cincinnati, develops provisional RfDs and CPSSs on request for contaminants not in IRIS or HEAST. These provisional values are labeled "e = EPA-ECAO provisional" in the table. It is possible they may be obsolete. If one of the "e" constants is important to a Superfund risk assessment, consider requesting, through a Regional risk assessor, a new provisional value.

Algorithms:

1. Age-adjusted factors: Because contact rates with tap water, ambient air, and residential soil are different for children and adults, carcinogenic risks during the first 30 years of life were calculated using age-adjusted factors. These factors approximated the integrated exposure from birth until age 30 by combining contact rates, body weights, and exposure durations for two age groups - small children and adults. The age-adjusted factor for soil was obtained from RAGS IB; the others were developed by analogy.

a. Air inhalation ([m³· y]/[kg· d]):

$$IFAdj_{air} = \frac{EDc \cdot IRAc}{BWc} + \frac{(EDtot - EDc) \cdot IRAa}{BWA}$$

b. Tap water ingestion ([L· y]/[kg· d]):

$$IFWadj = \frac{EDc \cdot IRWc}{BWc} + \frac{(EDtot - EDc) \cdot IRWa}{BWA}$$

c. Soil ingestion ([mg·y]/[kg·d]):

$$\text{IFSadj} = \frac{EDc \cdot IRS_c}{BWC} + \frac{(ED_{tot} - EDc) \cdot IRS_a}{BW_a}$$

2. Residential water use ($\mu\text{g/L}$). Volatilization terms were calculated only for compounds with "****" in the "VOC" column. Compounds having a Henry's Law constant greater than 10^5 were considered volatile. The list may be incomplete, but is unlikely to include false positives. The equations and the volatilization factor (VF, above) were obtained from RAGS IB. Oral potency slopes and reference doses were used for both oral and inhaled exposures for volatile compounds lacking inhalation values. Inhaled potency slopes were substituted for unavailable oral potency slopes only for volatile compounds; inhaled RfDs were substituted for unavailable oral RfDs for both volatile and non-volatile compounds.

a. Carcinogens: Calculations were based on combined childhood and adult exposure.

$$\frac{TR \cdot AT_c \cdot 1000 \frac{\text{ug}}{\text{mg}}}{EF_r \cdot ([VF \cdot IFA_{adj} \cdot CPS_i] + [IFW_{adj} \cdot CPS_o])}$$

b. Non-carcinogens: Calculations were based on adult exposure.

$$\frac{THQ \cdot BW_a \cdot AT_n \cdot 1000 \frac{\text{ug}}{\text{mg}}}{EF_r \cdot ED_{tot} \cdot \left(\frac{VF \cdot IRA_a}{RfDi} + \frac{IRW_a}{RfDo} \right)}$$

3. Air ($\mu\text{g/m}^3$). Oral potency slopes and references were used where inhalation values were not available.

a. Carcinogens: Calculations were based on combined childhood and adult exposure.

$$\frac{TR \cdot AT_c \cdot 1000 \frac{\text{ug}}{\text{mg}}}{EF_r \cdot IFA_{adj} \cdot CPS_i}$$

b. Non-carcinogens: Calculations were based on adult exposure.

$$\frac{THQ \cdot RfDi \cdot BW_a \cdot AT_n \cdot 1000 \frac{\text{ug}}{\text{mg}}}{EF_r \cdot ED_{tot} \cdot IRA_a}$$

EPA Region III Risk-Based Concentrations: R.L. Smith (7/11/94)

4. Fish (mg/kg):

a. Carcinogens: Calculations were based on adult exposure.

$$\frac{TR \cdot BWa \cdot ATc}{EFr \cdot EDtot \cdot \frac{IRF}{1000} \cdot CPSo}$$

b. Non-carcinogens: Calculations were based on adult exposure.

$$\frac{THQ \cdot RIDo \cdot BWa \cdot ATa}{EFr \cdot EDtot \cdot \frac{IRF}{1000} \cdot CPSo}$$

5. Soil commercial/industrial (mg/kg): The default exposure assumption that only 50% of incidental soil ingestion occurs at work has been omitted. Calculations were based on adult occupational exposure.

a. Carcinogens:

$$\frac{TR \cdot BWa \cdot ATc}{EFo \cdot EDo \cdot \frac{IRSg}{10^6} \cdot CPSo}$$

b. Non-carcinogens:

$$\frac{THQ \cdot RIDo \cdot BWa \cdot ATa}{EFo \cdot EDo \cdot \frac{IRSg}{10^6} \cdot CPSo}$$

6. Soil residential (mg/kg):

a. Carcinogens: Calculations were based on combined childhood and adult exposure.

$$\frac{TR \cdot ATc}{EFr \cdot \frac{IFSadj}{10^6} \cdot CPSo}$$

b. Non-carcinogens: Calculations were based on childhood exposure only.

$$\frac{THQ \cdot RIDo \cdot BWc \cdot ATD}{EFr \cdot EDc \cdot \frac{IRSc}{10^6} \cdot CPSo}$$

Source: ITRIS; h=HEAST; a=HEAT; n=Information from ITRIS or HEAT; e=Other EPA documents

Parts of RBC: c=carcinogenic effects; a=noncarcinogenic effects

Contaminant	CAS	RDX mg/kg/d	RDX mg/kg/d	CPSo mg/kg/d	CPSo mg/kg/d	Tap water mg/L	Ambient air mg/m ³	Fish mg/kg	Industrial soil mg/kg	Residential soil mg/kg
Acetophenone	30560191	4.00E-03 /	2.57E-03 /	8.70E-03 /	7.70E-03 /	7.7 a	0.72 a	0.36 a	130 a	72 a
Acetonechloride	34256821	2.00E-02 /				730 a	73 a	27 a	20000 a	16000 a
Acetone	67641	1.00E-01 /	2.86E-03 a			3700 a	370 a	140 a	100000 a	7800 a
Acetone cyanohydrin	75263	7.00E-02 h				2600 a	10 a	95 a	72000 a	51000 a
Acetonitrile	73078	6.00E-03 /	1.43E-02 a			230 a	92 a	8.1 a	56100 a	47000 a
Acetophenone	98362	1.00E-01 /	5.71E-05 a			600 a	0.042 a	0.021 a	140 a	100000 a
Acid fluorfen	62476399	1.10E-02 /				470 a	47 a	18 a	13000 a	10000 a
Aerolein	167023	2.00E-02 h	1.71E-06 /			730 a	0.071 a	27 a	20000 a	16000 a
Acrylamide	75061	2.00E-04 /	4.50E-00 /	4.35E+00 /	4.35E+00 /	5500 a	0.015 a	0.0014 a	0.0007 a	0.0014 a
Acrylic acid	79107	5.00E-01 /	1.00E-03 /			16000 a	3.7 a	600 a	510000 a	390000 a
Acrylonitrile	107111	1.00E-03 h	1.71E-04 /	5.40E-01 /	2.38E-01 /	612 a	0.025 a	0.0038 a	1.3 a	12 a
Alechlor	15972608	1.00E-02 /	8.00E-02 h			634 a	0.073 a	0.039 a	36 a	8 a
Ale	1596845	1.50E-01 /				5500 a	550 a	200 a	130000 a	120000 a
Aldicarb	1160053	1.00E-03 /				37 a	37 a	14 a	1000 a	78 a
Aldicarb sulfone	1645894	1.00E-03 /				160 a	18 a	14 a	1000 a	78 a
Aldrin	3090062	3.00E-03 /				1800 a	1 a	6.8 a	31000 a	39000 a
Allyl	74723646	2.50E-01 /				910 a	910 a	340 a	250000 a	200000 a
Allyl alcohol	107186	5.00E-03 /				160 a	18 a	6.8 a	5100 a	3900 a
Allyl chloride	107051	5.00E-02 a	2.86E-04 /			330 a	33 a	12 a	9200 a	7000 a
Aluminum phosphide	20919738	4.00E-04 /				2600 a	260 a	95 a	72000 a	55000 a
Andro	67485294	3.00E-04 /				11 a	1.1 a	0.41 a	310 a	23 a
Ametryn	834128	9.00E-03 a				330 a	33 a	12 a	9200 a	7000 a
m-Aminophenol	391275	7.00E-02 h				1000 a	100 a	34 a	250000 a	390000 a
4-Aminopyridine	404245	2.00E-03 h				673 a	0.073 a	0.027 a	220 a	160 a
Anilurez	33089511	2.50E-03 /				91 a	9.1 a	3.4 a	25000 a	20000 a
Ammonia	76644117		2.66E-02 /			1000 a	100 a	34 a	410 a	31 a
Ammonium sulfamate	7773050	2.00E-01 /				7300 a	730 a	270 a	200000 a	160000 a
Aniline	62133		2.86E-04 /	1.70E-03 /		10 a	1 a	0.59 a	500 a	110 a
Antimony and compounds	74407580	4.00E-04 /				15 a	15 a	0.54 a	410 a	31 a
Antimony pentoxide	1314609	5.00E-04 h				118 a	14 a	0.68 a	510 a	39 a
Antimony potassium tartrate	504610	9.00E-04 h				130 a	13 a	1.2 a	920 a	70 a
Antimony tetroxide	1332216	4.00E-04 h				15 a	15 a	0.54 a	410 a	31 a
Antimony trioxide	1309644	4.00E-04 h				115 a	13 a	0.54 a	410 a	31 a
Apollo	74115245	1.30E-02 h				970 a	47 a	18 a	13000 a	10000 a
Arsenite	1403578	5.00E-02 h		2.93E-02 /	2.49E-02 /	27 a	0.29 a	0.13 a	110 a	26 a
Arsenic	7440392	3.00E-04 /		1.71E-00 /	1.51E-01 /	11 a	1.1 a	0.41 a	310 a	23 a
Arsenic (as carcinogen)	7440392		1.43E-03 /			0.938 a	0.00041 a	0.0018 a	15 a	0.37 a
Arsine	76377148	9.00E-03 /				330 a	33 a	12 a	52000 a	700 a
Asulam	3337711	3.00E-02 /				1600 a	160 a	63 a	31000 a	3900 a
Atrazine	1912249	3.50E-02 /		2.22E-01 h		0.3 a	0.028 a	0.014 a	13 a	2.9 a

Source: I=IRIS h=HEAT a=HEAT all, w=Findchem from IRIS or HEAT, r=IRP/RCRA provision or Other EPA documents

Contaminant	RID#	RID# in AP4d	CAS#	CR30 mg/kg/d	CP31 kg/during ingestion	V IC	N Tef. value	Ambient air	Industrial air	Industrial soil	Residential soil
						ug/m3	ug/m3	ug/m3	ug/m3	ug/m3	ug/m3
Avermectin B1	63192553	4.00E-04 /	1.10E-01 /	1.08E-01 /	0.61 e	0.033 e	0.029 e	0.34 n	410 n	31 n	31 n
Azobenzene	103533	7.00E-02 /	1.43E-04 *	1.10E-01 /	2500 n	0.52 n	0.52 n	0.52 n	26 n	3.8 n	3.8 n
Berium and compounds	7440193	4.00E-03 /	1.43E-04 *	1.10E-01 /	1100 n	1.10 n	1.10 n	1.10 n	7200 n	5300 n	5300 n
Beryllium	114261	4.00E-03 /	3.00E-02 /	1.10E-01 /	910 n	0.91 n	0.91 n	0.91 n	410 n	310 n	310 n
Beythroid	43121433	3.00E-02 /	2.39E-02 /	1.10E-01 /	11000 n	1100 n	1100 n	1100 n	31000 n	2300 n	2300 n
Benfumit	1861401	3.00E-01 /	1.00E-01 /	1.11E-01 /	610 n	0.61 n	0.61 n	0.61 n	2600 n	2000 n	2000 n
Benomyl	17804152	3.00E-02 /	2.10E-03 /	1.10E-01 /	1800 n	180 n	180 n	180 n	51000 n	3900 n	3900 n
Bentazon	25057890	1.00E-03 /	1.00E-01 /	1.10E-01 /	91 n	0.91 n	0.91 n	0.91 n	2600 n	200 n	200 n
Benzaldehyde	100327	1.00E-01 /	1.11E-01 /	1.11E-01 /	2.90E-02 /	0.36 e	0.22 e	0.22 e	140 n	100000 n	23000 n
Benzene	71132	1.00E-01 /	1.00E-01 /	1.10E-01 /	1.30E+01 /	0.17 e	0.037 n	0.037 n	0.11 e	99 n	22 n
Benzethiol	1018785	1.00E-01 /	3.00E-03 /	1.10E-01 /	2.35E+02 /	0.00029 e	0.00027 e	0.00027 e	0.014 n	10 n	0.71 n
Benzidine	92873	3.00E-03 /	1.00E+00 /	1.10E-01 /	150000 n	14000 n	14000 n	14000 n	3400 n	310000 n	30000 n
Benzoic acid	654530	4.00E+00 /	98977	1.30E+01 /	1.30E+01 /	0.00352 e	0.00048 e	0.00024 e	0.22 e	0.22 e	0.039 n
Benzotrichloride	1003116	3.00E-01 /	1.00E-01 /	1.10E-01 /	11000 n	1100 n	1100 n	1100 n	310000 n	310000 n	310000 n
Benzyl alcohol	100447	1.00E-01 /	1.00E-01 /	1.10E-01 /	0.062 e	0.037 e	0.037 e	0.037 e	0.19 n	117 n	3.8 n
Benzyl chloride	7440417	1.00E-03 /	4.30E+00 /	1.10E-01 /	0.40E+00 /	0.016 e	0.00075 e	0.00073 e	0.057 e	0.057 e	0.13 n
Beryllium and compounds	1411662	1.00E-04 /	1.50E-02 /	1.10E-01 /	150 n	150 n	150 n	150 n	100 n	78 n	78 n
Bidrin	126137043	1.1-Biphenyl	572324	1.00E-02 /	7.00E-02 /	1.30E-02 /	0.26 e	0.18 e	0.043 e	3.41 e	9.1 e
Bis(2-chloroisopropyl)ether	396339129	4.00E-02 /	142881	2.20E+02 /	2.17E+02 /	0.000049 e	0.000029 e	0.000014 e	0.013 e	0.0029 e	0.0029 e
Bis(chloromethyl)ether	7440428	9.00E-02 /	7637072	1.00E-02 /	7.00E-02 /	0.96 e	0.095 e	0.049 e	0.41 e	9.1 e	9.1 e
Bis(2-chloro-1-methylethyl)ether	117817	2.00E-02 /	111444	1.40E-02 /	1.16E+00 /	4.8 e	0.45 e	0.23 e	200 e	46 e	46 e
Bis(chloromethyl)ether	80057	1.00E-02 /	73274	2.00E-02 /	1.10E-04 /	0.0092 e	0.0054 e	0.0029 e	0.26 e	0.58 e	0.58 e
Bisphenol A	80057	1.00E-02 /	7440428	9.00E-02 /	7.71E-03 n	1300 n	1.21 n	1.21 n	120 n	92000 n	7000 n
Boron (and borates)	7637072	2.00E-02 /	7440428	2.00E-04 n	6.20E-02 /	0.17 e	0.17 e	0.17 e	0.73 n	140 n	110 n
Boron trifluoride	75322	2.00E-02 /	101133	1.71E-03 n	1.10E-01 /	0.096 e	0.057 e	0.057 e	1.6 e	0.4 e	0.4 e
Bromodichloromethane	1591692	1.00E-02 /	1689945	1.80E-02 /	1.90E-03 /	3.35E-03 /	2.4 e	1.6 e	0.4 e	360 n	91 n
Bromoethene	2104963	5.00E-03 n	1689945	1.40E-03 /	1.43E-03 /	1.10E-01 /	0.096 e	0.057 e	0.68 n	1100 n	350 n
Bromoform (tetrachloromethane)	74839	1.40E-03 /	1065990	1.00E-01 /	9.80E-01 /	730 n	73 n	73 n	270 n	200000 n	16000 n
Bromophos	1689992	2.00E-02 /	71363	1.00E-01 /	1.00E-01 /	0.011 e	0.0064 e	0.0064 e	140 n	100000 n	16000 n
Bromoxynil	1065990	1.00E-01 /	2008415	5.00E-02 /	5.00E-02 /	3700 n	180 n	180 n	68 n	91000 n	3900 n
1,3-Butadiene	71363	1.00E-01 /	1359918	1.00E-02 /	61 n	61 n	61 n	61 n	14 n	10000 n	780 n
1-Butanol	83687	2.00E-01 /	2008415	5.00E-02 /	5.00E-02 /	7300 n	73 n	73 n	270 n	200000 n	16000 n
Butylbenzyl phthalate	1065990	1.00E-01 /	1359918	1.00E-02 /	61 n	61 n	61 n	61 n	14 n	10000 n	780 n
sec-Butylbenzene											

Source: 1=IRIS; 2=HEAST; 3=HEAST alt. - withdrawn from IRIS or HEAST; 4= EPA-RCRA prioritized or Other EPA documents

Contaminant	CAS	RDI			CPSD			CFSI			Ambient Air			Industrial Fugitive			Residential Soil		
		mg/kg/d	mg/kg/d	mg/kg/d	kg/d/mg	kg/d/mg	kg/d/mg	C	kg/d/mg	kg/d/mg	C	kg/d/mg	kg/d/mg	C	kg/d/mg	kg/d/mg	C	kg/d/mg	kg/d/mg
tert-Butylbenzene	104518	1.00E-02 *							61 n		37 n		14 n		10000 n		780 n		
Butylphthalyl butylglycolate	65701	1.00E+00 /							3700 n		3700 n		1400 n		100000 n		78000 n		
Casodytic acid	75605	3.00E-03 n							110 n		11 n		41 n		3100 n		230 n		
Cadmium and compounds	7440439	5.00E-04 /							13 n		0.00099 e		0.68 n		510 n		39 n		
Caprolactam	103602	5.00E-01 /							1800 n		1800 n		680 n		91000 n		3900 n		
Caprolactol	2425061	2.00E-03 /							73 n		0.73 n		0.37 n		330 n		74 n		
Captan	113062	1.30E-01 /							19 n		1.8 e		0.9 e		820 n		180 n		
Carberry	637252	1.00E-01 /							3700 n		3700 n		140 n		100000 n		7800 n		
Carbofuran	85748								34 e		0.21 e		0.16 e		140 n		32 n		
Carbon disulfide	1563662	5.00E-03 /							180 n		18 n		6.8 n		3100 n		190 n		
Carbon tetrachloride	75150	1.00E-01 /							21 n		10 n		140 n		100000 n		7800 n		
Carbosulfan	56235148	1.00E-02 /							0.16 e		0.12 e		0.024 e		22 n		4.9 n		
Carboxan	5224684	1.00E-01 /							3700 n		3700 n		14 n		100000 n		780 n		
Chloral	75876	2.00E-03 /							73 n		73 n		27 n		82000 n		160 n		
Chloramben	133904	1.50E-02 /							150 n		15 n		20 n		10000 n		1200 n		
Chloranil	118752								0.17 e		0.016 e		0.0074 e		71 n		1.6 n		
Chlordane	17749	6.00E-01 /							0.032 e		0.009 e		0.0024 e		22 n		0.49 n		
Chlorinated-ethyl	50982324	2.00E-02 /							730 n		73 n		27 n		200000 n		16000 n		
Chlorine	7781294	1.00E-01 /							2700 n		270 n		140 n		100000 n		7800 n		
Chlorine dioxide	1035944								21 n		0.21 n		0.15 n		71 n		5 n		
Chloroacetaldehyde	107206	6.90E-03 *							250 n		25 n		9.3 n		7100 n		540 n		
Chloroacetic acid	79118	2.00E-03 n							73 n		73 n		27 n		2000 n		160 n		
2-Chloracetophenone	512274								0.31 n		0.031 n		0.011 n		3 n		1 n		
4-Chloromilinc	106478	4.00E-03 /							150 n		15 n		1.4 n		3100 n		310 n		
Chlorobenzene	108507	2.00E-02 /							39 n		21 n		27 n		20000 n		1600 n		
Chlorobenzilate	510156	2.00E-02 /							0.25 e		0.023 e		0.012 e		11 n		2.4 n		
p-Chlorobenzoic acid	74113	2.00E-01 n							730 n		73 n		27 n		200000 n		16000 n		
4-Chloroboron trifluoride	98566	2.00E-02 n							730 n		73 n		27 n		200000 n		16000 n		
2-Chloro-1,3-butadiene	126998	2.00E-02 *							14 n		7.3 n		27 n		20000 n		16000 n		
1-Chlorobutane	109693	4.00E-01 n							2400 n		1300 n		940 n		410000 n		31000 n		
Chlorodifluoromethane	73456								52000 n		52000 n		340 n		2410000 n		31000 n		
Chloroethane	75003	4.00E-01 *							2600 n		10000 n		340 n		2410000 n		31000 n		
2-Chloromethyl vinyl ether	110734	2.50E-02 *							1.05E-02 n		1.05E-02 n		0.113 e		34 n		2000 n		
Chloroform	67663	1.00E-02 /							1.30E-02 n		1.30E-02 n		0.074 e		0.57 e		220 n		
Chloromethane	74873								1.4 n		0.99 e		0.24 e		170 n		49 n		
4-Chloro-2,2-methylaniline hydrochloride	3165933								0.15 e		0.014 e		0.0059 e		52 e		1.4 e		
4-Chloro-2-methylaniline	95692								0.12 e		0.011 e		0.0034 e		49 e		1.1 e		
Beta-Chloronaphthalene	91387	3.00E-02 /							2.50E-02 n		5.80E-01 n		0.002 n		160 n		35 n		
o-Chloronitrobenzene	18733								1.42 E-02 n		1.42 E-02 n		0.13 e		110 e		26 e		
p-Chloronitrobenzene	100092								1.87 E-02 n		1.87 E-02 n		0.18 e		160 n		35 n		
2-Chlorophenol	91371	5.00E-03 /							1.00E-02 n		1.00E-02 n		0.002 n		110 n		30 n		

Sources: 1=PRIS 2=HEAST 3=PRIS 4=PRIS or PRAT 5=PRAT 6=PRCA-RCRA prioritized 0=Other EPA documents

Contaminant	CAS	R'DO mg/l/d	R'DI mg/l/d	CPSI	CFSI kg/m ³	V kg/m ³	Basis of RBC: c=concentric effects n=nonconcentric effects		
							Ambient air µg/m ³	Ground water µg/L	Soil mg/kg
2-Chloropropane	7926	1.50E-02 /	2.80E-02 /	1.00E-02 /			100 n	100 n	100 n
Chlorobutanol	1897456	1.50E-02 /	2.80E-02 /	1.00E-02 /			6.1 n	0.57 n	0.29 n
c-Chlorotoluene	92498	2.00E-02 /					120 n	73 n	250 n
Chlorophorm	101213	2.00E-01 /					700 n	270 n	270 n
Chlorpyrifos	2971882	3.00E-03 /					110 n	11 n	41 n
Chlorviniles-methyl	3591130	1.00E-02 /					370 n	37 n	14 n
Chlorsulfuron	64002723	5.00E-02 /					1000 n	180 n	68 n
Chlorthiophos	60234164	4.00E-04 /					29 n	29 n	11 n
Chromium III and compounds	16063631	1.00E+00 /	5.70E-07 /				37000 n	0.0021 n	1400 n
Chromium VI and compounds	7440473	3.00E-03 /		4.20E-01 /			180 n	0.00013 n	6.5 n
Coal tar	8901189		2.20E-00 w					0.0021 n	1000 n
Cobalt	7402484	6.00E-02 *					2200 n	220 n	61 n
Coke Oven Emissions	8007452		2.17E-00 /					61 n	100000 n
Copper anti compounds	7440508	3.71E-02 /					1400 n	140 n	50 n
Crotonaldehyde	123739	1.00E-02 w	1.90E-00 h	1.90E-00 w			0.033 n	0.0031 n	0.0017 n
Cumene	938238	4.00E-02 /	2.57E-03 h				1500 n	9.4 n	3.15 n
Cyanides:									0.34 n
Barium cyanide	542671	1.00E-01 w					3700 n	370 n	140 n
Calcium cyanide	592018	4.00E-02 /					1500 n	150 n	54 n
Copper cyanide	544973	3.00E-03 /					180 n	18 n	6.8 n
Cyanazine	2172462	2.00E-03 h	8.40E-01 h				0.03 n	0.0073 n	0.0018 n
Cyanogen	460195	4.00E-02 /					1500 n	150 n	54 n
Cyanogen bromide	508663	9.00E-02 /					1300 n	130 n	120 n
Cyanogen chloride	508774	3.00E-02 /					1800 n	180 n	68 n
Free cyanide	37125	2.00E-02 /					730 n	73 n	27 n
Hydrogen cyanide	749018	2.00E-02 /					730 n	73 n	27 n
Potassium cyanide	151508	3.00E-02 /					1800 n	180 n	68 n
Potassium silver cyanide	5086616	2.00E-01 /					7300 n	730 n	3100 n
Silver cyanide	3086649	1.00E-01 /					1500 n	150 n	140 n
Sodium cyanide	143319	4.00E-02 /						1500 n	150 n
Zinc cyanide	357211	5.00E-02 /					1800 n	180 n	68 n
Cyclohexanone	108941	1.00E-00 /					30000 n	1800 n	6800 n
Cyclohexamine	108918	2.00E-01 /					7300 n	730 n	270 n
Cyhalothrin/Karnate	68083388	5.00E-03 /					180 n	18 n	6.8 n
Cypermethrin	5215078	1.00E-02 /					370 n	37 n	14 n
Cyromazine	66215278	7.50E-03 /					270 n	27 n	10 n
Dacthal	1861321	5.00E-01 /					1100 n	41 n	10000 n
Dalepon	759920	3.00E-02 /					110 n	41 n	2300 n
Dantol	39513418	1.00E-04 *					11 n	0.63 n	0.10 n
DDD	72348	2.40E-01 /					0.23 n	0.026 n	0.013 n
DDDE	72359	1.40E-01 /					0.2 n	0.018 n	0.0093 n
DDT	50293	3.40E-01 /					0.2 n	0.018 n	0.0093 n

Source: 1=IRIS R-REFST, 2=R-REFST alt., 3=IRIS from IRIS or HE4ST, 4=IRIS from IRIS or HE4ST, 5=Other EPA documents

Contaminant	CAS	RfD mg/kg/d	RfD mg/kg/d	CP50 kg/d/mg	CP50 kg/d/mg	Basis of RfD: <i>c</i> =carcinogenic effects; <i>a</i> =noncarcinogenic effects			Industrial solvent	Residential solvent
						Top water C	Top water C	Ambient air		
Decabromodiphenyl ether	1163195 80653483	1.00E-02 4.00E-03	1.00E-02 4.00E-03	6.10E-02 n	6.10E-02 n	61 n	37 n	14 n	10000 n	700 n
Dieldrin	2503164					1.5 n	0.15 n	0.054 n	41 n	31 n
Diazinon	3334119	9.00E-04	n			0.17 n	0.1 n	0.052 n	47 n	10 n
1,4-Dibromobenzene	106376	1.00E-02	/			33 n	33 n	12 n	1000 n	70 n
Dibromochloromethane	124481	2.00E-02	/	4.40E-02 /	4.40E-02 /	61 n	37 n	14 n	10000 n	700 n
1,2-Dibromo-3-chloropropane	96128	4.71E-03 /	1.40E-03 n	2.42E-03 n	0.044 n	0.21 n	0.0023 n	12 n	10 n	0.46 n
1,2-Dibromoethane	106934	5.71E-03 n	8.50E-01 /	7.70E-01 /	0.0073 n	0.0081 n	0.00037 n	0.034 n	0.0075 n	
Diisobutyl phthalate	84742	1.00E-01	/			3700 n	370 n	140 n	100000 n	7000 n
Dicamba	191889	3.00E-02	/			1100 n	110 n	41 n	31000 n	2300 n
1,2-Dichlorobenzene	95501	9.00E-02	/	5.71E-02 n	5.71E-02 n	370 n	210 n	120 n	92000 n	7000 n
1,3-Dichlorobenzene	341731	8.90E-02	n			540 n	320 n	120 n	91000 n	7000 n
1,4-Dichlorobenzene						0.44 n	0.26 n	0.13 n	120 n	27 n
1,3,5-Dichlorobenzene	106467			2.29E-01 /	2.40E-02 n	0.15 n	0.014 n	0.007 n	64 n	1.4 n
1,4-Dichloro-2-butene	91941			4.30E-01 /	9.30E-00 n	0.0011 n	0.00067 n	0.00037 n	0.00037 n	
Dichlorodifluoromethane	73718	2.00E-01	/	3.71E-02 n	3.71E-02 n	300 n	210 n	270 n	200000 n	16000 n
1,1-Dichloroethane	75343	1.00E-01	n	1.43E-01 n	2.88E-03 n	310 n	320 n	140 n	100000 n	7000 n
1,2-Dichloroethane (EDC)	107062	1.00E-02	n	9.10E-02 /	9.10E-02 /	0.12 n	0.069 n	0.035 n	31 n	7 n
1,1-Dichloroethylene	75354	9.00E-03	/	6.00E-03 /	1.73E-01 /	0.044 n	0.036 n	0.0053 n	14.8 n	1.1 n
1,2-Dichloroethylene (cis)	156392	1.00E-02	n			61 n	37 n	14 n	10000 n	700 n
1,2-Dichloroethylene (trans)	156605	2.00E-02	/			120 n	73 n	27 n	20000 n	16000 n
1,2-Dichloroethylene (mixture)	340590	9.00E-03	n			53 n	33 n	12 n	9200 n	700 n
2,4-Dichlorophenol	1208332	3.00E-03	/			110 n	11 n	4 n	3100 n	230 n
2,4-Dichlorophenoxyacetic Acid (2,4-D)	94717	1.00E-02	/			61 n	37 n	14 n	10000 n	700 n
4-(2,4-Dichlorophenoxy)butyric Acid	94826	8.00E-03	/			250 n	29 n	11 n	3200 n	630 n
1,2-Dichloropropane	78873	1.14E-03	/	6.80E-02	6.80E-02	0.16 n	0.092 n	0.046 n	42 n	9.4 n
2,3-Dichloropropanol	6162239	3.00E-03	/			110 n	11 n	4 n	3100 n	230 n
1,3-Dichloropropene	342736	3.00E-04	/	1.71E-03 /	1.71E-01 n	0.077 n	0.046 n	0.018 n	16 n	3.7 n
Dichlorvos	622737	5.00E-04	/	1.43E-04 /	2.90E-01 /	0.23 n	0.072 n	0.011 n	9.9 n	2.2 n
Dicofol	115322			4.40E-01	4.40E-01	0.19 n	0.014 n	0.0072 n	6.3 n	1.3 n
Dicyclopetadiene	77736	3.00E-02	n	5.71E-03 /	1.61E+01 /	0.42 n	0.21 n	41 n	31000 n	2300 n
Dieldrin	60571	5.00E-05	/	1.43E-03 /	1.60E+01 /	0.0042 n	0.0039 n	0.0022 n	0.018 n	0.004 n
Diesel emissions						29000 n	1100 n	100000 n	160000 n	160000 n
Diethyl phthalate	44652	2.00E-01	/	5.71E-03 n	5.71E-03 n	210 n	21 n	13 n	11000 n	850 n
Diethylene glycol, monobutyl ether	112345	2.00E+00	n			7300 n	7300 n	100000 n	110000 n	110000 n
Diethylene glycol, propylene ether	111900					40 n	40 n	13 n	24000 n	1300 n
Diethylformamide	617845	1.10E-02	n			56 n	32 n	12 n	100000 n	8000 n
Di(2-ethylhexyl)phthalate	103221	6.00E-01	/	1.20E-03 /	4.70E-03 n	0.00014 n	1.30E-06 n	6.70E-07 n	0.00061 n	0.00016 n
Diisobutylstibostrol	56331					730 n	290 n	27 n	20000 n	16000 n
Difenzoquat (Avenge)	43224486	8.00E-02	/			2900 n	110 n	110 n	120000 n	6300 n
Diisobutene	35367383	2.00E-02	/			730 n	73 n	27 n	20000 n	16000 n
Disopropyl methylphosphonate (DIMP)	1445736	8.00E-02	/			2900 n	290 n	110 n	82000 n	63000 n

Source: ITRIS. A=HEAST, a=HEAST only, n=Information from HAZ or HEAST, o=EP4-EC4O prioritizations, o=Other EPA documents.

Contaminant	CAS	RDO mg/kg/d	RDI mg/kg/d	CPS		V C	Ambient air	Indirect soil	Residential soil	Basis of RBC: c=concentric effects, n=nontoxicologic effects.	
				Eq. Dung	Eq. dung					Eq. water	Eq. soil
Dimethylpin...	53290847	2.00E-02 /				730 n	73 n	27 n	20000 n	1600 n	
Dimethoate	60315	2.00E-04 /				73 n	73 n	0.27 n	200 n	16 n	
3,3-Dimethoxybenzidine	119904		1.40E-02 n			4.8 n	0.45 n	0.23 n	200 n	46 n	
Dimethyl phthalate	131113	1.00E-01 n				30000 n	37000 n	14000 n	100000 n	70000 n	
Dimethyl terephthalate	120616	1.00E-01 /				3700 n	370 n	140 n	100000 n	7000 n	
Dimethylamine	124403		1.71E-06 *			0.21 n	0.021 n	0.01 n	100000 n	7000 n	
2,4-Dimethylamine hydrochloride	21436984		1.80E-01 n			0.12 n	0.011 n	0.0054 n	14.9 n	1.1 n	
2,4-Dimethylamine	94681		7.50E-01 n			0.09 n	0.0048 n	0.0042 n	13.8 n	0.83 n	
N,N-Dimethylaniline	121697	2.00E-03 /				73 n	73 n	2.7 n	200 n	160 n	
3,3-Dimethylbenzidine	119937		9.20E+00 n			0.0073 n	0.00068 n	0.00034 n	0.031 n	0.009 n	
N,N-Dimethylformamide	68122	1.00E-01 n	8.57E-03 /			3700 n	31 n	140 n	100000 n	7000 n	
1,1-Dimethylhydrazine	57147		2.60E+00	3.50E+00		0.026 n	0.0018 n	0.0012 n	11.8 n	0.23 n	
1,2-Dimethylhydrazine	940738		3.70E+01 *	3.70E+01 *		0.0016 n	0.00017 n	0.00009 n	24.077 n	0.017 n	
2,4-Dimethylphenol	103679	2.00E-02 /				730 n	73 n	27 n	20000 n	1500 n	
2,6-Dimethylphenol	576261	6.00E-04 /				22 n	22 n	0.81 n	7610 n	47 n	
3,4-Dimethylphenol	99658	1.00E-03 /				37 n	37 n	1.4 n	1000 n	78 n	
1,2-Dinitrobenzene	528290	4.00E-04 n				15 n	15 n	0.34 n	710 n	31 n	
1,4-Dinitrobenzene	99650	1.00E-04 /				3.7 n	0.37 n	0.14 n	100 n	7.8 n	
4,6-Dinitro-o-cresolphenol	100234	4.00E-04 n				73 n	73 n	2.7 n	2000 n	160 n	
2,4-Dinitrophenol	131893	2.00E-03 /				11 n	1.5 n	0.34 n	410 n	31 n	
Dinitrotoluene mixture	41285	2.00E-03 /				73 n	73 n	2.7 n	2000 n	160 n	
2,4-Dinitrotoluene	121142	2.00E-03 /				6.80E-01 /	0.0099 n	0.0092 n	0.0046 n	4.2 n	0.94 n
2,6-Dinitrotoluene	606202	1.00E-03 n				73 n	73 n	2.7 n	2000 n	160 n	
Dinoseb	88917	1.00E-03 /				37 n	3.7 n	1.4 n	1000 n	78 n	
di-n-Octyl phthalate	117840	2.00E-02 n				730 n	73 n	27 n	20000 n	1600 n	
1,4-Dioxane	122911		1.10E-02 /			6.1 n	0.57 n	0.29 n	7260 n	58 n	
Diphenamid	997317	3.00E-02 /				100 n	110 n	41 n	31000 n	2300 n	
Diphenylamine	122394	2.50E-02 /				910 n	91 n	34 n	20000 n	2000 n	
1,2-Diphenylhydrazine	122667		1.00E-01 /	7.70E-01 /		0.034 n	0.001 n	0.0039 n	1.36 n	0.1 n	
Diquat	85007	2.20E-03 /				80 n	8.7 n	3 n	2200 n	170 n	
Direct black 38	1937377		8.60E+00 n			0.0073 n	0.0037 n	0.0037 n	0.33 n	0.074 n	
Direct blue 6	26074462		8.10E+00 n			0.0063 n	0.0037 n	0.0036 n	0.33 n	0.079 n	
Direct brown 95	16071866		9.30E+00 n			0.0072 n	0.0037 n	0.0034 n	0.31 n	0.069 n	
Disulfoton	298044	4.00E-05 /				1.5 n	0.13 n	0.054 n	41 n	1 n	
1,4-Dithiane	595293	1.00E-02 /				370 n	37 n	14 n	10000 n	780 n	
Dijuron	330541	2.00E-03 /				73 n	73 n	2.7 n	2000 n	160 n	
Dodine	2439103	4.00E-03 /				150 n	15 n	3.4 n	2100 n	310 n	
Eadiomififen	115297	6.00E-03 n				220 n	22 n	2.1 n	18100 n	470 n	
Eddothall	145733	2.00E-02 /				730 n	73 n	27 n	20000 n	1600 n	
Endrin	722008	3.00E-04 /				11 n	11 n	0.41 n	310 n	23 n	
Epiclorohydrin	106898	2.00E-03 n	2.86E-04 /	9.90E-03 /		6.1 n	1 n	0.32 n	200 n	63 n	

Source: 1=IRIS; 2=HEAST; 3=HEAST ref.; 4=Information from IRIS or HEAST; 5=EP-AECIO promotional or Other EPA documents

Contaminant	CAS	RfD mg/kg/d	RDH mg/kg/d	CFS6 kg/dung	CFS1 kg/dung	V C	Parts of 86C: carcinogenic effects noncarcinogenic effects			
							Ambient air ppb/hair	Top water ppb/C	Fish mg/kg water	Industrial soil mg/kg
1,2-Epoxybutane	106387	5.00E-03	1.71E-03				210 n	21 n		
Etherben (2-chloroethyl phosphoric acid)	16472370	5.00E-03					100 n	14 n	6.1 n	3100 n
Ethion	563122	3.00E-04					18 n	1.8 n	0.58 n	310 n
2-Ethoxyethanol acetate	111159	1.00E-01					1100 n	110 n	410 n	31000 n
2-Ethoxyethanol	110803	4.00E-01					1800 n	210 n	340 n	41000 n
Ethyl acrylate	140683				4.60E-02	n	1.4 n	0.13 n	0.066 n	1.60 n
EPIC (S-Ethyl diisopropylcarbamate)	749944	2.50E-02					910 n	91 n	34 n	25000 n
Ethyl ether	60297	2.00E-01					1200 n	120 n	270 n	20000 n
Ethyl methacrylate	57632	9.00E-02					3100 n	310 n	310 n	16000 n
Ethyl acetate	141786	9.00E-01					3300 n	3300 n	120 n	92000 n
Ethylbenzene	100414	1.00E-01					1300 n	1300 n	140 n	50000 n
Ethylene cyanoboriran	109724	3.00E-01					11000 n	1100 n	410 n	7500 n
Ethylene diamine	107153	2.00E-02					730 n	73 n	27 n	20000 n
Ethylene glycol	107211	2.00E-00					7300 n	7300 n	2700 n	16000 n
Ethylene glycol, monobutyl ether	111762				5.71E-03	n	210 n	21 n	27 n	160000 n
Ethylen oxide	75218				1.02E-00	n	0.066 n	0.018 n	0.0031 n	2.8 n
Ethylen thimura (ETU)	96457	1.00E-03					0.57 n	0.013 n	0.027 n	5.6 n
Ethyl p-nitrophenyl phenylphosphotriocate	2104553	1.00E-05					0.37 n	0.017 n	0.014 n	10 n
Ethylnitrosourea	759739				1.40E-02	n	0.00048 n	0.000043 n	0.000023 n	0.02 n
Ethylphthalyl ethyl glycolate	84720	3.00E+00					11000 n	11000 n	4100 n	100000 n
Emboss	10120	1.00E-03					250 n	29 n	11 n	8200 n
Fenamiphos	22224926	2.50E-04					9.1 n	0.91 n	0.34 n	260 n
Fluorometharm	2164172	1.30E-02					470 n	47 n	11 n	1300 n
Fluoride	7782414	6.00E-02					2000 n	220 n	81 n	61000 n
Fluoridone	397366604	8.00E-02					2900 n	280 n	110 n	12000 n
Fluprimidol	56423913	2.00E-02					730 n	73 n	27 n	20000 n
Flutolanil	663129613	6.00E-02					2200 n	220 n	81 n	16000 n
Fluvalinate	694929545	1.00E-02					370 n	37 n	14 n	4700 n
Folpet	133073	1.00E-01					19 n	1.8 n	0.9 n	10000 n
Fomesafen	72178028				2.00E-01	n	0.31 n	0.03 n	0.017 n	11 n
Fonofos	944229	2.00E-03					110000 n	11000 n	4100 n	1000000 n
Formaldehyde	50000	2.00E-01					4.51E-01	7300 n	7300 n	2000 n
Formic Acid	64186	2.00E+00					0.0013	0.0013	0.00063	16000 n
Fosetyl-al	39146243	3.00E-00					1.37 n	1.37 n	1.3 n	1600 n
Furan	110009	1.00E-03					0.018 n	0.0016	0.00033	0.73 n
Furoxolidone	67454				3.80E-00	n	0.0013	0.0013	0.00063	0.17 n
Furfural	98011	3.00E-03					110 n	-12 n	4.1 n	13100 n
Furum	5318234	4.00E-04					0.0013	0.0013	0.00063	230 n
Furnexolox	60168050				3.00E-02	n	2.2 n	0.21 n	0.057 n	0.013 n
Gluconate-ammonium	771821822	4.00E-04					15 n	1.5 n	0.54 n	410 n
Oxydialdehydic	765344	4.00E-04					15 n	1.1 n	0.54 n	410 n
Glypoxate	1071836	1.00E-01					3700 n	370 n	140 n	100000 n

Source: 1/23 h - RCAST & RCAST all, w=Hydrogen from IRN or IEST, r=Hydrogen from EPA-5C10 protocol, o=Other EPA documents

Contaminant	CAS#	RDo mp/kg/d	RDo mg/kg/d	CPSo mg/dmg	V kg/dmg	V kg/dmg	Basis of RBC: c=carcinogen, o=other, n=noncarcinogen, e=acute.		
							Ambient soil	Fish soil	Industrial soil
Halocloform-methyl	69106-02	1.00E-01	1.30E-02				1.1 e	0.11 e	51 e
HCH (alpha)	79277773			6.30E-06 /	6.30E-06 /		4.70 e	4.70 e	1300 e
HCH (beta)	319846			1.80E-00 /	1.80E-00 /		0.011 e	0.00099 e	1000 e
HCH (gamma) Lindane	319847	3.00E-04 /					0.037 e	0.0035 e	0.1 e
HCH-technical	548599			1.30E-00 h			0.012 e	0.0043 e	0.34 e
Heptachlor	6018731			1.80E-00 /	1.798E-00 /		0.037 e	0.0035 e	0.49 e
Heptachlor epoxide	76443	5.00E-04 /		4.50E-00 /	4.50E-00 /		0.0223 e	0.0014 e	22 e
Hexabromobiphenyl	1024973	1.30E-03 /		9.10E-00 /	9.10E-00 /		0.0112 e	0.00069 e	0.31 e
Hexachlorobenzene	87821	2.00E-03 /					0.12 n	7.3 n	2.7 n
Hexachlorobutane	118741	6.00E-04 /		1.60E-00 /	1.61E-00 /		0.0066 e	0.0039 e	0.002 e
Hexachlorobutene	87863	2.00E-04 h		7.80E-02 /	7.70E-02 /		0.14 e	0.081 e	0.64 e
Hexachlorocyclopentadiene	77474	7.00E-03 /		2.00E-03 h			0.13 n	0.073 n	0.31 e
Hexachlorodibenz-p-dioxin mixture	19408743			6.20E+03 /	4.55E+03 /		0.000011 e	1.40E-06 e	1.10E-07 e
Hexachloroethane	67721	1.00E-03 /		1.40E-02 /	1.40E-02 /		0.75 e	0.43 e	0.23 e
Hexachlorophene	70304	3.00E-04 /					11 n	1.1 e	0.41 e
Heptahydro-1,3,5-trinitro-1,3,5-triazine	1211624	3.00E-03 /		1.10E-01 /			0.61 e	0.097 e	0.029 e
n-Hexane	110543	6.00E-02 h		3.71E-02 /			600 n	350 n	210 n
Heptazine	51235042	3.10E-02 /					1200 n	120 n	45 n
Hydrazine, hydrazine sulfate	307012			1.00E+00 /	1.71E+01 /		0.022 e	0.00037 e	0.0011 e
Hydrogen chloride	7647010			2.00E-03 /			73 n	73 n	0.93 e
Hydrogen sulfide	7783064	1.00E-03 /		2.51E-04 /			110 n	0.94 n	4.1 n
Hydroquinone	1223119	4.00E-02 h					1500 n	130 n	54 n
Inositol	35394440	1.30E-02 /					470 n	47 n	18 n
Isocitazquin	81339377	2.90E-01 /					9100 n	910 n	340 n
Iprodione	36773497	4.00E-02 /					1500 n	150 n	54 n
Isobutanol	78431	3.00E-01 /					1800 n	1100 n	410 n
Isophorone	78391	2.00E-01 /		9.50E-04 /			71 e	6.6 e	3.3 e
Isopropalin	13820130	1.50E-02 /					450 n	55 n	20 n
Isopropyl methyl phosphonic acid	1832544	1.00E-01 /					3700 n	370 n	140 n
Isotubiten	82558107	5.00E-03 /					1800 n	180 n	68 n
Kepone	143700			1.80E+01 *			0.037 e	0.00033 e	0.00018 e
Lactofen	77501634	2.00E-03 /					73 n	73 n	27 n
Lead (tetracetyl)	78002	1.00E-07 /					0.0037 e	0.00037 e	0.00014 e
Lanumox	339312	2.00E-03 /					730 n	73 n	27 n
Lithium	7439932	2.00E-02 *					3700 n	370 n	140 n
Lindane	121753	2.00E-02 /					18000 n	18000 n	510000 n
Melathion	83046996	2.00E-01 /					730 n	73 n	20000 n
Maleic anhydride	109316	1.00E-01 /					3700 n	370 n	100000 n
Maleic hydrazide	123331	9.00E-01 /					18000 n	600 n	15000 n
Malononitrile	1097772	2.00E-03 h					0.73 n	0.073 n	0.027 n
Mancorez	8018017	3.00E-02 h					1000 n	110 n	41 n
Menelb	12427362	3.00E-03 /					180 n	18 n	6.8 n

Source: I=IRIS A=HEAST C=HEATST d=Withdrawn from IRIS or HEATST e=EPA EC4.0 pre-revised or Other EPA documents

Contaminant	RDX mg/kg/d	RDX mg/kg/d	CPSO kg/d/mg	CPSO kg/d/mg	VOC kg/d/mg	Artificial air kg/d/mg	Basic of RBC: certain agent effects			Residential mg/kg
							CPSO kg/d/mg	VOC kg/d/mg	Artificial air kg/d/mg	
Manganese and compounds	7439983 910107 24307264	5.00E-01 / 1.43E-05 / 3.00E-02 /	1.43E-05 / 9.00E-05 / 3.00E-02 /		1.00 / 1.3 / 1.00 /	0.032 / 0.33 / 1.10 /	0.032 / 0.33 / 1.10 /	0.032 / 0.33 / 1.10 /	0.032 / 0.33 / 1.10 /	300
Mepiquat chloride	7439976 22967926	3.00E-04 / 3.00E-04 /	1.47E-05 / 1.00E-04 /		1.1 / 1.1 /	0.31 / 0.11 /	0.31 / 0.11 /	0.41 / 0.41 /	0.41 / 0.41 /	7
Mercury (inorganic)	150505	3.00E-03 /			11 /	1.1 /	1.1 /	1.1 /	1.1 /	23
Mercury (methyl)					11 /	1.1 /	1.1 /	1.1 /	1.1 /	23
Merphos					11 /	1.1 /	1.1 /	1.1 /	1.1 /	23
Merphos oxide	73488	3.00E-05 /			1.1 /	0.11 /	0.041 /	0.041 /	0.041 /	23
Metalaxyl	57817191	6.00E-02 /			2200 /	220 /	21 /	21 /	21 /	23
Methacrylonitrile	930373	1.00E-03 /			3.7 /	0.73 /	0.14 /	0.14 /	0.14 /	23
Methanidophos					910 /	91 /	14 /	14 /	14 /	23
Methanol	673561	5.00E-01 /			1800 /	1800 /	680 /	680 /	680 /	3.9
Methidathion					37 /	3.7 /	1.4 /	1.4 /	1.4 /	3.9
Methionyl					180 /	18 /	6.8 /	6.8 /	6.8 /	3.9
Methoxychlor					73 /	7.3 /	2.7 /	2.7 /	2.7 /	3.9
2-Methoxyacetophenone					37 /	2.1 /	1.4 /	1.4 /	1.4 /	3.9
2-Methoxy-5-nitroaniline	105984	1.00E-03 /	3.71E-03 /	4.50E-02 /	1.1 /	0.14 /	0.089 /	0.089 /	0.089 /	3.9
Methyl acetate	99992	792009	1.00E-00 /		37000 /	3700 /	1400 /	1400 /	1400 /	3.9
Methyl acrylate	96333	3.00E-02 /			1100 /	110 /	41 /	41 /	41 /	3.9
2-Methylaniline hydrochloride	6363213		1.80E-01 /	1.80E-01 /	0.37 /	0.035 /	0.018 /	0.018 /	0.018 /	3.9
2-Methyleniline	955334		2.40E-01 /	2.40E-01 /	0.38 /	0.026 /	0.013 /	0.013 /	0.013 /	3.9
Methyl chlorocarboxate	75221	1.00E-00 /			37000 /	3700 /	1400 /	1400 /	1400 /	3.9
4-(2-Methyl-4-chlorophenoxy) butyric acid	948119	1.00E-02 /			370 /	37 /	14 /	14 /	14 /	3.9
2-Methyl-4-chlorobenzoic acid	94746	5.00E-04 /			18 /	1.8 /	0.68 /	0.68 /	0.68 /	3.9
2-(2-Methyl-4-chlorophenoxy)propanoic acid	936532	1.00E-03 /	8.57E-01 /		37 /	3.7 /	1.4 /	1.4 /	1.4 /	3.9
Methylene bromide	1018772	1.00E-02 /			31000 /	3100 /	1000 /	1000 /	1000 /	3.9
Methylene chloride	74933	6.00E-02 /	8.57E-01 /	1.64E-03 /	4.1 /	3.8 /	0.42 /	0.42 /	0.42 /	3.9
4,4'-Methylene bis(2-chloroaniline)	101144	7.00E-04 /	1.30E-01 /	1.30E-01 /	0.52 /	0.048 /	0.024 /	0.024 /	0.024 /	4.9
4,4'-Methylenebis(benzeneamine)	101779		2.50E-01 /		0.27 /	0.025 /	0.013 /	0.013 /	0.013 /	4.9
4,4'-Methylene bis(N,N'-dimethyl)aniline	101611		4.60E-02 /		1.3 /	0.14 /	0.069 /	0.069 /	0.069 /	2.6
4,4'-Methylenebisphenyl isocyanate	101638		5.71E-06 /		0.035 /	0.021 /	0.007 /	0.007 /	0.007 /	2.6
Methyl ethyl ketone	78933	6.00E-01 /	2.86E-01 /	1.10E+00	22000 /	1000 /	310 /	310 /	310 /	19
Methyl hydrazine	60344	8.00E-02 /			0.031 /	0.007 /	0.0029 /	0.0029 /	0.0029 /	19
Methyl isobutyl ketone	108101	8.00E-02 /			2500 /	94 /	24 /	24 /	24 /	20
Methyl methacrylate	80526				2900 /	290 /	110 /	110 /	110 /	20
2-Methyl-5-nitroaniline	99758				1800 /	180 /	64 /	64 /	64 /	20
Methyl perbenzoate	290000				1800 /	180 /	64 /	64 /	64 /	20
2-Methylphenol (o-cresol)	954467	2.50E-04 /			120 /	12 /	4.8 /	4.8 /	4.8 /	20
3-Methylphenol (m-cresol)	103394	5.00E-02 /			120 /	12 /	4.8 /	4.8 /	4.8 /	20
4-Methylphenol (p-cresol)	106443	5.00E-03 /			60 /	6 /	3.1 /	3.1 /	3.1 /	20
Methyl styrene (mixture)	75013154	6.00E-03 /	1.14E-01 /							20

Source: f-17153 h-HEIST s-1424ST oh- 117153-00000 from IRIS or HEIST -EP1-8210 internal c=Other EP1 documents

Ratio of EPC _{residential} / EPC _{industrial} & nonresidential effects to EPC _{other} EP4 documents										
Contaminant	CAS	RPD _d mg/kg/d	RD _i mg/kg/d	CPS _d kg/kg	CPS _i kg/kg	Ambient air C mg/m ³	O ₃ wet air C mg/m ³	Fish mg/l ₆	Industrial soil mg/kg	Residential soil mg/kg
Methyl styrene (alpha)	98-63-9	7.00E-02 *				430 n	250 n	95 n	72000 n	4300 n
Methyl tert-butyl ether (MTBE)	103-04-4	3.00E-01 *	1.37E-01 *			600 n	3100 n	64 n	9100 n	390 n
Metolaser (Dual)	51218452	1.50E-01 *				5500 n	150 n	200 n	150000 n	12000 n
Metributrin	218071649	2.50E-02 *				910 n	91 n	34 n	26000 n	2000 n
Mirex	23151533	2.00E-04 *				0.037 e	0.0035 e	0.0018 e	1.6 e	0.13 e
Molinate	2212671	2.00E-03 *				73 n	73 n	2.7 n	2000 n	160 n
Molybdenum	74359987	3.00E-03 *				180 n	18 n	6.4 n	9100 n	390 n
Monochloramine	103599903	1.00E-01 *				3700 n	370 n	140 n	100000 n	7800 n
Naled	300753	2.00E-03 *				73 n	73 n	2.7 n	2000 n	160 n
2-Naphthylamine	915238		1.30E-03 *			8.00E-02 *	0.000048 e	0.000024 e	0.023 e	0.0049 e
Napropamide	132959997	1.00E-01 *				3700 n	370 n	140 n	100000 n	7800 n
Nickel refinery dust			2.40E-01 *			0.0075 e	0.0075 e			
Nickel (soluble salts)	74400720	2.00E-02 *				730 n	73 n	27 n	20000 n	16000 n
Nickel sulfide	120315722	1.50E-03 *				1.70E+00 f	0.0037 e	0.0037 e		
Nitramprin	1929824	1.50E-03 *				53 n	53 n	2 n	70500 n	120 n
Nitrate	147971538	1.60E+00 *				5800 n	1800 n	2200 n	100000 n	130000 n
Nitric Oxide	10102439	1.00E-01 *				3700 n	370 n	140 n	100000 n	7800 n
Nitrite	14797650	1.00E-01 *				3700 n	370 n	140 n	100000 n	7800 n
2-Nitroaniline	48744	6.00E-03 *	3.71E-03 n			2.2 n	0.21 n	0.081 n	2.61 n	4.7 n
3-Nitroaniline	950092	3.00E-03 *				110 n	11 n	4.1 n	3100 n	230 n
4-Nitroaniline	100016	3.00E-03 *				110 n	11 n	4.1 n	3100 n	230 n
Nitrobenzene	989931	1.00E-04 *	3.71E-04 *			3.4 n	2.1 n	0.68 e	3.510 n	39 n
Nitrotetraenoin	67209	7.00E-02 *				1.90E+00 h	9.40E+00 n	260 n	93 n	72000 n
Nitrofurazone	391870					0.045 c	0.0067 e	0.0021 c		3500 n
Nitrogen dioxide	10102440	1.00E+00 *				37000 n	3700 n	1400 n	100000 n	7800 n
Nitroguanidine	316887	1.00E-01 *				3700 n	370 n	140 n	100000 n	7800 n
4-Nitrophenol	100027	6.20E-02 *				2300 n	210 n	34 n	63000 n	4800 n
2-Nitropropane	794469		5.71E-03 *			9.40E+00 h	9.40E+00 h	210 n	0.00057 e	
N-Nitroso-di-isobutyleramine	924163		5.40E+00 f			3.60E+00 f	0.012 c	0.0011 e	0.00038 e	0.12 e
N-Nitrosoethanolamine	1116547		2.80E+00 f			1.41E+02 f	0.024 c	0.00722 e		0.23 e
N,N-Diisopropylethylamine	151183					1.50E+02 f	0.00043 e	0.000041 e	0.000021 e	0.00043 e
N,N-Dimethylethylamine	62749					4.10E+01 f	0.00013 e	0.000015 e	0.0000032 e	0.013 e
N,N-Diisopropylethylamine	84306					4.90E+03 f	0.014 e	0.0013 e	0.00044 e	0.13 e
N-Nitroso di-n-propylethamine	621647		7.00E+00 f			0.0056 e	0.00043 e	0.000043 e	0.41 e	0.091 e
N,N-Diisopropylethylamine	10359936		2.20E+01 f			0.00011 e	0.00022 e	0.000014 e	0.13 e	0.029 e
N,N-Diisopropylamine	930552		2.10E+00 f			0.012 c	0.0029 e	0.0015 e	7.14 e	0.3 e
n-Nitrotoluene	99081	1.00E-02 *				61 n	37 n	14 n	100000 n	7800 n
c-Nitrotoluene	81722	1.00E-02 *				61 n	37 n	14 n	100000 n	7800 n
p-Nitrotoluene	99990	1.00E-02 *				61 n	37 n	14 n	10000 n	7800 n
Nufluazom	27314132	4.00E-02 *				1500 n	150 n	54 n	41000 n	3100 n
NuStar	11309199	7.00E-04 *				35 n	24 n	9.95 n	720 n	31 n
Octabromodiphenyl ether	32316329	3.00E-03 *				110 n	11 n	4.1 n	3100 n	230 n

Source: 1=IRIS, 2=RHEAST, 3=RHEAST air, 4=Information from IRIS or HEAST, 5=Other EPA documents

Contaminant	CAS	RfD	RfD _{air}	CPSI				CPSI				Basis of RBC: environmental effects			
				mg/kg/d	mg/kg/d	kg/d/mg	kg/d/mg	%	Top water	Ambient air	Fish	%	soil	Industrial soil	Residential soil
Octahydro-1,3,5,7-tetratinro-1,3,5,7-tetrazocine	2691410	5.00E-02	/					1000 n	100 n	68 n	31000 n	3900 n			
Oxytolin	192169	2.00E-03	n					73 n	73 n	27 n	2000 n	160 n			
Oxadiazole	19044873	3.00E-02	/					1000 n	100 n	68 n	51000 n	3900 n			
Oxetan	19666309	5.00E-03	/					100 n	18 n	6.8 n	9100 n	3900 n			
Oxyfluorfen	23135220	2.40E-02	/					910 n	91 n	34 n	26000 n	2000 n			
Paslobenzazoles	42874603	3.00E-03	/					10 n	11 n	4.1 n	3100 n	230 n			
Paraquat	1910423	4.50E-03	/					470 n	47 n	18 n	13000 n	1000 n			
Parathion	56382	6.00E-03	n					160 n	16 n	6.1 n	4600 n	350 n			
Pebulate	1114712	5.00E-02	n					220 n	22 n	8.1 n	6100 n	470 n			
Pendimethalin	40487421	4.00E-02	/					1000 n	100 n	68 n	31000 n	3900 n			
Pentabromo-6-chloro cyclohexane	87943				2.30E-02	/		1500 n	150 n	54 n	41000 n	3100 n			
Pentabromodiphenyl ether	32534819	2.00E-03	/					29 n	0.27 n	0.14 n	120 c	28 c			
Pentachloronitrobenzene	603931	8.00E-04	/					73 n	73 n	27 n	2000 n	160 n			
Pentachlorophenol	878653	3.00E-02	/		2.60E-01	/		4.9 n	29 n	1.1 n	620 n	63 n			
Pentethrin	32543531	5.00E-02	/		1.20E-01	/		0.041 n	0.024 n	0.012 n	11 n	2.9 n			
Phenmedipham	13684634	2.50E-01	/					0.36 n	0.032 n	0.026 n	24 n	53 n			
Phenol	108952	6.00E-01	/					1800 n	180 n	61 n	51000 n	3900 n			
p-Naphthylmethamine	108452	6.00E-03	/					9100 n	910 n	340 n	260000 n	20000 n			
p-Phenylenediamine	106503	1.90E-01	/					6900 n	690 n	260 n	61000 n	4700 n			
Phenylmercuric acetate	67284	8.00E-05	/			1.94E-03	/	2.9 n	0.29 n	0.11 n	102 n	6.3 n			
2-Phenylphenol	90437							33 n	32 n	1.6 n	1100 n	330 n			
Phorate	298022	2.00E-04	n					73 n	0.73 n	0.27 n	200 n	16 n			
Phosmet	732116	2.00E-02	/					730 n	73 n	27 n	20000 n	1600 n			
Phosphine	7803112	3.00E-04	/		8.57E-06	/		11 n	0.031 n	0.41 n	310 n	23 n			
Phosphorus (white)	7723140	2.00E-05	/					0.73 n	0.073 n	0.027 n	20 n	16 n			
p-Phthalic acid	100210	1.00E-00	/					3700 n	370 n	140 n	100000 n	78000 n			
Phthalic anhydride	65449	2.00E-00	/		3.43E-01	/		0.0076 n	0.0077 n	0.0033 n	0.32 n	0.072 n			
Pidolensin	1918021	7.00E-02	/					2600 n	260 n	95 n	72000 n	5100 n			
Pirimiphos-methyl	29252537	1.00E-02	/					370 n	37 n	14 n	10000 n	780 n			
Polybrominated biphenyls (PCBs)	131463	7.00E-06	n			8.90E+00	/	0.0076 n	0.0087 n	0.0001 n	0.00041 n	0.003 n	0.37 n	0.063 n	
Arcofar 1016	12674112	7.00E-05	/			7.70E+00	/	0.0076 n	0.0087 n	0.0001 n	0.00043 n	0.003 n	0.37 n	0.063 n	
Polychlorinated terphenyls (PCTs)															
Polyfurane aromatic hydrocarbons															
Acenaphthene	81329	6.00E-02	/					220 n	22 n	81 n	61000 n	4700 n			
Anthracene	120122	3.00E-01	/					1100 n	110 n	410 n	310000 n	23000 n			
Benzolalaphyrene	503728							0.0092 n	0.0091 n	0.00043 n	0.39 n	0.088 n			
Benz[b]fluoranthene	201992							0.0092 n	0.01 n	0.0043 n	0.39 n	0.088 n			
Benz[k]fluoranthene	207059							0.0092 n	0.01 n	0.0043 n	0.39 n	0.088 n			
Benz[e]anthracene	563533							0.0092 n	0.01 n	0.0043 n	0.39 n	0.088 n			

Source: 1=RCR 2=HEA/ST 3=Predictions from R/S or HEA/ST 4=Other EPA documents

Contaminant	CAS	RID ^a mg/kg/d	RID ^b mg/kg/d	CPS ^c kg/m ³	V ^d kg/m ³	Ambient air			Industrial soil		Residential soil	
						CPS mg/m ³	V kg/m ³	PPM ^e	ppm ^f /kg	ppm ^f /kg	ppm ^f /kg	
Chrysene	213619			7.30E-03 *	6.10E-03 *	9.2 *	1 *	0.43 *	390 *	390 *	83 *	
Dibenz[ah]anthracene	53703			7.30E-00 *	6.10E+00 *	1300 *	130 *	0.001 *	0.00043 *	0.39 *	0.028 *	
Fluorene	206440	4.00E-02 *										
Indenol 2,3-dipyrone	86737	4.00E-02 *										
Naphthalene	193394			7.30E-01 *	6.10E-01 *	1300 *	130 *	0.092 *	0.01 *	0.0043 *	0.39 *	
Pyrene	91203	4.00E-02 *										
Prochloraz	129000	3.00E-02 *										
Profluralin	67747093	9.00E-03 *		1.30E-01 *	0.49 *	1100 *	110 *	0.042 *	0.021 *	19 *	4.3 *	
Prometon	26399360	6.00E-03 *										
Prometryn	1610180	1.30E-02 *										
Pronamide	7227196	4.00E-03 *										
Propachlor	23950345	7.50E-02 *										
Propam	1911167	1.30E-02 *										
Propargite	7095938	3.00E-03 *										
Propargyl alcohol	2312333	2.00E-02 *										
Propazine	107197	2.00E-03 *										
Propiconazole	1394072	2.00E-02 *										
Propylene glycol	122425	2.00E-02 *										
Propylene glycol, monomethyl ether	60207991	1.30E-02 *										
Propylene glycol, monomethyl ether	575556	2.00E+01 *										
Propylene glycol, monomethyl ether	52125538	7.00E-01 *										
Propylene glycol, monomethyl ether	107982	7.00E-01 *	1.71E-01 *									
Propylene oxide	75569	8.37E-03 *		2.40E-01 *	1.29E-02 *			0.28 *	0.49 *	0.013 *	2.7 *	
Pursuit	81335773	2.50E-01 *						9100 *	910 *	340 *	20000 *	
Pyrdu	51630381	2.50E-02 *						910 *	91 *	18 *	13000 *	
Pyridine	110861	1.00E-03 *						37 *	3.7 *	27 *	20000 *	
Quinalphos	13930338	5.00E-04 *						18 *	1.8 *	0.68 *	39 *	
Quinoline	91224			1.20E-01 *				0.0036 *	0.00032 *	0.0026 *	0.013 *	
Resmethrin	10463868	3.00E-02 *						1100 *	110 *	41 *	31000 *	
Ronnel	2398413	3.00E-02 *						1800 *	180 *	68 *	51000 *	
Rotenone	63794	4.00E-03 *						159 *	15 *	5.4 *	310 *	
Savay	7987050	2.50E-02 *						910 *	91 *	34 *	26000 *	
Selenious Acid	7782492	5.00E-03 *						180 *	18 *	6.8 *	31000 *	
Selentium	7783008	5.00E-03 *						160 *	16 *	6.8 *	39000 *	
Selentium	630104	5.00E-03 *						160 *	16 *	6.8 *	39000 *	
Selonydium	74051502	9.00E-02 *						3300 *	330 *	120 *	920000 *	
Silver and compounds	7440724	3.00E-03 *		1.20E-01 *				160 *	18 *	6.8 *	31000 *	
Samarazine	122349	5.00E-03 *		1.20E-01 *				0.96 *	0.052 *	0.026 *	24 *	
Sodium azide	26620223	4.00E-03 *						159 *	15 *	5.4 *	31000 *	
Sodium diethylidithiocarbamate	140165	3.00E-02 *		2.70E-01 *				0.29 *	0.023 *	0.012 *	24 *	
Sodium Duroacetate	62748	2.00E-03 *						0.73 *	0.073 *	0.027 *	20 *	
Sodium metavanadate	13718268	1.00E-03 *						37 *	3.7 *	1.4 *	10000 *	

Sources: I=IRIS; h=HEAST; a=HEAST off-shore platforms from IRIS or HEAST; o=EPAC/ACID information; o=Other EPA documents.

Contaminant	Basis of RBC: c=centrifugate effects; n=nonsolvent effects										Residential soil mg/kg
	CAS	RD ₀ mg/kg/d	RD ₁ mg/kg/d	CPS ₀	CPS ₁	Y kg/d/eng	O TBP units	Ambient air µg/m ³	Fish mg/kg	Industrial air mg/kg	
Stannous, stable	74407246	6.00E-01						22000 n	810 n	610000 n	47000 n
Styrene	37219	3.00E-04	1					11 n	1.1 n	0.41 n	310 n
Systhane	100423	2.00E-01	1	2.86E-01	1			1600 n	1000 n	270 n	200000 n
2,3,7,8-TCDD (dioxin)	BB671890	2.50E-02	1			1.56E-05	h	1.16E-03	h	34 n	260000 n
Tetabutyltin	1746015							910 n	91 n	2.00E-03	o
Temephos	34014161	7.00E-02	1					2600 n	250 n	91 n	0.000018
Terbacil	33833968	2.00E-02	h					730 n	73 n	22000 n	1600 n
Terbefos	39025112	1.30E-02	1					470 n	47 n	18 n	13000 n
Terbutryn	13071729	2.50E-05	h					0.91 n	0.091 n	0.034 n	26 n
1,2,4,5-Tetrachlorobenzene	886500	1.00E-03	1					37 n	3.7 n	1.4 n	1000 n
1,1,1,2-Tetrachloroethane	959413	3.00E-04	1					1.8 n	1.1 n	0.41 n	310 n
1,1,2,2-Tetrachloroethane	630206	3.00E-02	1			2.50E-02	1	2.59E-02	1	0.24 n	0.12 n
Tetrachloroethylene (PCE)	793449					2.00E-01	1	2.03E-01	1	0.031 n	0.016 n
2,3,4,6-Tetrachlorophenol	127184	1.00E-02	1			5.20E-02	1	1.1 n	3.1 n	0.061	53 n
o,p,p,Tetrachlorotoluene	389872	3.00E-02	1			2.00E-01	1	1100 n	110 n	41 n	31000 n
Tetrachlormxyphene	5216231					0.00053	o	0.00031	o	0.00016	o
Tetracyldibutylorthophosphate	961114	3.00E-02	1			2.4 n	2.4 n	0.26 n	0.13 n	120 n	27 n
Thallium oxide	3689241	3.00E-04	1			2.40E-02	h	18 n	1.8 n	0.68 n	310 n
Thallium	1314324	7.00E-03	w					26 n	0.26 n	0.093 n	72 n
Thallium acetate	363618	9.00E-05	1					3.3 n	0.33 n	0.12 n	92 n
Thallium carbonate	63131739	8.00E-05	1					2.9 n	0.29 n	0.11 n	82 n
Thallium chloride	77911220	8.00E-03	1					2.9 n	0.29 n	0.11 n	82 n
Thallium nitrate	10102451	9.00E-05	1					3.3 n	0.33 n	0.12 n	92 n
Thallium selenite	12039520	9.00E-05	w					3.3 n	0.33 n	0.12 n	92 n
Thallium sulfate	7446186	8.00E-05	1					2.9 n	0.29 n	0.11 n	82 n
Thiocencarb	28249776	1.00E-02	1					370 n	37 n	14 n	10000 n
2-(Tricyanomethyl)benzisoxazole	21564470	3.00E-02	h					1100 n	110 n	41 n	31000 n
Thiophanox	39196184	3.00E-04	h					11 n	1.1 n	0.41 n	310 n
Thiophenate-methyl	23364038	8.00E-02	1					2500 n	250 n	110 n	22000 n
Thiram	1372618	5.00E-03	1					180 n	18 n	6.8 n	3100 n
Tin end compounds	6.00E-01	h						22000 n	2200 n	810 n	610000 n
Toluene	1098113	2.00E-01	1	1.14E-01	1			730 n	420 n	270 n	200000 n
Toluene-2,4-diamine	918677	6.00E-01	h	3.20E-00	h			22000 n	2200 n	810 n	610000 n
Toluene-2,5-diamine	957073	2.00E-01	h					7300 n	730 n	270 n	200000 n
Toluene-2,6-diamine	8723403							0.35 n	0.033 n	0.017 n	13 n
p-Toluidine	106490							0.061	0.0036	0.0029	13 n
Torpophene	8901512							270 n	27 n	10 n	7700 n
Tralomethrin	668441256	7.50E-03	1					470 n	47 n	18 n	13000 n
Triallate	2363175	1.30E-02	1					370 n	37 n	14 n	10000 n
Triasulfuron	82057503	1.00E-02	1					30 n	18 n	6.8 n	760 n
1,2,4-Tribromobenzene	615543	3.00E-03	1					30 n	18 n	6.8 n	350 n

Source: 1=IRIS & HEAT; 2=HEAT; 3=Information from IUS or IEST; 4=Other EPA documents

Basis of RBC: c=carcinogenic effects n=nocarcinogenic effects

Contaminant	CAS	RBC mg/kg/d	RBC mg/kg/d	CFS	CFS kg/d/m ³	V C mg/L	Ambient air µg/m ³	Residential soil mg/kg
Tributyltin oxide (TBT)	96359	3.00E-03 /		2.90E-02 h		1.1 n	0.041 n	31 n
2,4,6-Trichloroaniline hydrochloride	3166392			3.40E-02 h		2.3 n	0.22 n	99 n
2,4,6-Trichlorobenzene	6349333	1.00E-02 /	1.71E-02 h			0.2 n	0.11 n	22 n
1,2,4-Trichlorobutane	120821	9.00E-02 w	2.86E-01 w			190 n	210 n	14 n
1,1,1-Trichloroethane	71556			9.00E-03 /	1.60E-02 f	1300 n	1000 n	120 n
1,1,2-Trichloroethane	79003	4.00E-03 /		1.10E-02 f		0.19 n	0.11 n	0.93 n
Trichloroethylene (TCE)	79816	6.00E-03 *	1.00E-01 *	6.00E-03 *	1.60E-02 f	1.6 n	1.0	0.29 n
Trichlorofluoromethane	73694	3.00E-01 /	2.00E-01 *			1300 n	730 n	410 n
2,4,5-Trichlorophenol	91934	1.00E-01 /				3700 n	370 n	140 n
2,4,6-Trichlorophenol	88062			1.10E-02 f	1.09E-02 f	6.1 n	0.57 e	260 n
2,4,5-Trichlorophenoxyacetic acid	93765	1.00E-02 /				370 n	37 n	14 n
2-(2,4,5-Trichlorophenoxy)propionic acid	93721	8.00E-03 /				290 n	29 n	11 n
1,1,2-Trichloropropane	598776	4.00E-03 /				30 n	18 n	6.8 n
1,2,3-Trichloropropene	96184	6.00E-03 /		7.00E-00 /		0.0013 e	0.00089 e	0.00043 e
1,2,3-Trichloropropene	96195	5.00E-03 h				30 n	18 n	6.8 n
1,1,2-Trichloro-1,2,2-trifluoromethane	76131	3.00E-01 /	8.97E-00 h			59000 n	31000 n	41000 n
Tridiphatic Triethylamine	3011802	3.00E-03 /	2.00E-03 /			110 n	11 n	4.1 n
Trifluralin	1562094	7.50E-03 /	7.70E-03 f			6.7 n	0.51 n	0.41 n
1,2,4-Trimethylbenzene	95636	3.00E-04 *				3 n	1.8 n	0.63 n
1,2,5-Trimethylbenzene	108678	4.00E-04 *				2.4 n	1.5 n	0.54 n
Trimethyl phosphate	312561			3.70E-02 h		1.8 n	0.17 e	0.083 e
1,3,5-Trinitrobenzene	99334	5.00E-05 /				1.8 n	0.18 n	0.068 n
Trinitrophenylmethyltrimine	479458	1.00E-02 h				370 n	37 n	14 n
2,4,6-Trinitrotoluene	1189637	5.00E-04 /		3.00E-02 f		2.2 n	0.21 e	0.11 e
Uranium (soluble salts)	7440611		3.00E-03 /			110 n	11 n	4.1 n
Vanadium	7440622		7.00E-03 h			260 n	26 n	9.5 n
Vanadium pentoxide	1314621	9.00E-03 /				330 n	33 n	12 n
Vanadium sulfate	36907423	2.00E-02 h				730 n	73 n	27 n
Verbenam	1929777	1.00E-03 /				37 n	3.7 n	1.4 n
Vinclozolin	30471448	2.30E-02 /				910 n	91 n	34 n
Vinyl acetate	108054	1.00E+00 h	9.71E-02 f			3700 n	210 n	1400 n
Vinyl bromide	593602	1.71E-04 /	6.37E-04 /			5.2 n	3.1 n	
Vinyl chloride	73014	1.11E-04 /	1.11E-04 /	3.00E-01 h**	3.00E-01 h**	0.019 e	0.021 e	0.0017 e
Warfarin	81812	3.00E-03 h	3.00E-03 h			11 n	1.1 n	0.41 n
m-Xylene	103325	1.700E+00 h	1.700E+00 h	2.00E-01 *	2.00E-01 *	1.400 n	730 n	270 n
p-Xylene	954775	2.00E+00 h	1.700E+00 h			520 n	310 p	100000 n
Xylylene (mixed)	1084743	1.700E+00 h	1.700E+00 h			12000 n	7300 n	2700 n
Zinc	1330207	3.1E-02 /	3.1E-02 /			11000 n	1100 n	100000 n
Zinc phosphide	7440666	2.80E+04 /	2.80E+04 /			11 n	410 n	310 n
Zinc	1314447	3.00E-04 /				10000 n	68 n	31000 n
Zinc	12122677	9.00E+02 /				10000 n	100 n	3900 n

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