

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; H = HEAST; W = WHO; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; F = See FAQ #28; c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User's Guide); s = Concentration may exceed Csat (See User's Guide); SSL values are based on DAF=1

Contaminant	CAS No.	Toxicity and Chemical-specific Information											Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
		SFO	k _e	IUR	k _e	RfDo	k _e	RfCi	k _e	v	muta-	GIABS	ABS	PEF	VF	Csat	Ingestion	Dermal	Inhalation	Total	Ingestion	Dermal	Inhalation	Total
		(mg/kg-day) ⁻¹	(ug/m ³) ⁻¹	(mg/kg-day)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(m ³ /kg)	(m ³ /kg)	(mg/kg)	(m ³ /kg)	(m ³ /kg)	(m ³ /kg)	(m ³ /kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)
ALAR	1596-84-5	1.8E-02	C	5.1E-06	C	1.5E-01	I				1	0.1	1.4E+09				3.5E+01	1.1E+02	6.5E+05	2.7E+01	1.2E+04	4.2E+04	9.2E+03	
Acephate	30560-19-1	8.7E-03	I			4.0E-03	I				1	0.1	1.4E+09				7.3E+01	2.3E+02		5.6E+01	3.1E+02	1.1E+03	2.4E+02	
Acetaldehyde	75-07-0			2.2E-06	I			9.0E-03	I	V	1		1.4E+09	9.5E+03	1.1E+05				1.1E+01	1.1E+01		8.9E+01	8.9E+01	
Acetochlor	34256-82-1					2.0E-02	I				1	0.1	1.4E+09								1.6E+03	5.6E+03	1.2E+03	
Acetone	67-64-1					9.0E-01	I	3.1E+01	A	V	1		1.4E+09	1.4E+04	1.1E+05						7.0E+04	4.4E+05	6.1E+04	
Acetone Cyanohydrin	75-86-5					3.0E-03	P	6.0E-02	P	V	1		1.4E+09	2.6E+04	1.1E+05						2.3E+02	1.6E+03	2.0E+02	
Acetonitrile	75-05-8							6.0E-02	I	V	1		1.4E+09	1.4E+04	1.3E+05							8.7E+02	8.7E+02	
Acetophenone	98-86-2					1.0E-01	I			V	1		1.4E+09	6.2E+04	2.3E+03						7.8E+03		7.8E+03	
Acetylaminofluorene, 2-	53-96-3	3.8E+00	C	1.3E-03	C						1	0.1	1.4E+09				1.7E-01	5.3E-01	2.5E+03	1.3E-01				
Acrolein	107-02-8					5.0E-04	I	2.0E-05	I	V	1		1.4E+09	7.8E+03	2.5E+04						3.9E+01	1.6E-01	1.6E-01	
Acrylamide	79-06-1	4.5E+00	I	1.3E-03	I	2.0E-04	I				1	0.1	1.4E+09				1.4E-01	4.5E-01	2.5E+03	1.1E-01	1.6E+01	5.6E+01	1.2E+01	
Acrylic Acid	79-10-7					5.0E-01	I	1.0E-03	I		1	0.1	1.4E+09								3.9E+04	1.4E+05	1.4E+06	
Acrylonitrile	107-13-1	5.4E-01	I	6.8E-05	I	4.0E-02	A	2.0E-03	I	V	1		1.4E+09	8.2E+03	1.1E+04						3.1E+03	1.7E+01	1.7E+01	
Adiponitrile	111-69-3							6.0E-03	P		1	0.1	1.4E+09									8.5E+06	8.5E+06	
Alachlor	15972-60-8	5.8E-02	C			1.0E-02	I				1	0.1	1.4E+09				1.1E+01	3.6E+01		8.7E+00	7.8E+02	2.8E+03	6.1E+02	
Aldicarb	116-06-3					1.0E-03	I				1	0.1	1.4E+09								7.8E+01	2.8E+02	6.1E+01	
Aldicarb Sulfone	1646-88-4					1.0E-03	I				1	0.1	1.4E+09								7.8E+01	2.8E+02	6.1E+01	
Aldrin	309-00-2	1.7E+01	I	4.9E-03	I	3.0E-05	I				1	0.1	1.4E+09				3.8E-02	1.2E-01	6.8E+02	2.9E-02	2.3E+00	8.4E+00	1.8E+00	
Allyl	74223-64-6					2.5E-01	I				1	0.1	1.4E+09								2.0E+04	7.0E+04	1.5E+04	
Allyl Alcohol	107-18-6					5.0E-03	I	3.0E-04	P		1	0.1	1.4E+09								3.9E+02	1.4E+03	3.1E+02	
Allyl Chloride	107-05-1	2.1E-02	C	6.0E-06	C			1.0E-03	I	V	1		1.4E+09	1.8E+03	1.5E+03						7.0E-01	1.8E+00	1.8E+00	
Aluminum	7429-90-5					1.0E+00	P	5.0E-03	P		1		1.4E+09								7.8E+04	7.1E+06	7.7E+04	
Aluminum Phosphide	20859-73-8					4.0E-04	I				1		1.4E+09								3.1E+01		3.1E+01	
Amdro	67485-29-4					3.0E-04	I				1	0.1	1.4E+09								2.3E+01	8.4E+01	1.8E+01	
Ametryn	834-12-8					9.0E-03	I				1	0.1	1.4E+09								7.0E+02	2.5E+03	5.5E+02	
Aminobiphenyl, 4-	92-67-1	2.1E+01	C	6.0E-03	C						1	0.1	1.4E+09				3.0E-02	9.6E-02	5.5E+02	2.3E-02				
Aminophenol, m-	591-27-5					8.0E-02	P				1	0.1	1.4E+09								6.3E+03	2.2E+04	4.9E+03	
Aminophenol, p-	123-30-8					2.0E-02	P				1	0.1	1.4E+09								1.6E+03	5.6E+03	1.2E+03	
Amitraz	33089-61-1					2.5E-03	I				1	0.1	1.4E+09								2.0E+02	7.0E+02	1.5E+02	
Ammonium Perchlorate	7790-98-9					7.0E-04	I				1		1.4E+09								5.5E+01		5.5E+01	
Ammonium Sulfamate	7773-06-0					2.0E-01	I				1		1.4E+09								1.6E+04		1.6E+04	
Aniline	62-53-3	5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I		1	0.1	1.4E+09				1.1E+02	3.5E+02	2.1E+06	8.5E+01	5.5E+02	2.0E+03	1.4E+06	
Antimony (metallic)	7440-36-0					4.0E-04	I				0.15		1.4E+09								3.1E+01		3.1E+01	
Antimony Pentoxide	1314-60-9					5.0E-04	H				0.15		1.4E+09								3.9E+01		3.9E+01	
Antimony Potassium Tartrate	11071-15-1					9.0E-04	H				0.15		1.4E+09								7.0E+01		7.0E+01	
Antimony Tetroxide	1332-81-6					4.0E-04	H				0.15		1.4E+09								3.1E+01		3.1E+01	
Antimony Trioxide	1309-64-4							2.0E-04	I		0.15		1.4E+09									2.8E+05	2.8E+05	
Apollo	74115-24-5					1.3E-02	I				1	0.1	1.4E+09								1.0E+03	3.6E+03	7.9E+02	
Aramite	140-57-8	2.5E-02	I	7.1E-06	I	5.0E-02	H				1	0.1	1.4E+09				2.6E+01	8.1E+01	4.7E+05	1.9E+01	3.9E+03	1.4E+04	3.1E+03	
Arsenic, Inorganic	7440-38-2	1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C		1	0.03	1.4E+09				4.3E-01	4.5E+00	7.7E+02	3.9E-01	2.3E+01	2.8E+02	2.1E+04	
Arsine	7784-42-1							5.0E-05	I		1		1.4E+09									7.1E+04	7.1E+04	
Assure	76578-14-8					9.0E-03	I				1	0.1	1.4E+09								7.0E+02	2.5E+03	5.5E+02	
Asulam	3337-71-1					5.0E-02	I				1	0.1	1.4E+09								3.9E+03	1.4E+04	3.1E+03	
Atrazine	1912-24-9	2.3E-01	C			3.5E-02	I				1	0.1	1.4E+09				2.8E+00	8.8E+00		2.1E+00	2.7E+03	9.8E+03	2.1E+03	
Avermectin B1	65195-55-3					4.0E-04	I				1	0.1	1.4E+09								3.1E+01	1.1E+02	2.4E+01	
Azobenzene	103-33-3	1.1E-01	I	3.1E-05	I					V	1		1.4E+09	4.2E+05							3.3E+01	4.9E+00		
Barium	7440-39-3					2.0E-01	I	5.0E-04	H		0.07		1.4E+09								1.6E+04		1.5E+04	
Baygon	114-26-1					4.0E-03	I				1	0.1	1.4E+09								3.1E+02	1.1E+03	2.4E+02	
Bayleton	43121-43-3					3.0E-02	I				1	0.1	1.4E+09								2.3E+03	8.4E+03	1.8E+03	
Baythroid	68359-37-5					2.5E-02	I				1	0.1	1.4E+09								2.0E+03	7.0E+03	1.5E+03	
Benefin	1861-40-1					3.0E-01	I				1	0.1	1.4E+09								2.3E+04	8.4E+04	1.8E+04	
Benomyl	17804-35-2					5.0E-02	I				1	0.1	1.4E+09								3.9E+03	1.4E+04	3.1E+03	
Bentazon	25057-89-0					3.0E-02	I				1	0.1	1.4E+09								2.3E+03	8.4E+03	1.8E+03	
Benzaldehyde	100-52-7					1.0E-01	I			V	1		1.4E+09	3.2E+04	1.9E+03						7.8E+03		7.8E+03	
Benzene	71-43-2	5.5E-02	I	7.8E-06	I	4.0E-03	I	3.0E-02	I	V	1		1.4E+09	4.0E+03	2.0E+03		1.2E+01		1.3E+00	1.1E+00	3.1E+02		1.3E+02	
Benzenethiol	108-98-5					1.0E-05	H			V	1		1.4E+09	2.2E+04	1.4E+03						7.8E-01		7.8E-01	
Benzidine	92-87-5	2.3E+02	I	6.7E-02	I	3.0E-03	I			M	1	0.1	1.4E+09				6.5E-04	2.2E-03	1.9E+01	5.0E-04	2.3E+02	8.4E+02	1.8E+02	
Benzoic Acid	65-85-0					4.0E+00	I																	

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Contaminant	CAS No.	Toxicity and Chemical-specific Information										Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1								
		SFO	k _e	IUR	k _e	RfDo	k _e	RfCi	k _e	v _o	muta-	GIABS	ABS	PEF	VF	Csat	Ingestion	Dermal	Inhalation	Total	Ingestion	Dermal	Inhalation	Total
Analyte		(mg/kg-day) ⁻¹	(ug/m ³) ⁻¹	(mg/kg-day)	y	(mg/m ³)	y	(mg/m ³)	y	gen			(m ³ /kg)	(m ³ /kg)	(mg/kg)	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg
Bisphenol A	80-05-7			5.0E-02	I						1	0.1	1.4E+09							3.9E+03	1.4E+04			3.1E+03
Boron And Borates Only	7440-42-8			2.0E-01	I	2.0E-02	H				1		1.4E+09							1.6E+04		2.8E+07		1.6E+04
Boron Trifluoride	7637-07-2					7.0E-04	H				1		1.4E+09									9.9E+05		9.9E+05
Bromate	15541-45-4	7.0E-01	I	4.0E-03	I						1		1.4E+09			9.1E-01			9.1E-01	3.1E+02				3.1E+02
Bromobenzene	108-86-1			2.0E-02	P	1.0E-02	P	V			1		1.4E+09	9.6E+03	7.7E+02					1.6E+03		1.0E+02		9.4E+01
Bromodichloromethane	75-27-4	6.2E-02	I	3.7E-05	C	2.0E-02	I		V		1		1.4E+09	4.4E+03	9.9E+02	1.0E+01		2.9E-01	2.8E-01	1.6E+03				1.6E+03
Bromoform	75-25-2	7.9E-03	I	1.1E-06	I	2.0E-02	I				1	0.1	1.4E+09			8.1E+01	2.6E+02	3.0E+06	6.1E+01	1.6E+03	5.6E+03			1.2E+03
Bromomethane	74-83-9			1.4E-03	I	5.0E-03	I	V			1		1.4E+09	1.6E+03	3.6E+03					1.1E+02		8.5E+00		7.9E+00
Bromophos	2104-96-3			5.0E-03	H						1	0.1	1.4E+09							3.9E+02	1.4E+03			3.1E+02
Bromoxynil	1689-84-5			2.0E-02	I						1	0.1	1.4E+09							1.6E+03	5.6E+03			1.2E+03
Bromoxynil Octanoate	1689-99-2			2.0E-02	I						1	0.1	1.4E+09							1.6E+03	5.6E+03			1.2E+03
Butadiene, 1,3-	106-99-0	3.4E+00	C	3.0E-05	I			2.0E-03	I	V	1		1.4E+09	9.4E+02	6.9E+02	1.9E-01		7.7E-02	5.4E-02			2.0E+00		2.0E+00
Butanol, N-	71-36-3			1.0E-01	I						1	0.1	1.4E+09							7.8E+03	2.8E+04			6.1E+03
Butyl Benzyl Phthlate	85-68-7	1.9E-03	P	2.0E-01	I						1	0.1	1.4E+09			3.4E+02	1.1E+03		2.6E+02	1.6E+04	5.6E+04			1.2E+04
Butyl alcohol, sec-	78-92-2			2.0E+00	P	3.0E+01	P				1		1.4E+09							1.6E+05		4.3E+10		1.6E+05
Butylate	2008-41-5			5.0E-02	I						1	0.1	1.4E+09							3.9E+03	1.4E+04			3.1E+03
Butylphthalyl Butylglycolate	85-70-1			1.0E+00	I						1	0.1	1.4E+09							7.8E+04	2.8E+05			6.1E+04
Cacodylic Acid	75-60-5			2.0E-02	A						1	0.1	1.4E+09							1.6E+03	5.6E+03			1.2E+03
Cadmium (Diet)	7440-43-9			4.2E-03	C	1.0E-03	I	1.0E-05	A		0.025	0.001	1.4E+09					7.9E+02	7.9E+02	7.8E+01	7.0E+02	1.4E+04		7.0E+01
Caprolactam	105-60-2			5.0E-01	I						1	0.1	1.4E+09							3.9E+04	1.4E+05			3.1E+04
Captafol	2425-06-1	1.5E-01	C	4.3E-05	C	2.0E-03	I				1	0.1	1.4E+09			4.3E+00	1.3E+01	7.7E+04	3.2E+00	1.6E+02	5.6E+02			1.2E+02
Captan	133-06-2	2.3E-03	C	6.6E-07	C	1.3E-01	I				1	0.1	1.4E+09			2.8E+02	8.8E+02	5.0E+06	2.1E+02	1.0E+04	3.6E+04			7.9E+03
Carbaryl	63-25-2			1.0E-01	I						1	0.1	1.4E+09							7.8E+03	2.8E+04			6.1E+03
Carbofuran	1563-66-2			5.0E-03	I						1	0.1	1.4E+09							3.9E+02	1.4E+03			3.1E+02
Carbon Disulfide	75-15-0			1.0E-01	I	7.0E-01	I	V			1		1.4E+09	1.0E+03	2.6E+02					7.8E+03		7.3E+02		6.7E+02
Carbon Tetrachloride	56-23-5	1.3E-01	I	1.5E-05	I	7.0E-04	I	1.9E-01	A	V	1		1.4E+09	1.6E+03	4.8E+02	4.9E+00		2.7E-01	2.5E-01	5.5E+01		3.2E+02		4.7E+01
Carbosulfan	55285-14-8			1.0E-02	I						1	0.1	1.4E+09							7.8E+02	2.8E+03			6.1E+02
Carboxin	5234-68-4			1.0E-01	I						1	0.1	1.4E+09							7.8E+03	2.8E+04			6.1E+03
Chloral Hydrate	302-17-0			1.0E-01	I						1	0.1	1.4E+09							7.8E+03	2.8E+04			6.1E+03
Chloramben	133-90-4			1.5E-02	I						1	0.1	1.4E+09							1.2E+03	4.2E+03			9.2E+02
Chloranil	118-75-2	4.0E-01	H								1	0.1	1.4E+09			1.6E+00	5.0E+00		1.2E+00					
Chlordane	12789-03-6	3.5E-01	I	1.0E-04	I	5.0E-04	I	7.0E-04	I		0.04		1.4E+09			1.8E+00	1.4E+01	3.3E+04	1.6E+00	3.9E+01	3.5E+02	9.9E+05		3.5E+01
Chlordecone (Kepone)	143-50-0	1.6E+01	C	4.6E-03	C	5.0E-04	A				1	0.1	1.4E+09			4.0E-02	1.3E-01	7.2E+02	3.0E-02	3.9E+01	1.4E+02			3.1E+01
Chlorfenvinphos	470-90-6			7.0E-04	A						1	0.1	1.4E+09							5.5E+01	2.0E+02			4.3E+01
Chlorimuron, Ethyl-	90982-32-4			2.0E-02	I						1	0.1	1.4E+09							1.6E+03	5.6E+03			1.2E+03
Chlorine	7782-50-5			1.0E-01	I	1.5E-04	A				1		1.4E+09							7.8E+03		2.1E+05		7.5E+03
Chlorine Dioxide	10049-04-4			3.0E-02	I	2.0E-04	I				1		1.4E+09							2.3E+03		2.8E+05		2.3E+03
Chlorite (Sodium Salt)	7758-19-2			3.0E-02	I						1		1.4E+09							2.3E+03				2.3E+03
Chloro-1,1-difluoroethane, 1-	75-68-3					5.0E+01	I	V			1		1.4E+09	1.1E+03	1.2E+03							5.9E+04		5.9E+04
Chloro-1,3-butadiene, 2-	126-99-8			2.0E-02	H	7.0E-03	H	V			1		1.4E+09	1.2E+03	8.2E+02					1.6E+03		8.7E+00		8.6E+00
Chloro-2-methylaniline HCl, 4-	3165-93-3	4.6E-01	H								1	0.1	1.4E+09			1.4E+00	4.4E+00		1.1E+00					
Chloro-2-methylaniline, 4-	95-69-2	2.7E-01	C	7.7E-05	C						1	0.1	1.4E+09			2.4E+00	7.5E+00	4.3E+04	1.8E+00					
Chloroacetic Acid	79-11-8			2.0E-03	H						1	0.1	1.4E+09							1.6E+02	5.6E+02			1.2E+02
Chloroacetophenone, 2-	532-27-4					3.0E-05	I				1	0.1	1.4E+09									4.3E+04		4.3E+04
Chloroaniline, p-	106-47-8	2.0E-01	P	4.0E-03	I						1	0.1	1.4E+09			3.2E+00	1.0E+01		2.4E+00	3.1E+02	1.1E+03			2.4E+02
Chlorobenzene	108-90-7			2.0E-02	I	5.0E-02	P	V			1		1.4E+09	7.4E+03	8.6E+02					1.6E+03		3.9E+02		3.1E+02
Chlorobenzilate	510-15-6	1.1E-01	C	3.1E-05	C	2.0E-02	I				1	0.1	1.4E+09			5.8E+00	1.8E+01	1.1E+05	4.4E+00	1.6E+03	5.6E+03			1.2E+03
Chlorobenzotrifluoride, 4-	98-56-6			3.0E-03	P	3.0E-01	P	V			1		1.4E+09	7.9E+03	5.5E+02					2.3E+02		2.5E+03		2.1E+02
Chlorobutane, 1-	109-69-3			4.0E-02	P						1		1.4E+09	2.0E+03	7.9E+02					3.1E+03				3.1E+03
Chlorodifluoromethane	75-45-6					5.0E+01	I	V			1		1.4E+09	1.0E+03	1.7E+03							5.3E+04		5.3E+04
Chloroform	67-66-3	3.1E-02	C	2.3E-05	I	1.0E-02	I	9.8E-02	A	V	1		1.4E+09	2.9E+03	2.7E+03	2.1E+01		3.1E-01	3.0E-01	7.8E+02			3.0E+02	2.2E+02
Chloromethane	74-87-3					9.0E-02	I	V			1		1.4E+09	1.3E+03	1.4E+03							1.2E+02		1.2E+02
Chloromethyl Methyl Ether	107-30-2	2.4E+00	C	6.9E-04	C						1		1.4E+09	5.3E+03	8.1E+03	2.7E-01		1.9E-02	1.8E-02					
Chloronaphthalene, Beta-	91-58-7			8.0E-02	I						1		1.4E+09	9.4E+04	2.1E+02					6.3E+03				6.3E+03
Chloronitrobenzene, o-	88-73-3	9.7E-03	P	1.0E-03	P	7.0E-05	P				1	0.1	1.4E+09			6.6E+01								

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; H = HEAST; W = WHO; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; F = See FAQ #28; c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User's Guide); s = Concentration may exceed Csat (See User's Guide); SSL values are based on DAF=1

Contaminant	Toxicity and Chemical-specific Information														Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1					
	CAS No.	SFO (mg/kg-day) ⁻¹	k e (ug/m ³) ⁻¹	IUR (mg/kg-day)	k e (mg/kg-day)	RfDo (mg/m ³)	k e (mg/m ³)	RfCi (mg/m ³)	k e (mg/m ³)	v o l u t a b i l i t y	m u t a g e n	GIABS	ABS	PEF (m ³ /kg)	VF (m ³ /kg)	Csat (mg/kg)	Ingestion mg/kg	Dermal mg/kg	Inhalation mg/kg	Total mg/kg	Ingestion mg/kg	Dermal mg/kg	Inhalation mg/kg	Total mg/kg
AnalYTE																								
Cresol, m-	108-39-4					5.0E-02	I					1	0.1	1.4E+09							3.9E+03	1.4E+04		3.1E+03
Cresol, o-	95-48-7					5.0E-02	I					1	0.1	1.4E+09							3.9E+03	1.4E+04		3.1E+03
Cresol, p-	106-44-5					5.0E-03	H					1	0.1	1.4E+09							3.9E+02	1.4E+03		3.1E+02
Cresols	1319-77-3					1.0E-01	A	6.0E-01	C	V		1		1.4E+09	3.9E+05	7.1E+04					7.8E+03		2.4E+05	7.6E+03
Crotonaldehyde, trans-	123-73-9	1.9E+00	H							V		1		1.4E+09	2.2E+04	2.4E+04	3.4E-01			3.4E-01				
Cumene	98-82-8					1.0E-01	I	4.0E-01	I	V		1		1.4E+09	7.2E+03	3.1E+02					7.8E+03		3.0E+03	2.2E+03
Cyanazine	21725-46-2	8.4E-01	H			2.0E-03	H					1	0.1	1.4E+09			7.6E-01	2.4E+00		5.8E-01	1.6E+02	5.6E+02		1.2E+02
Cyanides																								
-Calcium Cyanide	592-01-8					4.0E-02	I					1		1.4E+09							3.1E+03			3.1E+03
-Copper Cyanide	544-92-3					5.0E-03	I					1		1.4E+09							3.9E+02			3.9E+02
-Cyanide (CN-)	57-12-5					2.0E-02	I					1		1.4E+09							1.6E+03			1.6E+03
-Cyanogen	460-19-5					4.0E-02	I			V		1		1.4E+09							3.1E+03			3.1E+03
-Cyanogen Bromide	506-68-3					9.0E-02	I			V		1		1.4E+09							7.0E+03			7.0E+03
-Cyanogen Chloride	506-77-4					5.0E-02	I			V		1		1.4E+09							3.9E+03			3.9E+03
-Hydrogen Cyanide	74-90-8					2.0E-02	I	3.0E-03	I	V		1		1.4E+09							1.6E+03	4.3E+06		1.6E+03
-Potassium Cyanide	151-50-8					5.0E-02	I					1		1.4E+09							3.9E+03			3.9E+03
-Potassium Silver Cyanide	506-61-6					2.0E-01	I				0.04	1		1.4E+09							1.6E+04			1.6E+04
-Silver Cyanide	506-64-9					1.0E-01	I				0.04	1		1.4E+09							7.8E+03			7.8E+03
-Sodium Cyanide	143-33-9					4.0E-02	I					1		1.4E+09							3.1E+03			3.1E+03
-Thiocyanate	463-56-9					2.0E-04	P			V		1		1.4E+09	7.0E+03	5.6E+03				1.6E+01				1.6E+01
-Zinc Cyanide	557-21-1					5.0E-02	I					1		1.4E+09							3.9E+03			3.9E+03
Cyclohexane	110-82-7							6.0E+00	I	V		1		1.4E+09	1.2E+03	1.2E+02						7.2E+03		7.2E+03
Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	2.3E-02	H									1	0.1	1.4E+09			2.8E+01	8.8E+01		2.1E+01				
Cyclohexanone	108-94-1					5.0E+00	I					1	0.1	1.4E+09							3.9E+05	1.4E+06		3.1E+05
Cyclohexylamine	108-91-8					2.0E-01	I					1	0.1	1.4E+09							1.6E+04	5.6E+04		1.2E+04
Cyhalothrin/karate	68085-85-8					5.0E-03	I					1	0.1	1.4E+09							3.9E+02	1.4E+03		3.1E+02
Cypermethrin	52315-07-8					1.0E-02	I					1	0.1	1.4E+09							7.8E+02	2.8E+03		6.1E+02
Cyromazine	66215-27-8					7.5E-03	I					1	0.1	1.4E+09							5.9E+02	2.1E+03		4.6E+02
DDD	72-54-8	2.4E-01	I	6.9E-05	C							1	0.1	1.4E+09			2.7E+00	8.4E+00	4.8E+04	2.0E+00				
DDE, p,p'-	72-55-9	3.4E-01	I	9.7E-05	C							1	0.1	1.4E+09			1.9E+00	6.0E+00	3.4E+04	1.4E+00				
DDT	50-29-3	3.4E-01	I	9.7E-05	I	5.0E-04	I					1	0.03	1.4E+09			1.9E+00	2.0E+01	3.4E+04	1.7E+00	3.9E+01	4.7E+02		3.6E+01
Dacthal	1861-32-1					1.0E-02	I					1	0.1	1.4E+09							7.8E+02	2.8E+03		6.1E+02
Dalapon	75-99-0					3.0E-02	I					1	0.1	1.4E+09							2.3E+03	8.4E+03		1.8E+03
Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'-(BDE-209)	1163-19-5	7.0E-04	I			7.0E-03	I					1	0.1	1.4E+09			9.1E+02	2.9E+03		6.9E+02	5.5E+02	2.0E+03		4.3E+02
Demeton	8065-48-3					4.0E-05	I					1	0.1	1.4E+09							3.1E+00	1.1E+01		2.4E+00
Di(2-ethylhexyl)adipate	103-23-1	1.2E-03	I			6.0E-01	I					1	0.1	1.4E+09			5.3E+02	1.7E+03		4.0E+02	4.7E+04	1.7E+05		3.7E+04
Diallate	2303-16-4	6.1E-02	H									1	0.1	1.4E+09			1.0E+01	3.3E+01		8.0E+00				
Diazinon	333-41-5					7.0E-04	A					1	0.1	1.4E+09							5.5E+01	2.0E+02		4.3E+01
Dibromo-3-chloropropane, 1,2-	96-12-8	8.0E-01	P	6.0E-03	P	2.0E-04	P	2.0E-04	I	V	M	1		1.4E+09	3.6E+04	1.1E+03	1.9E-01		5.8E-03	5.6E-03	1.6E+01		7.6E+00	5.1E+00
Dibromobenzene, 1,4-	106-37-6					1.0E-02	I					1	0.1	1.4E+09							7.8E+02	2.8E+03		6.1E+02
Dibromochloromethane	124-48-1	8.4E-02	I	2.7E-05	C	2.0E-02	I			V		1	0.1	1.4E+09	8.8E+03	8.5E+02	7.6E+00	2.4E+01	8.0E-01	7.0E-01	1.6E+03	5.6E+03		1.2E+03
Dibromoethane, 1,2-	106-93-4	2.0E+00	I	6.0E-04	I	9.0E-03	I	9.0E-03	I	V		1		1.4E+09	9.5E+03	1.4E+03	3.2E-01		3.9E-02	3.4E-02	7.0E+02		8.9E+01	7.9E+01
Dibromomethane (Methylene Bromide)	74-95-3					1.0E-02	H			V		1		1.4E+09	6.2E+03	3.0E+03					7.8E+02			7.8E+02
Dibutyl Phthalate	84-74-2					1.0E-01	I					1	0.1	1.4E+09							7.8E+03	2.8E+04		6.1E+03
Dibutyltin Compounds	NA					3.0E-04	P					1	0.1	1.4E+09							2.3E+01	8.4E+01		1.8E+01
Dicamba	1918-00-9					3.0E-02	I					1	0.1	1.4E+09							2.3E+03	8.4E+03		1.8E+03
Dichloro-2-butene, 1,4-	764-41-0			4.2E-03	P					V		1		1.4E+09	3.4E+03	6.1E+02				2.0E-03	2.0E-03			
Dichloro-2-butene, cis-1,4-	1476-11-5			4.2E-03	P					V		1	0.1	1.4E+09	3.6E+03	6.1E+02				2.1E-03	2.1E-03			
Dichloro-2-butene, trans-1,4-	110-57-6			4.2E-03	P					V		1	0.1	1.4E+09	1.3E+04	5.8E+02				7.3E-03	7.3E-03			
Dichloroacetic Acid	79-43-6	5.0E-02	I			4.0E-03	I					1	0.1	1.4E+09			1.3E+01	4.0E+01		9.7E+00	3.1E+02	1.1E+03		2.4E+02
Dichlorobenzene, 1,2-	95-50-1					9.0E-02	I	2.0E-01	H	V		1		1.4E+09	1.3E+04	2.2E+02					7.0E+03		2.8E+03	2.0E+03
Dichlorobenzene, 1,4-	106-46-7	5.4E-03	C	1.1E-05	C	7.0E-02	A	8.0E-01	I	V		1		1.4E+09	1.2E+04		1.2E+02		2.7E+00	2.6E+00	5.5E+03		1.0E+04	3.5E+03
Dichlorobenzidine, 3,3'-	91-94-1	4.5E-01	I	3.4E-04	C							1	0.1	1.4E+09			1.4E+00	4.5E+00	9.7E+03	1.1E+00				
Dichlorodifluoromethane	75-71-8					2.0E-01	I	2.0E-01	H	V		1		1.4E+09	9.0E+02	8.5E+02					1.6E+04		1.9E+02	1.9E+02
Dichloroethane, 1,1-	75-34-3	5.7E-03	C	1.6E-06	C	2.0E-01	P			V		1		1.4E+09	2.3E+03	1.8E+03	1.1E+02		3.5E+00	3.4E+00	1.6E+04			1.6E+04
Dichloroethane, 1,2-	107-06-2	9.1E-02	I	2.6E-05																				

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; H = HEAST; W = WHO; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; F = See FAQ #28; c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User's Guide); s = Concentration may exceed Csat (See User's Guide); SSL values are based on DAF=1

Contaminant	CAS No.	Toxicity and Chemical-specific Information											Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
		SFO	k _e	IUR	k _e	RfDo	k _e	RfCi	k _e	v _o	muta-	GIABS	ABS	PEF	VF	Csat	Ingestion	Dermal	Inhalation	Total	Ingestion	Dermal	Inhalation	Total
Analyte		(mg/kg-day) ⁻¹	(ug/m ³) ⁻¹	(mg/kg-day)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(m ³ /kg)	(m ³ /kg)	(mg/kg)	(m ³ /kg)	(m ³ /kg)	(m ³ /kg)	(m ³ /kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)
Diethylene Glycol Monobutyl Ether	112-34-5			1.0E-02	P	2.0E-02	P	1	0.1	1.4E+09										7.8E+02	2.8E+03	2.8E+07	6.1E+02	
Diethylene Glycol Monoethyl Ether	111-90-0			6.0E-02	P	3.0E-03	P	1	0.1	1.4E+09										4.7E+03	1.7E+04	4.3E+06	3.7E+03	
Diethylformamide	617-84-5			1.0E-03	P			1	0.1	1.4E+09										7.8E+01	2.8E+02		6.1E+01	
Diethylstilbestrol	56-53-1	3.5E+02	C	1.0E-01	C			1	0.1	1.4E+09						1.8E-03	5.8E-03	3.3E+01	1.4E-03					
Difenzozat	43222-48-6			8.0E-02	I			1	0.1	1.4E+09										6.3E+03	2.2E+04		4.9E+03	
Diflubenzuron	35367-38-5			2.0E-02	I			1	0.1	1.4E+09										1.6E+03	5.6E+03		1.2E+03	
Difluoroethane, 1,1-	75-37-6					4.0E+01	I V	1		1.4E+09	1.3E+03	1.5E+03										5.3E+04	5.3E+04	
Diisopropyl Ether	108-20-3					4.0E-01	P V	1		1.4E+09	2.9E+03	1.6E+03										1.2E+03	1.2E+03	
Diisopropyl Methylphosphonate	1445-75-6			8.0E-02	I			1		1.4E+09	2.8E+04	4.3E+02								6.3E+03			6.3E+03	
Dimethipin	55290-64-7			2.0E-02	I			1	0.1	1.4E+09										1.6E+03	5.6E+03		1.2E+03	
Dimethoate	60-51-5			2.0E-04	I			1	0.1	1.4E+09										1.6E+01	5.6E+01		1.2E+01	
Dimethoxybenzidine, 3,3'	119-90-4	1.4E-02	H					1	0.1	1.4E+09						4.6E+01	1.4E+02		3.5E+01					
Dimethyl methylphosphonate	756-79-6	1.7E-03	P	6.0E-02	P			1	0.1	1.4E+09						3.8E+02	1.2E+03		2.9E+02	4.7E+03	1.7E+04		3.7E+03	
Dimethylamino azobenzene [p-]	60-11-7	4.6E+00	C	1.3E-03	C			1	0.1	1.4E+09						1.4E-01	4.4E-01	2.5E+03	1.1E-01					
Dimethylaniline HCl, 2,4-	21436-96-4	5.8E-01	H					1	0.1	1.4E+09						1.1E+00	3.5E+00		8.4E-01					
Dimethylaniline, 2,4-	95-68-1	7.5E-01	H					1	0.1	1.4E+09						8.5E-01	2.7E+00		6.5E-01					
Dimethylaniline, N,N-	121-69-7			2.0E-03	I			1		1.4E+09	3.3E+04	8.2E+02								1.6E+02			1.6E+02	
Dimethylbenzidine, 3,3'	119-93-7	1.1E+01	P					1	0.1	1.4E+09						5.8E-02	1.8E-01		4.4E-02					
Dimethylformamide	68-12-2			1.0E-01	P	3.0E-02	I	1	0.1	1.4E+09						1.2E-03	3.7E-03	2.1E+01	8.8E-04	7.8E+03	2.8E+04	4.3E+07	6.1E+03	
Dimethylhydrazine, 1,2-	540-73-8	5.5E+02	C	1.6E-01	C			1	0.1	1.4E+09										1.6E+03	5.6E+03		1.2E+03	
Dimethylphenol, 2,4-	105-67-9			2.0E-02	I			1	0.1	1.4E+09										4.7E+01	1.7E+02		3.7E+01	
Dimethylphenol, 2,6-	576-26-1			6.0E-04	I			1	0.1	1.4E+09										7.8E+01	2.8E+02		6.1E+01	
Dimethylphenol, 3,4-	95-65-8			1.0E-03	I			1	0.1	1.4E+09										7.8E+03			7.8E+03	
Dimethylterephthalate	120-61-6			1.0E-01	I			1		1.4E+09	2.4E+04	6.1E+00								7.8E+00	2.8E+01		6.1E+00	
Dinitro-o-cresol, 4,6-	534-52-1			1.0E-04	P			1	0.1	1.4E+09										1.6E+02	5.6E+02		1.2E+02	
Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5			2.0E-03	I			1	0.1	1.4E+09										7.8E+00	2.8E+01		6.1E+00	
Dinitrobenzene, 1,2-	528-29-0			1.0E-04	P			1	0.1	1.4E+09										7.8E+00	2.8E+01		6.1E+00	
Dinitrobenzene, 1,3-	99-65-0			1.0E-04	I			1	0.1	1.4E+09										7.8E+00	2.8E+01		6.1E+00	
Dinitrobenzene, 1,4-	100-25-4			1.0E-04	P			1	0.1	1.4E+09										7.8E+00	2.8E+01		6.1E+00	
Dinitrophenol, 2,4-	51-28-5			2.0E-03	I			1	0.1	1.4E+09										1.6E+02	5.6E+02		1.2E+02	
Dinitrotoluene Mixture, 2,4/2,6-	25321-14-6	6.8E-01	I					1	0.1	1.4E+09						9.4E-01	3.0E+00		7.1E-01					
Dinitrotoluene, 2,4-	121-14-2	3.1E-01	C	8.9E-05	C	2.0E-03	I	1	0.102	1.4E+09						2.1E+00	6.4E+00	3.7E+04	1.6E+00	1.6E+02	5.5E+02		1.2E+02	
Dinitrotoluene, 2,6-	606-20-2			1.0E-03	P			1	0.099	1.4E+09										7.8E+01	2.8E+02		6.1E+01	
Dinitrotoluene, 2-Amino-4,6-	35572-78-2			2.0E-03	I			1	0.006	1.4E+09										1.6E+02	9.3E+03		1.5E+02	
Dinitrotoluene, 4-Amino-2,6-	19406-51-0			2.0E-03	I			1	0.009	1.4E+09										1.6E+02	6.2E+03		1.5E+02	
Dinoseb	88-85-7			1.0E-03	I			1	0.1	1.4E+09										7.8E+01	2.8E+02		6.1E+01	
Dioxane, 1,4-	123-91-1	1.1E-02	I	7.7E-06	C	1.0E-01	A	3.6E+00	A	1	0.1	1.4E+09				5.8E+01	1.8E+02	4.3E+05	4.4E+01	7.8E+03	2.8E+04	5.1E+09	6.1E+03	
Dioxins																								
-Hexachlorodibenzo-p-dioxin, Mixture	NA	6.2E+03	I	1.3E+00	I			1	0.03	1.4E+09						1.0E-04	1.1E-03	2.5E+00	9.4E-05					
-TCDD, 2,3,7,8-	1746-01-6	1.3E+05	C	3.8E+01	C	1.0E-09	A	4.0E-08	C	1	0.03	1.4E+09				4.9E-06	5.2E-05	8.7E-02	4.5E-06	7.8E-05	9.3E-04	5.7E+01	7.2E-05	
Diphenamid	957-51-7			3.0E-02	I			1	0.1	1.4E+09										2.3E+03	8.4E+03		1.8E+03	
Diphenyl Sulfone	127-63-9			3.0E-03	P			1	0.1	1.4E+09										2.3E+02	8.4E+02		1.8E+02	
Diphenylamine	122-39-4			2.5E-02	I			1	0.1	1.4E+09										2.0E+03	7.0E+03		1.5E+03	
Diphenylhydrazine, 1,2-	122-66-7	8.0E-01	I	2.2E-04	I			1	0.1	1.4E+09						8.0E-01	2.5E+00	1.5E+04	6.1E-01					
Diquat	85-00-7			2.2E-03	I			1	0.1	1.4E+09										1.7E+02	6.1E+02		1.3E+02	
Direct Black 38	1937-37-7	7.4E+00	C	2.1E-03	C			1	0.1	1.4E+09						8.6E-02	2.7E-01	1.6E+03	6.6E-02					
Direct Blue 6	2602-46-2	7.4E+00	C	2.1E-03	C			1	0.1	1.4E+09						8.6E-02	2.7E-01	1.6E+03	6.6E-02					
Direct Brown 95	16071-86-6	6.7E+00	C	1.9E-03	C			1	0.1	1.4E+09						9.5E-02	3.0E-01	1.7E+03	7.2E-02					
Disulfoton	298-04-4			4.0E-05	I			1	0.1	1.4E+09										3.1E+00	1.1E+01		2.4E+00	
Dithiane, 1,4-	505-29-3			1.0E-02	I			1	0.1	1.4E+09										7.8E+02	2.8E+03		6.1E+02	
Diuron	330-54-1			2.0E-03	I			1	0.1	1.4E+09										1.6E+02	5.6E+02		1.2E+02	
Dodine	2439-10-3			4.0E-03	I			1	0.1	1.4E+09										3.1E+02	1.1E+03		2.4E+02	
EPTC	759-94-4			2.5E-02	I			1		1.4E+09	1.6E+05	6.2E+02								2.0E+03			2.0E+03	
Endosulfan	115-29-7			6.0E-03	I			1	0.1	1.4E+09										4.7E+02	1.7E+03		3.7E+02	
Endothall	145-73-3			2.0E-02	I			1	0.1	1.4E+09										1.6E+03	5.6E+03		1.2E+03	
Endrin	72-20-8			3.0E-04	I			1	0.1	1.4E+09										2.3E+01	8.4E+01		1.8E+01	
Epichlorohydrin	106-89-8	9.9E-03	I	1.2E-06	I	6.0E-03	P	1.0E-03	I V	1	1.4E+09	1.8E+04	8.4E+03			6.5E+01		3.6E+01	2.3E+01	4.7E+02		1.9E+01	1.8E+01	
Epoxybutane, 1,2-	106-88-7					2.0E-02	I V	1		1.4E+09	7.4E+03													

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; H = HEAST; W = WHO; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; F = See FAQ #28; c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User's Guide); s = Concentration may exceed Csat (See User's Guide); SSL values are based on DAF=1

Contaminant	CAS No.	Toxicity and Chemical-specific Information										Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1								
		SFO	ke	IUR	ke	RfDo	ke	RfCi	ke	Vol	muta-	GIABS	ABS	PEF	VF	Csat	Ingestion	Dermal	Inhalation	Total	Ingestion	Dermal	Inhalation	Total
Analyte		(mg/kg-day) ⁻¹	(ug/m ³) ⁻¹	(mg/kg-day)	y	(mg/m ³)	y	(mg/m ³)	y	gen			(m ³ /kg)	(m ³ /kg)	(mg/kg)	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg
Ethylene Glycol	107-21-1			2.0E+00	I	4.0E-01	C				1	0.1	1.4E+09								1.6E+05	5.6E+05	5.7E+08	1.2E+05
Ethylene Glycol Monobutyl Ether	111-76-2			5.0E-01	I	1.3E+01	I				1	0.1	1.4E+09								3.9E+04	1.4E+05	1.8E+10	3.1E+04
Ethylene Oxide	75-21-8	3.1E-01	C	8.8E-05	C			3.0E-02	C	V	1		1.4E+09	6.3E+03	1.1E+05	2.1E+00		1.7E-01	1.6E-01				2.0E+02	2.0E+02
Ethylene Thiourea	96-45-7	4.5E-02	C	1.3E-05	C	8.0E-05	I				1	0.1	1.4E+09			1.4E+01	4.5E+01	2.5E+05	1.1E+01		6.3E+00	2.2E+01		4.9E+00
Ethylphthalyl Ethyl Glycolate	84-72-0			3.0E+00	I						1	0.1	1.4E+09								2.3E+05	8.4E+05		1.8E+05
Express	101200-48-0			8.0E-03	I						1	0.1	1.4E+09								6.3E+02	2.2E+03		4.9E+02
Fenamiphos	22224-92-6			2.5E-04	I						1	0.1	1.4E+09								2.0E+01	7.0E+01		1.5E+01
Fenpropathrin	39515-41-8			2.5E-02	I						1	0.1	1.4E+09								2.0E+03	7.0E+03		1.5E+03
Fluometuron	2164-17-2			1.3E-02	I						1	0.1	1.4E+09								1.0E+03	3.6E+03		7.9E+02
Fluorine (Soluble Fluoride)	7782-41-4			6.0E-02	I						1		1.4E+09								4.7E+03			4.7E+03
Fluridone	59756-60-4			8.0E-02	I						1	0.1	1.4E+09								6.3E+03	2.2E+04		4.9E+03
Flurprimidol	56425-91-3			2.0E-02	I						1	0.1	1.4E+09								1.6E+03	5.6E+03		1.2E+03
Flutolanil	66332-96-5			6.0E-02	I						1	0.1	1.4E+09								4.7E+03	1.7E+04		3.7E+03
Fluvalinate	69409-94-5			1.0E-02	I						1	0.1	1.4E+09								7.8E+02	2.8E+03		6.1E+02
Folpet	133-07-3	3.5E-03	I			1.0E-01	I				1	0.1	1.4E+09			1.8E+02	5.8E+02			1.4E+02	7.8E+03	2.8E+04		6.1E+03
Fomesafen	72178-02-0	1.9E-01	I								1	0.1	1.4E+09			3.4E+00	1.1E+01			1.4E+00	7.8E+03	2.8E+04		6.1E+03
Fonofos	944-22-9			2.0E-03	I						1	0.1	1.4E+09								1.6E+02	5.6E+02		1.2E+02
Formaldehyde	50-00-0			1.3E-05	I	2.0E-01	I	9.8E-03	A		1	0.1	1.4E+09					2.5E+05	2.5E+05		1.6E+04	5.6E+04	1.4E+07	1.2E+04
Formic Acid	64-18-6			2.0E+00	H	3.0E-03	P				1	0.1	1.4E+09								1.6E+05	5.6E+05	4.3E+06	1.2E+05
Fosetyl-AL	39148-24-8			3.0E+00	I						1	0.1	1.4E+09								2.3E+05	8.4E+05		1.8E+05
Furans																								
-Furan	110-00-9			1.0E-03	I				V		1		1.4E+09	2.9E+03	6.8E+03						7.8E+01			7.8E+01
Furazolidone	67-45-8	3.8E+00	H								1	0.1	1.4E+09			1.7E-01	5.3E-01			1.3E-01				
Furfural	98-01-1			3.0E-03	I	5.0E-02	H				1	0.1	1.4E+09								2.3E+02	8.4E+02	7.1E+07	1.8E+02
Furium	531-82-8	1.5E+00	C	4.3E-04	C						1	0.1	1.4E+09			4.3E-01	1.3E+00	7.7E+03	3.2E-01		2.3E+02	8.4E+02		1.8E+02
Furmecycloz	60568-05-0	3.0E-02	I	8.6E-06	C						1	0.1	1.4E+09			2.1E+01	6.7E+01	3.8E+05	1.6E+01					
Glufosinate, Ammonium	77182-82-2			4.0E-04	I						1	0.1	1.4E+09								3.1E+01	1.1E+02		2.4E+01
Glycidyl	765-34-4			4.0E-04	I	1.0E-03	H				1	0.1	1.4E+09								3.1E+01	1.1E+02	1.4E+06	2.4E+01
Glyphosate	1071-83-6			1.0E-01	I						1	0.1	1.4E+09								7.8E+03	2.8E+04		6.1E+03
Goal	42874-03-3			3.0E-03	I						1	0.1	1.4E+09								2.3E+02	8.4E+02		1.8E+02
Guthion	86-50-0			3.0E-03	A	1.0E-02	A				1	0.1	1.4E+09								2.3E+02	8.4E+02	1.4E+07	1.8E+02
Haloxypol, Methyl	69806-40-2			5.0E-05	I						1	0.1	1.4E+09								3.9E+00	1.4E+01		3.1E+00
Harmony	79277-27-3			1.3E-02	I						1	0.1	1.4E+09								1.0E+03	3.6E+03		7.9E+02
Heptachlor	76-44-8	4.5E+00	I	1.3E-03	I	5.0E-04	I				1	0.1	1.4E+09			1.4E-01	4.5E-01	2.5E+03	1.1E-01		3.9E+01	1.4E+02		3.1E+01
Heptachlor Epoxide	1024-57-3	9.1E+00	I	2.6E-03	I	1.3E-05	I				1	0.1	1.4E+09			7.0E-02	2.2E-01	1.3E+03	5.3E-02		1.0E+00	3.6E+00		7.9E-01
Hexabromobenzene	87-82-1			2.0E-03	I						1	0.1	1.4E+09								1.6E+02	5.6E+02		1.2E+02
Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BDE-153)	68631-49-2			2.0E-04	I						1		1.4E+09								1.6E+01			1.6E+01
Hexachlorobenzene	118-74-1	1.6E+00	I	4.6E-04	I	8.0E-04	I				1	0.1	1.4E+09			4.0E-01	1.3E+00	7.2E+03	3.0E-01		6.3E+01	2.2E+02		4.9E+01
Hexachlorobutadiene	87-68-3	7.8E-02	I	2.2E-05	I	1.0E-03	P				1	0.1	1.4E+09			8.2E+00	2.6E+01	1.5E+05	6.2E+00		7.8E+01	2.8E+02		6.1E+01
Hexachlorocyclohexane, Alpha-	319-84-6	6.3E+00	I	1.8E-03	I	8.0E-03	A				1	0.1	1.4E+09			1.0E-01	3.2E-01	1.8E+03	7.7E-02		6.3E+02	2.2E+03		4.9E+02
Hexachlorocyclohexane, Beta-	319-85-7	1.8E+00	I	5.3E-04	I						1	0.1	1.4E+09			3.5E-01	1.1E+00	6.2E+03	2.7E-01					
Hexachlorocyclohexane, Gamma-(Lindane)	58-89-9	1.1E+00	C	3.1E-04	C	3.0E-04	I				1	0.04	1.4E+09			5.8E-01	4.6E+00	1.1E+04	5.2E-01		2.3E+01	2.1E+02		2.1E+01
Hexachlorocyclohexane, Technical	608-73-1	1.8E+00	I	5.1E-04	I						1	0.1	1.4E+09			3.5E-01	1.1E+00	6.5E+03	2.7E-01					
Hexachlorocyclopentadiene	77-47-4			6.0E-03	I	2.0E-04	I				1	0.1	1.4E+09								4.7E+02	1.7E+03	2.8E+05	3.7E+02
Hexachloroethane	67-72-1	1.4E-02	I	4.0E-06	I	1.0E-03	I				1	0.1	1.4E+09			4.6E+01	1.4E+02	8.3E+05	3.5E+01		7.8E+01	2.8E+02		6.1E+01
Hexachlorophene	70-30-4			3.0E-04	I						1	0.1	1.4E+09								2.3E+01	8.4E+01		1.8E+01
Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	1.1E-01	I	3.0E-03	I						1	0.015	1.4E+09			5.8E+00	1.2E+02		5.5E+00		2.3E+02	5.6E+03		2.3E+02
Hexamethylene Diisocyanate, 1,6-	822-06-0					1.0E-05	I	V			1		1.4E+09	3.6E+05	4.1E+03								3.7E+00	3.7E+00
Hexane, N-	110-54-3			6.0E-02	H	7.0E-01	I	V			1		1.4E+09	9.0E+02	1.4E+02						4.7E+03		6.5E+02	5.7E+02
Hexanedioic Acid	124-04-9			2.0E+00	P						1	0.1	1.4E+09								1.6E+05	5.6E+05		1.2E+05
Hexazinone	51235-04-2			3.3E-02	I						1	0.1	1.4E+09								2.6E+03	9.2E+03		2.0E+03
Hydrazine	302-01-2	3.0E+00	I	4.9E-03	I			2.0E-04	C		1		1.4E+09			2.1E-01		6.8E+02	2.1E-01				2.8E+05	2.8E+05
Hydrazine Sulfate	10034-93-2	3.0E+00	I	4.9E-03	I						1		1.4E+09			2.1E-01		6.8E+02	2.1E-01					
Hydrogen Chloride	7647-01-0					2.0E-02	I				1		1.4E+09										2.8E+07	2.8E+07
Hydrogen Fluoride	7664-39-3			4.0E-02	C	1.4E-02	C				1		1.4E+09									3.1E+03	2.0E+07	3.1E+03
Hydrogen Sulfide	7783-06-4																							

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; H = HEAST; W = WHO; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; F = See FAQ #28; c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User's Guide); s = Concentration may exceed Csat (See User's Guide); SSL values are based on DAF=1

Contaminant	CAS No.	Toxicity and Chemical-specific Information											Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1										
		SFO	k _e	IUR	k _e	RfDo	k _e	RfCi	k _e	v	muta-	GIABS	ABS	PEF	VF	Csat	Ingestion	Dermal	Inhalation	Total	Ingestion	Dermal	Inhalation	Total			
Analyte		(mg/kg-day) ⁻¹	(ug/m ³) ⁻¹	(mg/kg-day)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(m ³ /kg)	(m ³ /kg)	(mg/kg)	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg			
Lead Compounds																											
-Lead and Compounds	7439-92-1					L					1		1.4E+09											4.0E+02			
-Tetraethyl Lead	78-00-2					1.0E-07	I			1	0.1	1.4E+09									7.8E-03	2.8E-02		6.1E-03			
Linuron	330-55-2					2.0E-03	I			1	0.1	1.4E+09									1.6E+02	5.6E+02		1.2E+02			
Lithium	7439-93-2					2.0E-03	P			1		1.4E+09									1.6E+02			1.6E+02			
Lithium Perchlorate	7791-03-9					7.0E-04	I			1		1.4E+09									5.5E+01			5.5E+01			
Londax	83055-99-6					2.0E-01	I			1	0.1	1.4E+09									1.6E+04	5.6E+04		1.2E+04			
MCPA	94-74-6					5.0E-04	I			1	0.1	1.4E+09									3.9E+01	1.4E+02		3.1E+01			
MCPB	94-81-5					1.0E-02	I			1	0.1	1.4E+09									7.8E+02	2.8E+03		6.1E+02			
MCPD	93-65-2					1.0E-03	I			1	0.1	1.4E+09									7.8E+01	2.8E+02		6.1E+01			
Malathion	121-75-5					2.0E-02	I			1	0.1	1.4E+09									1.6E+03	5.6E+03		1.2E+03			
Maleic Anhydride	108-31-6					1.0E-01	I	7.0E-04	C	1	0.1	1.4E+09									7.8E+03	2.8E+04	9.9E+05	6.1E+03			
Maleic Hydrazide	123-33-1					5.0E-01	I			1	0.1	1.4E+09									3.9E+04	1.4E+05		3.1E+04			
Malononitrile	109-77-3					1.0E-04	P			1	0.1	1.4E+09									7.8E+00	2.8E+01		6.1E+00			
Mancozeb	8018-01-7					3.0E-02	H			1	0.1	1.4E+09									2.3E+03	8.4E+03		1.8E+03			
Maneb	12427-38-2					5.0E-03	I			1	0.1	1.4E+09									3.9E+02	1.4E+03		3.1E+02			
Manganese (Water)	7439-96-5					2.4E-02	I	5.0E-05	I	0.04		1.4E+09									1.9E+03		7.1E+04	1.8E+03			
Mephosfolan	950-10-7					9.0E-05	H			1	0.1	1.4E+09									7.0E+00	2.5E+01		5.5E+00			
Mepiquat Chloride	24307-26-4					3.0E-02	I			1	0.1	1.4E+09									2.3E+03	8.4E+03		1.8E+03			
Mercury Compounds																											
-Mercuric Chloride	7487-94-7					3.0E-04	I			0.07		1.4E+09									2.3E+01			2.3E+01			
-Mercuric Sulfide	1344-48-5					3.0E-04	I			1		1.4E+09									2.3E+01			2.3E+01			
-Mercury (elemental)	7439-97-6					1.6E-04	C	3.0E-04	I	V	1	1.4E+09	2.1E+04	3.1E+00							1.3E+01		6.7E+00	4.3E+00			
-Mercury, Inorganic Salts	NA					3.0E-04	I			0.07		1.4E+09									2.3E+01			2.3E+01			
-Methyl Mercury	22967-92-6					1.0E-04	I			1		1.4E+09									7.8E+00			7.8E+00			
-Phenylmercuric Acetate	62-38-4					8.0E-05	I			1	0.1	1.4E+09									6.3E+00	2.2E+01		4.9E+00			
Merphos	150-50-5					3.0E-05	I			1	0.1	1.4E+09									2.3E+00	8.4E+00		1.8E+00			
Merphos Oxide	78-48-8					3.0E-05	I			1	0.1	1.4E+09									2.3E+00	8.4E+00		1.8E+00			
Metalaxyl	57837-19-1					6.0E-02	I			1	0.1	1.4E+09									4.7E+03	1.7E+04		3.7E+03			
Methacrylonitrile	126-98-7					1.0E-04	I	7.0E-04	H	V	1	1.4E+09	7.3E+03	4.5E+03							7.8E+00		5.3E+00	3.2E+00			
Methamidophos	10265-92-6					5.0E-05	I			1	0.1	1.4E+09									3.9E+00	1.4E+01		3.1E+00			
Methanol	67-56-1					5.0E-01	I	4.0E+00	C	1	0.1	1.4E+09									3.9E+04	1.4E+05	5.7E+09	3.1E+04			
Methidathion	950-37-8					1.0E-03	I			1	0.1	1.4E+09									7.8E+01	2.8E+02		6.1E+01			
Methomyl	16752-77-5					2.5E-02	I			1	0.1	1.4E+09									2.0E+03	7.0E+03		1.5E+03			
Methoxy-5-nitroaniline, 2-	99-59-2	4.9E-02	C	1.4E-05	C					1	0.1	1.4E+09				1.3E+01	4.1E+01	2.4E+05	9.9E+00								
Methoxychlor	72-43-5					5.0E-03	I			1	0.1	1.4E+09									3.9E+02	1.4E+03		3.1E+02			
Methoxyethanol Acetate, 2-	110-49-6					2.0E-03	H	9.0E-02	C	1	0.1	1.4E+09									1.6E+02	5.6E+02	1.3E+08	1.2E+02			
Methoxyethanol, 2-	109-86-4					3.0E-03	P	2.0E-02	I	1	0.1	1.4E+09									2.3E+02	8.4E+02	2.8E+07	1.8E+02			
Methyl Acetate	79-20-9					1.0E+00	H		V	1	1.4E+09	8.8E+03	2.9E+04								7.8E+04			7.8E+04			
Methyl Acrylate	96-33-3					3.0E-02	H		V	1	1.4E+09	7.6E+03	6.9E+03								2.3E+03			2.3E+03			
Methyl Ethyl Ketone (2-Butanone)	78-93-3					6.0E-01	I	5.0E+00	I	V	1	1.4E+09	1.3E+04	2.8E+04							4.7E+04	6.7E+04		2.8E+04			
Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1					8.0E-02	H	3.0E+00	I	V	1	1.4E+09	1.1E+04	3.2E+03							6.3E+03		3.5E+04	5.3E+03			
Methyl Methacrylate	80-62-6					1.4E+00	I	7.0E-01	I	V	1	1.4E+09	6.8E+03	2.5E+03							1.1E+05		4.9E+03	4.7E+03			
Methyl Parathion	298-00-0					2.5E-04	I			1	0.1	1.4E+09									2.0E+01	7.0E+01		1.5E+01			
Methyl Phosphonic Acid	993-13-5					2.0E-02	P			1	0.1	1.4E+09									1.6E+03	5.6E+03		1.2E+03			
Methyl Styrene (Mixed Isomers)	25013-15-4					6.0E-03	H	4.0E-02	H	V	1	1.4E+09	7.6E+03	4.5E+02							4.7E+02		3.2E+02	1.9E+02			
Methyl methanesulfonate	66-27-3	9.9E-02	C	2.8E-05	C					1	0.1	1.4E+09				6.5E+00	2.0E+01	1.2E+05	4.9E+00								
Methyl tert-Butyl Ether (MTBE)	1634-04-4					1.8E-03	C	2.6E-07	C			1.4E+09	4.7E+03	6.9E+03		3.5E+02	2.0E+01	4.4E+01	3.9E+01				1.5E+04	1.5E+04			
Methyl-5-Nitroaniline, 2-	99-55-8					3.3E-02	H			1	0.1	1.4E+09				1.9E+01	6.1E+01				1.5E+01						
Methylaniline Hydrochloride, 2-	636-21-5					1.3E-01	C	3.7E-05	C	1	0.1	1.4E+09				4.9E+00	1.6E+01	8.9E+04	3.7E+00								
Methylarsonic acid	124-58-3					1.0E-02	A			1	0.1	1.4E+09									7.8E+02	2.8E+03		6.1E+02			
Methylcholanthrene, 3-	56-49-5	2.2E+01	C	6.3E-03	C					1	0.1	1.4E+09				2.9E-02	9.2E-02	5.3E+02	2.2E-02								
Methylene Chloride	75-09-2					7.5E-03	I	4.7E-07	I	6.0E-02	I	1.0E+00	A	V	1	1.4E+09	2.4E+03	3.5E+03			8.5E+01		1.1E+01	4.7E+03	2.6E+03	1.7E+03	
Methylene-bis(2-chloroaniline), 4,4'-	101-14-4					1.0E-01	P	4.3E-04	C	2.0E-03	P			M	1	0.1	1.4E+09				1.5E+00	5.1E+00	3.0E+03	1.2E+00	1.6E+02	5.6E+02	1.2E+02
Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1					4.6E-02	I	1.3E-05	C	1	0.1	1.4E+09				1.4E+01	4.4E+01	2.5E+05	1.1E+01								
Methylenebisbenzenamine, 4,4'-	101-77-9					1.6E+00	C	4.6E-04	C			2.0E-02	C	1	0.1	1.4E+09	4.0E-01	1.3E+00	7.2E+03	3.0E-01				2.8E+07	2.8E+07		
Methylenediphenyl Diisocyanate	101-68-8							6.0E-04	I	1	0.1	1.4E+09											8.5E+05	8.5E+05			
Methylstyrene, Alpha-	98-83-9					7.0E-02																					

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; H = HEAST; W = WHO; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; F = See FAQ #28; c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User's Guide); s = Concentration may exceed Csat (See User's Guide); SSL values are based on DAF=1

Contaminant	CAS No.	Toxicity and Chemical-specific Information													Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1					
		SFO	k _e	IUR	k _e	RfDo	k _e	RfCi	k _e	v _o	muta-	GIABS	ABS	PEF	VF	Csat	Ingestion	Dermal	Inhalation	Total	Ingestion	Dermal	Inhalation	Total
		(mg/kg-day) ⁻¹	y	(ug/m ³) ⁻¹	y	(mg/kg-day)	y	(mg/m ³)	y	o	gen			(m ³ /kg)	(m ³ /kg)	(mg/kg)	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg
Nickel Sub sulfide	12035-72-2	1.7E+00	C	4.8E-04	I						0.04			1.4E+09			3.8E-01		6.9E+03	3.8E-01				
Nitrate	14797-55-8					1.6E+00	I						1.4E+09								1.3E+05			1.3E+05
Nitrite	14797-65-0					1.0E-01	I						1.4E+09								7.8E+03			7.8E+03
Nitroaniline, 2-	88-74-4			3.0E-03	P	1.0E-04	P	1	0.1	1.4E+09			1.4E+09								2.3E+02	8.4E+02	1.4E+05	1.8E+02
Nitroaniline, 4-	100-01-6	2.0E-02	P			4.0E-03	P	6.0E-03	P	1	0.1	1.4E+09					3.2E+01	1.0E+02		2.4E+01	3.1E+02	1.1E+03	8.5E+06	2.4E+02
Nitrobenzene	98-95-3			4.0E-05	I	2.0E-03	I	9.0E-03	I	V	1	1	1.4E+09	7.3E+04	2.6E+03				4.4E+00	4.4E+00	1.6E+02	6.8E+02	1.3E+02	3.3E+02
Nitrofurantoin	67-20-9					7.0E-02	H				1	0.1	1.4E+09								5.5E+03	2.0E+04		4.3E+03
Nitrofurazone	59-87-0	1.3E+00	C	3.7E-04	C						1	0.1	1.4E+09				4.9E-01	1.6E+00	8.9E+03	3.7E-01				
Nitroglycerin	55-63-0	1.7E-02	P			1.0E-04	P				1	0.1	1.4E+09				3.8E+01	1.2E+02		2.9E+01	7.8E+00	2.8E+01		6.1E+00
Nitroguanidine	556-88-7					1.0E-01	I				1	0.1	1.4E+09								7.8E+03	2.8E+04		6.1E+03
Nitromethane	75-52-5			9.0E-06	P			2.0E-02	P	V	1	1	1.4E+09	1.7E+04	1.7E+04				4.7E+00	4.7E+00			3.6E+02	3.6E+02
Nitropropane, 2-	79-46-9			2.7E-03	H			2.0E-02	I	V	1	1	1.4E+09	1.3E+04	4.3E+03				1.2E-02	1.2E-02			3.6E+02	2.8E+02
Nitroso-N-ethylurea, N-	759-73-9	2.7E+01	C	7.7E-03	C						1	0.1	1.4E+09				2.4E-02	7.5E-02	4.3E-02	1.8E-02				
Nitroso-N-methylurea, N-	684-93-5	1.2E+02	C	3.4E-02	C						1	0.1	1.4E+09				5.3E-03	1.7E-02	9.7E+01	4.0E-03				
Nitroso-di-N-butylamine, N-	924-16-3	5.4E+00	I	1.6E-03	I						1	1	1.4E+09	2.8E+05	1.3E+04				4.3E-01	9.3E-02				
Nitroso-di-N-propylamine, N-	621-64-7	7.0E+00	I	2.0E-03	C						1	0.1	1.4E+09				9.1E-02	2.9E-01	1.7E+03	6.9E-02				
Nitrosodiethanolamine, N-	1116-54-7	2.8E+00	I	8.0E-04	C						1	0.1	1.4E+09				2.3E-01	7.2E-01	4.1E+03	1.7E-01				
Nitrosodiethylamine, N-	55-18-5	1.5E+02	I	4.3E-02	I						M	1	0.1	1.4E+09			9.9E-04	3.4E-03	3.0E+01	7.7E-04				
Nitrosodimethylamine, N-	62-75-9	5.1E+01	I	1.4E-02	I	8.0E-06	P				M	1	0.1	1.4E+09			2.9E-03	9.9E-03	9.3E+01	2.3E-03	6.3E-01	2.2E+00		4.9E-01
Nitrosodiphenylamine, N-	86-30-6	4.9E-03	I	2.6E-06	C						1	0.1	1.4E+09				1.3E+02	4.1E+02	1.3E+06	9.9E+01				
Nitrosomethylethylamine, N-	10595-95-6	2.2E+01	I	6.3E-03	C						1	0.1	1.4E+09				2.9E-02	9.2E-02	5.3E+02	2.2E-02				
Nitrosomorpholine [N-]	59-89-2	6.7E+00	C	1.9E-03	C						1	0.1	1.4E+09				9.5E-02	3.0E-01	1.7E+03	7.2E-02				
Nitrosopiperidine [N-]	100-75-4	9.4E+00	C	2.7E-03	C						1	0.1	1.4E+09				6.8E-02	2.2E-01	1.2E+03	5.2E-02				
Nitrosopyrrolidine, N-	930-55-2	2.1E+00	I	6.1E-04	I						1	0.1	1.4E+09				3.0E-01	9.6E-01	5.4E+03	2.3E-01				
Nitrotoluene, m-	99-08-1					2.0E-02	P				1	0.1	1.4E+09								1.6E+03	5.6E+03		1.2E+03
Nitrotoluene, o-	88-72-2	2.2E-01	P			9.0E-04	P				V	1	1.4E+09	1.4E+05	1.3E+03					2.9E+00	7.0E+01			7.0E+01
Nitrotoluene, p-	99-99-0	1.6E-02	P			4.0E-03	P				1	0.1	1.4E+09				4.0E+01	1.3E+02		3.0E+01	3.1E+02	1.1E+03		2.4E+02
Norflurazon	27314-13-2					4.0E-02	I				1	0.1	1.4E+09								3.1E+03	1.1E+04		2.4E+03
Nustar	85509-19-9					7.0E-04	I				1	0.1	1.4E+09								5.5E+01	2.0E+02		4.3E+01
Octabromodiphenyl Ether	32536-52-0					3.0E-03	I				1	0.1	1.4E+09								2.3E+02	8.4E+02		1.8E+02
Octahydro-1,3,5,7-tetra-1,3,5,7-tetra (HMx)	2691-41-0					5.0E-02	I				1	0.006	1.4E+09								3.9E+03	2.3E+05		3.8E+03
Octamethylpyrophosphoramide	152-16-9					2.0E-03	H				1	0.1	1.4E+09								1.6E+02	5.6E+02		1.2E+02
Oryzalin	19044-88-3					5.0E-02	I				1	0.1	1.4E+09								3.9E+03	1.4E+04		3.1E+03
Oxadiazon	19666-30-9					5.0E-03	I				1	0.1	1.4E+09								3.9E+02	1.4E+03		3.1E+02
Oxamyl	23135-22-0					2.5E-02	I				1	0.1	1.4E+09								2.0E+03	7.0E+03		1.5E+03
Paclobutrazol	76738-62-0					1.3E-02	I				1	0.1	1.4E+09								1.0E+03	3.6E+03		7.9E+02
Paraquat Dichloride	1910-42-5					4.5E-03	I				1	0.1	1.4E+09								3.5E+02	1.3E+03		2.7E+02
Parathion	56-38-2					6.0E-03	H				1	0.1	1.4E+09								4.7E+02	1.7E+03		3.7E+02
Pebulate	1114-71-2					5.0E-02	H				1	0.1	1.4E+09								3.9E+03	1.4E+04		3.1E+03
Pendimethalin	40487-42-1					4.0E-02	I				1	0.1	1.4E+09								3.1E+03	1.1E+04		2.4E+03
Pentabromodiphenyl Ether	32534-81-9					2.0E-03	I				1	0.1	1.4E+09								1.6E+02	5.6E+02		1.2E+02
Pentabromodiphenyl ether, 2,2',4,4',5- (BDE-99)	60348-60-9					1.0E-04	I				1	0.1	1.4E+09								7.8E+00			7.8E+00
Pentachlorobenzene	608-93-5					8.0E-04	I				1	0.1	1.4E+09								6.3E+01	2.2E+02		4.9E+01
Pentachloroethane	76-01-7	9.0E-02	P								1	0.1	1.4E+09				7.1E+00	2.2E+01		5.4E+00				
Pentachloronitrobenzene	82-68-8	2.6E-01	H			3.0E-03	I				1	0.1	1.4E+09				2.5E+00	7.8E+00		1.9E+00	2.3E+02	8.4E+02		1.8E+02
Pentachlorophenol	87-86-5	1.2E-01	I	4.6E-06	C	3.0E-02	I				1	0.25	1.4E+09				5.3E+00	6.7E+00	7.2E+05	3.0E+00	2.3E+03	3.4E+03		1.4E+03
Perchlorate and Perchlorate Salts	14797-73-0					7.0E-04	I				1	1	1.4E+09								5.5E+01			5.5E+01
Permethrin	52645-53-1					5.0E-02	I				1	0.1	1.4E+09								3.9E+03	1.4E+04		3.1E+03
Phenacetin	62-44-2	2.2E-03	C	6.3E-07	C						1	0.1	1.4E+09				2.9E+02	9.2E+02	5.3E+06	2.2E+02				
Phenmedipham	13684-63-4					2.5E-01	I				1	0.1	1.4E+09								2.0E+04	7.0E+04		1.5E+04
Phenol	108-95-2					3.0E-01	I	2.0E-01	C		1	0.1	1.4E+09								2.3E+04	8.4E+04	2.8E+08	1.8E+04
Phenylenediamine, m-	108-45-2					6.0E-03	I				1	0.1	1.4E+09								4.7E+02	1.7E+03		3.7E+02
Phenylenediamine, o-	95-54-5	4.7E-02	H								1	0.1	1.4E+09				1.4E+01	4.3E+01		1.0E+01				
Phenylenediamine, p-	106-50-3					1.9E-01	H				1	0.1	1.4E+09								1.5E+04	5.3E+04		1.2E+04
Phenylphenol, 2-	90-43-7	1.9E-03	H								1	0.1	1.4E+09				3.3E+02	1.0E+03		2.5E+02				
Phorate	298-02-2					2.0E-04	H				1	0.1	1.4E+09								1.6E+01			

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; H = HEAST; W = WHO; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; F = See FAQ #28; c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User's Guide); s = Concentration may exceed Csat (See User's Guide); SSL values are based on DAF=1

Contaminant	CAS No.	Toxicity and Chemical-specific Information										Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
		SFO	k _e	IUR	k _e	RfDo	k _e	RfCi	k _e	v _o	muta-	GIABS	ABS	PEF	VF	Csat	Ingestion	Dermal	Inhalation	Total	Ingestion	Dermal	Inhalation
Analyte		(mg/kg-day) ⁻¹	(ug/m ³) ⁻¹	(mg/kg-day)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(m ³ /kg)	(m ³ /kg)	(mg/kg)			(m ³ /kg)	(m ³ /kg)	(mg/kg)	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg
-Aroclor 1242	53469-21-9	2.0E+00	I	5.7E-04	I						1	0.14	1.4E+09			3.2E-01	7.2E-01	5.8E+03	2.2E-01				
-Aroclor 1248	12672-29-6	2.0E+00	I	5.7E-04	I						1	0.14	1.4E+09			3.2E-01	7.2E-01	5.8E+03	2.2E-01				
-Aroclor 1254	11097-69-1	2.0E+00	I	5.7E-04	I	2.0E-05	I				1	0.14	1.4E+09			3.2E-01	7.2E-01	5.8E+03	2.2E-01	1.6E+00	4.0E+00		1.1E+00
-Aroclor 1260	11096-82-5	2.0E+00	I	5.7E-04	I						1	0.14	1.4E+09			3.2E-01	7.2E-01	5.8E+03	2.2E-01				
-Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	1.3E+01	C	3.8E-03	C						1	0.14	1.4E+09			4.9E-02	1.1E-01	8.7E+02	3.4E-02				
-Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167)	52663-72-6	1.3E+01	C	3.8E-03	C						1	0.14	1.4E+09			4.9E-02	1.1E-01	8.7E+02	3.4E-02				
-Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	6.5E+02	C	1.9E-02	C						1	0.14	1.4E+09			9.8E-04	2.2E-03	1.7E+02	6.8E-04				
-Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 156)	38380-08-4	6.5E+02	C	1.9E-02	C						1	0.14	1.4E+09			9.8E-04	2.2E-03	1.7E+02	6.8E-04				
-Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	1.3E+01	C	3.8E-03	C						1	0.14	1.4E+09			4.9E-02	1.1E-01	8.7E+02	3.4E-02				
-Pentachlorobiphenyl, 2',3,4,4',5- (PCB 123)	65510-44-3	1.3E+01	C	3.8E-03	C						1	0.14	1.4E+09			4.9E-02	1.1E-01	8.7E+02	3.4E-02				
-Pentachlorobiphenyl, 2,3,4,4',5- (PCB 118)	31508-00-6	1.3E+01	C	3.8E-03	C						1	0.14	1.4E+09			4.9E-02	1.1E-01	8.7E+02	3.4E-02				
-Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	1.3E+01	C	3.8E-03	C						1	0.14	1.4E+09			4.9E-02	1.1E-01	8.7E+02	3.4E-02				
-Pentachlorobiphenyl, 2,3,4,4',5- (PCB 114)	74472-37-0	6.5E+02	C	1.9E-02	C						1	0.14	1.4E+09			9.8E-04	2.2E-03	1.7E+02	6.8E-04				
-Pentachlorobiphenyl, 3,3',4,4',5- (PCB 126)	57465-28-8	1.3E+04	C	3.8E+00	C						1	0.14	1.4E+09			4.9E-05	1.1E-04	8.7E-01	3.4E-05				
-Polychlorinated Biphenyls (high risk)	1336-36-3	2.0E+00	I	5.7E-04	C						1	0.14	1.4E+09			3.2E-01	7.2E-01	5.8E+03	2.2E-01				
-Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	1.3E+01	C	3.8E-03	C						1	0.14	1.4E+09			4.9E-02	1.1E-01	8.7E+02	3.4E-02				
-Tetrachlorobiphenyl, 3,4,4',5- (PCB 81)	70362-50-4	1.3E+01	C	3.8E-03	C						1	0.14	1.4E+09			4.9E-02	1.1E-01	8.7E+02	3.4E-02				
Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9							6.0E-04	I		1	0.1	1.4E+09									8.5E+05	8.5E+05
Polynuclear Aromatic Hydrocarbons (PAHs)																							
-Acenaphthene	83-32-9					6.0E-02	I			V	1	0.13	1.4E+09	1.7E+05						4.7E+03	1.3E+04		3.4E+03
-Anthracene	120-12-7					3.0E-01	I			V	1	0.13	1.4E+09	6.3E+05						2.3E+04	6.4E+04		1.7E+04
-Benz[a]anthracene	56-55-3	7.3E-01	I	1.1E-04	C					M	1	0.13	1.4E+09			2.0E-01	5.3E-01	1.2E+04	1.5E-01				
-Benzo[a]pyrene	50-32-8	7.3E+00	I	1.1E-03	C					M	1	0.13	1.4E+09			2.0E-02	5.3E-02	1.2E+03	1.5E-02				
-Benzo[b]fluoranthene	205-99-2	7.3E-01	I	1.1E-04	C					M	1	0.13	1.4E+09			2.0E-01	5.3E-01	1.2E+04	1.5E-01				
-Benzo[k]fluoranthene	207-08-9	7.3E-02	I	1.1E-04	C					M	1	0.13	1.4E+09			2.0E+00	5.3E+00	1.2E+04	1.5E+00				
-Chrysene	218-01-9	7.3E-03	I	1.1E-05	C					M	1	0.13	1.4E+09			2.0E+01	5.3E+01	1.2E+05	1.5E+01				
-Dibenz[a,h]anthracene	53-70-3	7.3E+00	I	1.2E-03	C					M	1	0.13	1.4E+09			2.0E-02	5.3E-02	1.1E+03	1.5E-02				
-Dimethylbenz[a]anthracene, 7,12-	57-97-6	2.5E+02	C	7.1E-02	C						1	0.13	1.4E+09			2.6E-03	6.2E-03	4.7E+01	1.8E-03				
-Fluoranthene	206-44-0					4.0E-02	I				1	0.13	1.4E+09							3.1E+03	8.6E+03		2.3E+03
-Fluorene	86-73-7					4.0E-02	I			V	1	0.13	1.4E+09	3.4E+05						3.1E+03	8.6E+03		2.3E+03
-Indeno[1,2,3-cd]pyrene	193-39-5	7.3E-01	I	1.1E-04	C					M	1	0.13	1.4E+09			2.0E-01	5.3E-01	1.2E+04	1.5E-01				
-Methylnaphthalene, 1-	90-12-0	2.9E-02	P			7.0E-02	A			V	1	1.4E+09	6.9E+04	4.6E+02					2.2E+01			5.5E+03	5.5E+03
-Methylnaphthalene, 2-	91-57-6					4.0E-03	I			V	1	1.4E+09	6.8E+04	4.4E+02								3.1E+02	3.1E+02
-Naphthalene	91-20-3					2.0E-02	I	3.0E-03	I	V	1	0.13	1.4E+09	5.4E+04				3.9E+00	3.9E+00	1.6E+03	4.3E+03	1.7E+02	1.5E+02
-Pyrene	129-00-0					3.0E-02	I			V	1	0.13	1.4E+09	2.9E+06						2.3E+03	6.4E+03		1.7E+03
Potassium Perchlorate	7778-74-7					7.0E-04	I				1	1.4E+09								5.5E+01			5.5E+01
Prochloraz	67747-09-5	1.5E-01	I			9.0E-03	I				1	0.1	1.4E+09			4.3E+00	1.3E+01			7.0E+02	2.5E+03		5.5E+02
Profluralin	26399-36-0					6.0E-03	H				1	0.1	1.4E+09							4.7E+02	1.7E+03		3.7E+02
Prometon	1610-18-0					1.5E-02	I				1	0.1	1.4E+09							1.2E+03	4.2E+03		9.2E+02
Prometryn	7287-19-6					4.0E-03	I				1	0.1	1.4E+09							3.1E+02	1.1E+03		2.4E+02
Propachlor	1918-16-7					1.3E-02	I				1	0.1	1.4E+09							1.0E+03	3.6E+03		7.9E+02
Propanil	709-98-8					5.0E-03	I				1	0.1	1.4E+09							3.9E+02	1.4E+03		3.1E+02
Propargite	2312-35-8					2.0E-02	I				1	0.1	1.4E+09							1.6E+03	5.6E+03		1.2E+03
Propargyl Alcohol	107-19-7					2.0E-03	I				1	0.1	1.4E+09							1.6E+02	5.6E+02		1.2E+02
Propazine	139-40-2					2.0E-02	I				1	0.1	1.4E+09							1.6E+03	5.6E+03		1.2E+03
Propham	122-42-9					2.0E-02	I				1	0.1	1.4E+09							1.6E+03	5.6E+03		1.2E+03
Propiconazole	60207-90-1					1.3E-02	I				1	0.1	1.4E+09							1.0E+03	3.6E+03		7.9E+02
Propionaldehyde	123-38-6							8.0E-03	I	V	1	1.4E+09	1.0E+04	3.6E+04								8.6E+01	8.6E+01
Propylene Glycol	57-55-6					2.0E+01	P				1	0.1	1.4E+09							1.6E+06	5.6E+06		1.2E+06
Propylene Glycol Dinitrate	6423-43-4							2.7E-04	A	V	1	1.4E+09	2.1E+05	1.4E+03								6.0E+01	6.0E+01
Propylene Glycol Monoethyl Ether	1569-02-4					7.0E-01	H				1	0.1	1.4E+09							5.5E+04	2.0E+05		4.3E+04
Propylene Glycol Monomethyl Ether	107-98-2					7.0E-01	H	2.0E+00	I		1	0.1	1.4E+09							5.5E+04	2.0E+05	2.8E+09	4.3E+04
Propylene Oxide	75-56-9	2.4E-01	I	3.7E-06	I			3.0E-02	I	V	1	1.4E+09	9.6E+03	6.8E+04					2.7E+00		6.3E+00	1.9E+00	3.0E+02
Pursuit	81335-77-5					2.5E-01	I				1	0.1	1.4E+09							2.0E+04	7.0E+04		1.5E+04
Pydrin	51630-58-1					2.5E-02	I				1	0.1	1.4E+09							2.0E+03	7.0E+03		1.5E+03
Pyridine	110-86-1					1.0E-03	I			V	1	1.4E+09	4.5E+04	3.0E+05						7.8E+01			7.8E+01
Quinalphos	13593-03-8					5.0E-04																	

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; H = HEAST; W = WHO; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; F = See FAQ #28; c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User's Guide); s = Concentration may exceed Csat (See User's Guide); SSL values are based on DAF=1

Contaminant	CAS No.	Toxicity and Chemical-specific Information											Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1								
		SFO	k _e	IUR	k _e	RfDo	k _e	RfCi	k _e	v	muta-	GIABS	ABS	PEF	VF	Csat	Ingestion	Dermal	Inhalation	Total	Ingestion	Dermal	Inhalation	Total	
Analyte		(mg/kg-day) ⁻¹	(ug/m ³) ⁻¹	(mg/kg-day)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(mg/m ³)	(m ³ /kg)	(m ³ /kg)	(mg/kg)	(m ³ /kg)	(m ³ /kg)	(m ³ /kg)	(m ³ /kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg)		
Sodium Azide	26628-22-8			4.0E-03	I						1		1.4E+09										3.1E+02		
Sodium Diethyldithiocarbamate	148-18-5	2.7E-01	H	3.0E-02	I						1	0.1	1.4E+09				2.4E+00	7.5E+00		1.8E+00			2.3E+03	8.4E+03	3.1E+02
Sodium Fluoride	7681-49-4			5.0E-02	A						1		1.4E+09										3.9E+03		
Sodium Fluoroacetate	62-74-8			2.0E-05	I						1	0.1	1.4E+09									1.6E+00	5.6E+00	3.9E+03	
Sodium Metavanadate	13718-26-8			1.0E-03	H						1		1.4E+09										7.8E+01		
Sodium Perchlorate	7601-89-0			7.0E-04	I						1		1.4E+09										5.5E+01		
Stirofos (Tetrachlorovinphos)	961-11-5	2.4E-02	H	3.0E-02	I						1	0.1	1.4E+09				2.7E+01	8.4E+01		2.0E+01			2.3E+03	8.4E+03	
Strontium, Stable	7440-24-6			6.0E-01	I						1		1.4E+09										4.7E+04		
Strychnine	57-24-9			3.0E-04	I						1	0.1	1.4E+09										2.3E+01	8.4E+01	
Styrene	100-42-5			2.0E-01	I	1.0E+00	I	V			1		1.4E+09	1.1E+04	1.0E+03								1.6E+04	1.1E+04	
Sulfonfylbis(4-chlorobenzene), 1,1'-	80-07-9			5.0E-03	P						1	0.1	1.4E+09										3.9E+02	1.4E+03	
Systhane	88671-89-0			2.5E-02	I						1	0.1	1.4E+09										2.0E+03	7.0E+03	
TCMTB	21564-17-0			3.0E-02	H						1	0.1	1.4E+09										2.3E+03	8.4E+03	
Tebuthiuron	34014-18-1			7.0E-02	I						1	0.1	1.4E+09										5.5E+03	2.0E+04	
Temephos	3383-96-8			2.0E-02	H						1	0.1	1.4E+09										1.6E+03	5.6E+03	
Terbacil	5902-51-2			1.3E-02	I						1	0.1	1.4E+09										1.0E+03	3.6E+03	
Terbufos	13071-79-9			2.5E-05	H						1	0.1	1.4E+09										2.0E+00	7.0E+00	
Terbutryn	886-50-0			1.0E-03	I						1	0.1	1.4E+09										7.8E+01	2.8E+02	
Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1			1.0E-04	I						1		1.4E+09										7.8E+00		
Tetrachlorobenzene, 1,2,4,5-	95-94-3			3.0E-04	I						1	0.1	1.4E+09										2.3E+01	8.4E+01	
Tetrachloroethane, 1,1,1,2-	630-20-6	2.6E-02	I	7.4E-06	I	3.0E-02	I		V		1		1.4E+09	6.5E+03	7.5E+02	2.5E+01		2.1E+00	2.0E+00				2.3E+03		
Tetrachloroethane, 1,1,2,2-	79-34-5	2.0E-01	I	5.8E-05	I	4.0E-03	P		V		1		1.4E+09	1.7E+04	2.1E+03	3.2E+00		7.2E-01	5.9E-01				3.1E+02		
Tetrachloroethylene	127-18-4	5.4E-01	C	5.9E-06	C	1.0E-02	I	2.7E-01	A	V	1		1.4E+09	2.6E+03	1.8E+02	1.2E+00		1.1E+00	5.7E-01				7.5E+02		
Tetrachlorophenol, 2,3,4,6-	58-90-2			3.0E-02	I						1	0.1	1.4E+09										2.3E+03	8.4E+03	
Tetrachlorotoluene, p- alpha, alpha, alpha-	5216-25-1	2.0E+01	H								1	0.1	1.4E+09				3.2E-02	1.0E-01		2.4E-02					
Tetraethyl Dithiopyrophosphate	3689-24-5			5.0E-04	I						1	0.1	1.4E+09										3.9E+01	1.4E+02	
Tetrafluoroethane, 1,1,1,2-	811-97-2							8.0E+01	I	V	1		1.4E+09	1.4E+03	8.2E+02								1.1E+05		
Tetryl (Trinitrophenylmethylnitramine)	479-45-8			4.0E-03	P						1	0.1	1.4E+09										3.1E+02	1.1E+03	
Thallium (I) Nitrate	10102-45-1			9.0E-05	I						1		1.4E+09										7.0E+00		
Thallium (Soluble Salts)	7440-28-0			6.5E-05	I						1		1.4E+09										5.1E+00		
Thallium Acetate	563-68-8			9.0E-05	I						1		1.4E+09										7.0E+00		
Thallium Carbonate	6533-73-9			8.0E-05	I						1		1.4E+09										6.3E+00		
Thallium Chloride	7791-12-0			8.0E-05	I						1		1.4E+09										6.3E+00		
Thallium Sulfate	7446-18-6			8.0E-05	I						1		1.4E+09										6.3E+00		
Thiobencarb	28249-77-6			1.0E-02	I						1	0.1	1.4E+09										7.8E+02	2.8E+03	
Thiofanox	39196-18-4			3.0E-04	H						1	0.1	1.4E+09										2.3E+01	8.4E+01	
Thiophanate, Methyl	23564-05-8			8.0E-02	I						1	0.1	1.4E+09										6.3E+03	2.2E+04	
Thiram	137-26-8			5.0E-03	I						1	0.1	1.4E+09										3.9E+02	1.4E+03	
Tin	7440-31-5			6.0E-01	H						1		1.4E+09										4.7E+04		
Titanium Tetrachloride	7550-45-0							1.0E-04	A		1		1.4E+09												
Toluene	108-88-3			8.0E-02	I	5.0E+00	I	V			1		1.4E+09	4.9E+03	9.3E+02								6.3E+03		
Toluene diisocyanate mixture (TDI)	26471-62-5	3.9E-02	C	1.1E-05	C			7.0E-05	I	V	1		1.4E+09	7.5E+05	2.1E+03	1.6E+01		1.6E+02	1.5E+01				5.4E+01		
Toluene-2,4-diamine	95-80-7	3.8E+00	C	1.1E-03	C						1	0.1	1.4E+09				1.7E-01	5.3E-01	3.0E+03	1.3E-01					
Toluene-2,5-diamine	95-70-5			6.0E-01	H						1	0.1	1.4E+09										4.7E+04	1.7E+05	
Toluene-2,6-diamine	823-40-5			3.0E-02	P						1	0.1	1.4E+09										2.3E+03	8.4E+03	
Toluidine, o- (Methylaniline, 2-)	95-53-4	1.8E-01	C	5.1E-05	C						1	0.1	1.4E+09				3.5E+00	1.1E+01	6.5E+04	2.7E+00					
Toluidine, p-	106-49-0	1.9E-01	H								1	0.1	1.4E+09				3.4E+00	1.1E+01		2.6E+00					
Toxaphene	8001-35-2	1.1E+00	I	3.2E-04	I						1	0.1	1.4E+09				5.8E-01	1.8E+00	1.0E+04	4.4E-01					
Tralomethrin	66841-25-6			7.5E-03	I						1	0.1	1.4E+09										5.9E+02	2.1E+03	
Tri-n-butyltin	688-73-3			3.0E-04	A						1	0.1	1.4E+09										2.3E+01	8.4E+01	
Triallate	2303-17-5			1.3E-02	I						1	0.1	1.4E+09										1.0E+03	3.6E+03	
Triasulfuron	82097-50-5			1.0E-02	I						1	0.1	1.4E+09										7.8E+02	2.8E+03	
Tribromobenzene, 1,2,4-	615-54-3			5.0E-03	I						1	0.1	1.4E+09										3.9E+02	1.4E+03	
Tributyl Phosphate	126-73-8	9.2E-03	P			2.0E-01	P				1	0.1	1.4E+09				6.9E+01	2.2E+02		5.3E+01			1.6E+04	5.6E+04	
Tributyltin Compounds	NA			3.0E-04	P						1	0.1	1.4E+09										2.3E+01	8.4E+01	
Tributyltin Oxide	56-35-9			3.0E-04	I						1	0.1	1.4E+09										2.3E+01	8.4E+01	
Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1			3.0E+01	I	3.0E+01	H	V			1		1.4E+09	1.4E+03	9.4E+02								2.3E+06	4.4E+04	
Trichloroaniline HCl, 2,4,6-	33663-50-2	2.9E-02	H								1	0.1	1.4E+09				2.2E+01	7.0E+01		1.7E+01					
Trichloroaniline, 2,4,6-	634-93-5	3.4E-02	H								1	0.1	1.4E+09				1.9E+01	6.0E+01		1.4E+01					
Trichlorobenzene, 1,2,4-	120-82-1	3.6E-03	C			1.0E-02	I	4.0E-03	P	V	1		1.4E+09	2.4E+04	2.2E										

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; H = HEAST; W = WHO; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; F = See FAQ #28; c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User's Guide); s = Concentration may exceed Csat (See User's Guide); SSL values are based on DAF=1

Contaminant	CAS No.	Toxicity and Chemical-specific Information											Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1			
		SFO (mg/kg-day) ⁻¹	k e IUR (ug/m ³) ⁻¹	k e RfDo (mg/kg-day)	k e RfCi (mg/m ³)	k e v o l u t a g e n	GIABS	ABS	PEF (m ³ /kg)	VF (m ³ /kg)	Csat (mg/kg)	Ingestion mg/kg	Dermal mg/kg	Inhalation mg/kg	Total mg/kg	Ingestion mg/kg	Dermal mg/kg	Inhalation mg/kg	Total mg/kg	
Tridiphane	58138-08-2			3.0E-03	I		1	0.1	1.4E+09						2.3E+02	8.4E+02		1.8E+02		
Triethylamine	121-44-8					7.0E-03	I V	1	1.4E+09	2.3E+04	5.5E+04						1.7E+02	1.7E+02		
Trifluralin	1582-09-8	7.7E-03	I	7.5E-03	I		1	0.1	1.4E+09				8.3E+01	2.6E+02	6.3E+01	5.9E+02	2.1E+03	4.6E+02		
Trimethyl Phosphate	512-56-1	3.7E-02	H				1	0.1	1.4E+09				1.7E+01	5.5E+01	1.3E+01					
Trimethylbenzene, 1,2,4-	95-63-6					7.0E-03	P V	1	1.4E+09	9.2E+03	2.5E+02						6.7E+01	6.7E+01		
Trimethylbenzene, 1,3,5-	108-67-8			5.0E-02	P	6.0E-03	P V	1	1.4E+09	7.7E+03	2.1E+02				3.9E+03		4.8E+01	4.7E+01		
Trinitrobenzene, 1,3,5-	99-35-4			3.0E-02	I		1	0.019	1.4E+09						2.3E+03	4.4E+04		2.2E+03		
Trinitrotoluene, 2,4,6-	118-96-7	3.0E-02	I	5.0E-04	I		1	0.032	1.4E+09				2.1E+01	2.1E+02	1.9E+01			3.6E+01		
Triphenylphosphine Oxide	791-28-6			2.0E-02	P		1	0.1	1.4E+09						1.6E+03	5.6E+03		1.2E+03		
Tris(2-chloroethyl)phosphate	115-96-8	1.4E-02	P	3.0E-01	P		1	0.1	1.4E+09				4.6E+01	1.4E+02	3.5E+01	2.3E+04	8.4E+04	1.8E+04		
Tris(2-ethylhexyl)phosphate	78-42-2	3.2E-03	P	1.0E-01	P		1	0.1	1.4E+09				2.0E+02	6.3E+02	1.5E+02	7.8E+03	2.8E+04	6.1E+03		
Uranium (Soluble Salts)	NA			3.0E-03	I	3.0E-04	A		1.4E+09						2.3E+02		4.3E+05	2.3E+02		
Vanadium Pentoxide	1314-62-1			8.3E-03	P	9.0E-03	I	7.0E-06	P	0.026					4.0E+02	4.0E+02	7.0E+02	6.6E+02		
Vanadium Sulfate	36907-42-3			2.0E-02	H				0.026						1.6E+03		9.9E+03	1.6E+03		
Vanadium and Compounds	NA			5.0E-03	I				1						3.9E+02			3.9E+02		
Vanadium, Metallic	7440-62-2			7.0E-03	H				0.026						5.5E+02			5.5E+02		
Vernolate	1929-77-7			1.0E-03	I				1	0.1					7.8E+01	2.8E+02		6.1E+01		
Vinclozolin	50471-44-8			2.5E-02	I				1	0.1					2.0E+03	7.0E+03		1.5E+03		
Vinyl Acetate	108-05-4			1.0E+00	H	2.0E-01	I V	1	1.4E+09	4.8E+03	2.8E+03						7.8E+04	1.0E+03		
Vinyl Bromide	593-60-2					3.0E-03	I V	1	1.4E+09	1.5E+03	1.7E+03						4.7E+00	4.7E+00		
Vinyl Chloride	75-01-4	7.2E-01	I	4.4E-06	I	3.0E-03	I	1.0E-01	I V M	1			9.3E-02		1.7E-01	6.0E-02	2.3E+02	1.1E+02		
Warfarin	81-81-2			3.0E-04	I				1	0.1					2.3E+01	8.4E+01		1.8E+01		
Xylene, Mixture	1330-20-7			2.0E-01	I	1.0E-01	I V	1	1.4E+09	5.9E+03	3.0E+02				1.6E+04		6.2E+02	6.0E+02		
Xylene, P-	106-42-3					7.0E-01	C V	1	1.4E+09	6.4E+03	4.5E+02						4.7E+03	4.7E+03		
Xylene, m-	108-38-3			2.0E+00	H	7.0E-01	C V	1	1.4E+09	6.3E+03	4.4E+02				1.6E+05		4.6E+03	4.5E+03		
Xylene, o-	95-47-6			2.0E+00	H	7.0E-01	C V	1	1.4E+09	7.4E+03	3.0E+02				1.6E+05		5.4E+03	5.3E+03		
Zinc (Metallic)	7440-66-6			3.0E-01	I				1						2.3E+04			2.3E+04		
Zinc Phosphide	1314-84-7			3.0E-04	I				1						2.3E+01			2.3E+01		
Zineb	12122-67-7			5.0E-02	I				1	0.1					3.9E+03	1.4E+04		3.1E+03		