Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	GIABS	ABS_d	C_{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
				3.00E-04	O					1	0.1		Acephate	30560-19-1	1.90E+01	n	2.50E+02	n					6.00E+00	n	
		2.20E-06	I			9.00E-03	I	V		1		1.07E+05	Acetaldehyde	75-07-0	8.20E+01	n	3.40E+02	n	9.40E+00	n	3.90E+01	n	1.90E+01	n	
				2.00E-02	I					1	0.1		Acetochlor	34256-82-1	1.30E+03	n	1.60E+04	n					3.50E+02	n	
				9.00E-01	I			V		1		1.14E+05	Acetone	67-64-1	7.00E+04	n	1.10E+06	nms					1.80E+04	n	
						2.00E-03	X			1	0.1		Acetone Cyanohydrin	75-86-5	2.80E+06	nm	1.20E+07	nm	2.10E+00	n	8.80E+00	n			
						6.00E-02	I	V		1		1.28E+05	Acetonitrile	75-05-8	8.10E+02	n	3.40E+03	n	6.30E+01	n	2.60E+02	n	1.30E+02	n	
				1.00E-01	I			V		1		2.52E+03	Acetophenone	98-86-2	7.80E+03	ns	1.20E+05	nms					1.90E+03	n	
3.80E+00	C	1.30E-03	C							1	0.1		Acetylaminofluorene, 2-	53-96-3	1.40E+01	c	6.00E+01	c	2.20E-01	c	9.40E-01	c	1.60E+00	c	
				5.00E-04	I	2.00E-05	I	V		1		2.27E+04	Acrolein	107-02-8	1.40E-01	n	6.00E-01	n	2.10E-02	n	8.80E-02	n	4.20E-02	n	
5.00E-01	I	1.00E-04	I	2.00E-03	I	6.00E-03	I		M	1	0.1		Acrylamide	79-06-1	2.40E+01	c**	4.60E+02	c**	1.00E+00	c**	1.20E+01	c**	5.00E+00	c**	
				5.00E-01	I	2.00E-04	P	V		1		1.09E+05	Acrylic Acid	79-10-7	2.00E+01	n	8.30E+01	n	2.10E-01	n	8.80E-01	n	4.20E-01	n	
5.40E-01	I	6.80E-05	I	1.00E-02	A	2.00E-03	I	V		1		1.13E+04	Acrylonitrile	107-13-1	1.60E+01	n	6.70E+01	n	2.10E+00	n	8.80E+00	n	4.10E+00	n	
						6.00E-03	P			1	0.1		Adiponitrile	111-69-3	8.50E+06	nm	3.60E+07	nm	6.30E+00	n	2.60E+01	n			
5.60E-02	C			1.00E-02	I					1	0.1		Alachlor	15972-60-8	6.30E+02	n	4.10E+03	c**					1.10E+02	c**	2.00E+00
				1.00E-03	I					1	0.1		Aldicarb	116-06-3	6.30E+01	n	8.20E+02	n					2.00E+01	n	3.00E+00
				1.00E-03	I					1	0.1		Aldicarb Sulfone	1646-88-4	6.30E+01	n	8.20E+02	n					2.00E+01	n	2.00E+00
										1	0.1		Aldicarb sulfoxide	1646-87-3											4.00E+00
1.70E+01	I	4.90E-03	I	3.00E-05	I			V		1			Aldrin	309-00-2	2.30E+00	n	1.80E+01	c**	5.70E-02	c	2.50E-01	c	9.20E-02	c**	
				4.00E-03	P	1.00E-04	X	V		1		1.11E+05	Allyl Alcohol	107-18-6	3.50E+00	n	1.50E+01	n	1.00E-01	n	4.40E-01	n	2.10E-01	n	
2.10E-02	C	6.00E-06	C			1.00E-03	I	V		1		1.42E+03	Allyl Chloride	107-05-1	1.70E+00	n	6.90E+00	n	1.00E+00	n	4.40E+00	n	2.10E+00	n	
				1.00E+00	P	5.00E-03	P			1			Aluminum	7429-90-5	7.70E+04	n	1.10E+06	nm	5.20E+00	n	2.20E+01	n	2.00E+04	n	
				4.00E-04	I					1			Aluminum Phosphide	20859-73-8	3.10E+01	n	4.70E+02	n					8.00E+00	n	
				9.00E-03	I					1	0.1		Ametryn	834-12-8	5.70E+02	n	7.40E+03	n					1.50E+02	n	
2.10E+01	C	6.00E-03	C							1	0.1		Aminobiphenyl, 4-	92-67-1	2.60E+00	c	1.10E+01	c	4.70E-02	c	2.00E-01	c	3.00E-01	c	
				8.00E-02	P					1	0.1		Aminophenol, m-	591-27-5	5.10E+03	n	6.60E+04	n					1.60E+03	n	
				4.00E-03	X					1	0.1		Aminophenol, o-	95-55-6	2.50E+02	n	3.30E+03	n					7.90E+01	n	
				2.00E-02	P					1	0.1		Aminophenol, p-	123-30-8	1.30E+03	n	1.60E+04	n					4.00E+02	n	
				2.50E-03	I					1	0.1		Amitraz	33089-61-1	1.60E+02	n	2.10E+03	n					8.20E+00	n	
						5.00E-01	I	V		1			Ammonia	7664-41-7					5.20E+02	n	2.20E+03	n			
				2.00E-03	X					1	0.1		Ammonium Picrate	131-74-8	1.30E+02	n	1.60E+03	n					4.00E+01	n	
				2.00E-01	I					1			Ammonium Sulfamate	7773-06-0	1.60E+04	n	2.30E+05	nm					4.00E+03	n	
						3.00E-03	X	V		1		1.37E+04	Amyl Alcohol, tert-	75-85-4	8.20E+01	n	3.40E+02	n	3.10E+00	n	1.30E+01	n	6.30E+00	n	
5.70E-03	I	1.60E-06	C	7.00E-03	P	1.00E-03	I			1	0.1		Aniline	62-53-3	4.40E+02	n	5.70E+03	n	1.00E+00	n	4.40E+00	n	1.40E+02	n	
4.00E-02	P			2.00E-03	X					1	0.1		Anthraquinone, 9,10-	84-65-1	1.30E+02	n	1.60E+03	n					3.00E+01	n	
				4.00E-04	I	3.00E-04	A			0.15			Antimony (metallic)	7440-36-0	3.10E+01	n	4.70E+02	n	3.10E-01	n	1.30E+00	n	7.80E+00	n	6.00E+00
				5.00E-04	H					0.15			Antimony Pentoxide	1314-60-9	3.90E+01	n	5.80E+02	n					9.70E+00	n	
				4.00E-04	H					0.15			Antimony Tetroxide	1332-81-6	3.10E+01	n	4.70E+02	n					7.80E+00	n	
						2.00E-04	I			0.15			Antimony Trioxide	1309-64-4	2.80E+05	nm	1.20E+06	nm	2.10E-01	n	8.80E-01	n			
1.50E+00	I	4.30E-03	I	3.00E-04	I	1.50E-05	C			1	0.03		Arsenic, Inorganic	7440-38-2	3.50E+01	nR	3.00E+02	c**R	1.60E-02	n	6.60E-02	n	5.20E+00	c**	1.00E+01
				3.50E-06	C	5.00E-05	I			1			Arsine	7784-42-1	2.70E-01	n	4.10E+00	n	5.20E-02	n	2.20E-01	n	7.00E-02	n	
										1			Asbestos (units in fibers)	1332-21-4											7.0E+06(G)
				3.60E-02	O					1	0.1		Asulam	3337-71-1	2.30E+03	n	3.00E+04	n					7.20E+02	n	
2.30E-01	C			3.00E-03	A					1	0.1		Atrazine	1912-24-9	1.90E+02	n	1.00E+03	c**					3.00E+01	c**	3.00E+00
8.80E-01	C	2.50E-04	C							1	0.1		Auramine	492-80-8	6.20E+01	c	2.60E+02	c	1.10E+00	c	4.90E+00	c	7.80E+00	c	
				4.00E-04	I					1	0.1		Avermectin B1	65195-55-3	2.50E+01	n	3.30E+02	n					8.00E+00	n	
				3.00E-03	A	1.00E-02	A			1	0.1		Azinphos-methyl	86-50-0	1.90E+02	n	2.50E+03	n	1.00E+01	n	4.40E+01	n	5.60E+01	n	
1.10E-01	I	3.10E-05	I					V		1			Azobenzene	103-33-3	5.60E+02	c	2.60E+03	c	9.10E+00	c	4.00E+01	c	1.20E+01	c	
				1.00E+00	P	7.00E-06	P			1	0.1		Azodicarbonamide	123-77-3	8.60E+03	n	4.00E+04	n	7.30E-03	n	3.10E-02	n	2.00E+04	n	
				2.00E-01	I	5.00E-04	H			0.07			Barium	7440-39-3	1.50E+04	n	2.20E+05	nm	5.20E-01	n	2.20E+00	n	3.80E+03	n	2.00E+03
				5.00E-03	O			V		1			Benfluralin	1861-40-1	3.90E+02	n	5.80E+03	n					2.80E+01	n	
				5.00E-02	I					1	0.1		Benomyl	17804-35-2	3.20E+03	n	4.10E+04	n					9.70E+02	n	
				2.00E-01	I					1	0.1		Bensulfuron-methyl	83055-99-6	1.30E+04	n	1.60E+05	nm					3.90E+03	n	
				3.00E-02	I					1	0.1		Bentazon	25057-89-0	1.90E+03	n	2.50E+04	n					5.70E+02	n	
4.00E-03	P			1.00E-01	I			V		1		1.16E+03	Benzaldehyde	100-52-7	7.80E+03	ns	8.20E+04	c**s					1.90E+03	c**	
5.50E-02	I	7.80E-06	I	4.00E-03	I	3.00E-02	I	V		1		1.82E+03	Benzene	71-43-2	8.20E+01	n	4.20E+02	n	3.10E+01	n	1.30E+02	n	3.30E+01	n	5.00E+00
1.00E-01	X			3.00E-04	X					1	0.1		Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1	1.90E+01	n	2.50E+02	n					6.00E+00	n	
				1.00E-03	P			V		1		1.26E+03	Benzenethiol	108-98-5	7.80E+01	n	1.20E+03	n					1.70E+01	n	
2.30E+02	I	6.70E-02	I	3.00E-03	I				M	1	0.1		Benzidine	92-87-5	5.30E-02	c	1.00E+00	c	1.50E-03	c	1.80E-02	c	1.10E-02	c	
				4.00E+00	I					1	0.1		Benzoic Acid	65-85-0	2.50E+05	nm	3.30E+06	nm					7.50E+04	n	
1.30E+01	I							V		1		3.24E+02	Benzotrichloride	98-07-7	5.30E+00	c	2.50E+01	c					3.00E-01	c	
				1.00E-01	P					1	0.1		Benzyl Alcohol	100-51-6	6.30E+03	n	8.20E+04	n					2.00E+03	n	
1.70E-01	I	4.90E-05	C	2.00E-03	P	1.00E-03	P	V		1		1.46E+03	Benzyl Chloride	100-44-7	2.30E+01	n	1.10E+02	n	1.00E+00	n	4.40E+00	n	2.00E+00	n	
		2.40E-03	I	2.00E-03	I	2.00E-05	I			0.007			Beryllium and compounds	7440-41-7	1.60E+02	n	2.30E+03	n	2.10E-02	n	8.80E-02	n	2.50E+01	n	4.00E+00
				9.00E-03	P					1	0.1		Bifenox	42576-02-3	5.70E+02	n	7.40E+03	n					1.00E+02	n	
				1.50E-02	I					1	0.1		Biphenthrin	82657-04-3	9.50E+02	n	1.20E+04	n					3.00E+02	n	
8.00E-03	I			5.00E-01	I	4.00E-04	X	V		1			Biphenyl, 1,1'-	92-52-4	4.70E+01	n	2.00E+02	n	4.20E-01	n	1.80E+00	n	8.30E-01	n	
				4.00E-02	I			V		1		1.02E+03	Bis(2-chloro-1-methylethyl) ether	108-60-1	3.10E+03	ns	4.70E+04	ns					7.10E+02	n	
				3.00E-03	P					1	0.1		Bis(2-chloroethoxy)methane	111-91-1	1.90E+02	n	2.50E+03	n					5.90E+01	n	
1.10E+00	I	3.30E-04	I					V		1		5.05E+03	Bis(2-chloroethyl)ether	111-44-4	2.30E+01	c	1.00E+02	c	8.50E-01	c	3.70E+00	c	1.40E+00	c	
2.20E+02	I	6.20E-02	I					V		1		4.22E+03	Bis(chloromethyl)ether	542-88-1	8.30E-03	c	3.60E-02	c	4.50E-03	c	2.00E-02	c	7.20E-03	c	
				5.00E-02	I					1	0.1		Bisphenol A	80-05-7	3.20E+03	n	4.10E+04	n					7.70E+02	n	

RMLS TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	GIABS	ABS_d	C_{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
				2.00E-01	I	2.00E-02	H			1			Boron And Borates Only	7440-42-8	1.60E+04	n	2.30E+05	nm	2.10E+01	n	8.80E+01	n	4.00E+03	n	
				2.00E+00	P	2.00E-02	P	V		1			Boron Trichloride	10294-34-5	1.60E+05	nm	2.30E+06	nm	2.10E+01	n	8.80E+01	n	4.20E+01	n	
				4.00E-02	C	1.30E-02	C	V		1			Boron Trifluoride	7637-07-2	3.10E+03	n	4.70E+04	n	1.40E+01	n	5.70E+01	n	2.60E+01	n	
7.00E-01	I			4.00E-03	I					1			Bromate	15541-45-4	9.90E+01	c**	4.70E+02	c*					1.10E+01	c**	1.00E+01
				1.00E-04	X	6.00E-05	X	V		1		2.38E+03	Bromo-2-chloroethane, 1-	107-04-0	3.50E-01	n	1.50E+00	n	6.30E-02	n	2.60E-01	n	1.20E-01	n	
				3.00E-04	X			V		1		8.96E+02	Bromo-3-fluorobenzene, 1-	1073-06-9	2.30E+01	n	3.50E+02	n					4.90E+00	n	
				3.00E-04	X			V		1		3.23E+02	Bromo-4-fluorobenzene, 1-	460-00-4	2.30E+01	n	3.50E+02	ns					4.60E+00	n	
										1	0.1		Bromoacetic acid	79-08-3											6.0E+01(G)
				8.00E-03	I	6.00E-02	I	V		1		6.79E+02	Bromobenzene	108-86-1	2.90E+02	n	1.80E+03	ns	6.30E+01	n	2.60E+02	n	6.20E+01	n	
						4.00E-02	X	V		1		4.04E+03	Bromochloromethane	74-97-5	1.50E+02	n	6.30E+02	n	4.20E+01	n	1.80E+02	n	8.30E+01	n	
6.20E-02	I	3.70E-05	C	8.00E-03	P			V		1		9.32E+02	Bromodichloromethane	75-27-4	2.90E+01	c*	1.30E+02	c*	7.60E+00	c	3.30E+01	c	1.30E+01	c*	8.0E+01(G)
7.90E-03	I	1.10E-06	I	2.00E-02	I			V		1		9.15E+02	Bromoform	75-25-2	1.60E+03	ns	8.60E+03	c**s	2.60E+02	c	1.10E+03	c	3.30E+02	c**	8.0E+01(G)
				1.40E-03	I	5.00E-03	I	V		1		3.59E+03	Bromomethane	74-83-9	6.80E+00	n	3.00E+01	n	5.20E+00	n	2.20E+01	n	7.50E+00	n	
				5.00E-03	H			V		1			Bromophos	2104-96-3	3.90E+02	n	5.80E+03	n					3.50E+01	n	
						1.00E-01	A	V		1		9.66E+02	Bromopropane, 1-	106-94-5	2.20E+02	n	9.40E+02	n	1.00E+02	n	4.40E+02	n	2.10E+02	n	
1.00E-01	O			1.50E-02	O					1	0.1		Bromoxynil	1689-84-5	5.30E+02	c**	2.20E+03	c**					6.10E+01	c**	
1.00E-01	O			1.50E-02	O			V		1			Bromoxynil Octanoate	1689-99-2	6.70E+02	c**	3.20E+03	c**					2.40E+01	c**	
6.00E-01	C	3.00E-05	I			2.00E-03	I	V		1		6.67E+02	Butadiene, 1,3-	106-99-0	1.80E+00	n	7.60E+00	n	2.10E+00	n	8.80E+00	n	4.20E+00	n	
				1.00E-01	I			V		1		7.64E+03	Butanol, N-	71-36-3	7.80E+03	ns	1.20E+05	nms					2.00E+03	n	
5.00E-04	I			4.00E-01	I	5.00E+00	I	V		1			Butyl Alcohol, t-	75-65-0	2.60E+04	n	2.70E+05	nm	5.20E+03	n	2.20E+04	n	4.50E+03	n	
				2.00E+00	P	3.00E+01	P	V		1		2.13E+04	Butyl alcohol, sec-	78-92-2	1.30E+05	nms	1.50E+06	nms	3.10E+04	n	1.30E+05	n	2.40E+04	n	
				5.00E-02	I			V		1			Butylate	2008-41-5	3.90E+03	n	5.80E+04	n					4.60E+02	n	
2.00E-04	C	5.70E-08	C							1	0.1		Butylated hydroxyanisole	25013-16-5	2.70E+05	cm	1.10E+06	cm	4.90E+03	c	2.20E+04	c	1.50E+04	c	
3.60E-03	P			3.00E-01	P					1	0.1		Butylated hydroxytoluene	128-37-0	1.50E+04	c**	6.40E+04	c**					3.40E+02	c**	
				5.00E-02	P			V		1		1.08E+02	Butylbenzene, n-	104-51-8	3.90E+03	ns	5.80E+04	ns					1.00E+03	n	
				1.00E-01	X			V		1		1.45E+02	Butylbenzene, sec-	135-98-8	7.80E+03	ns	1.20E+05	nms					2.00E+03	n	
				1.00E-01	X			V		1		1.83E+02	Butylbenzene, tert-	98-06-6	7.80E+03	ns	1.20E+05	nms					6.90E+02	n	
				2.00E-02	A					1	0.1		Cacodylic Acid	75-60-5	1.30E+03	n	1.60E+04	n					4.00E+02	n	
		1.80E-03	I	1.00E-04	A	1.00E-05	A			0.025	0.001		Cadmium (Diet)	7440-43-9	7.10E+00	n	1.00E+02	n	1.00E-02	n	4.40E-02	n			
		1.80E-03	I	1.00E-04	A	1.00E-05	A			0.05	0.001		Cadmium (Water)	7440-43-9					1.00E-02	n	4.40E-02	n	1.80E+00	n	5.00E+00
				5.00E-01	I	2.20E-03	C			1	0.1		Caprolactam	105-60-2	3.10E+04	n	4.00E+05	nm	2.30E+00	n	9.60E+00	n	9.90E+03	n	
1.50E-01	C	4.30E-05	C	2.00E-03	I					1	0.1		Captafol	2425-06-1	1.30E+02	n	1.50E+03	c**	6.50E+00	c	2.90E+01	c	3.20E+01	n	
2.30E-03	C	6.60E-07	C	1.30E-01	I					1	0.1		Captan	133-06-2	8.20E+03	n	1.00E+05	c**	4.30E+02	c	1.90E+03	c	2.40E+03	n	
				1.00E-01	I					1	0.1		Carbaryl	63-25-2	6.30E+03	n	8.20E+04	n					1.80E+03	n	
				5.00E-03	I					1	0.1		Carbofuran	1563-66-2	3.20E+02	n	4.10E+03	n					9.40E+01	n	4.00E+01
				1.00E-01	I	7.00E-01	I	V		1		7.38E+02	Carbon Disulfide	75-15-0	7.70E+02	ns	3.50E+03	ns	7.30E+02	n	3.10E+03	n	8.10E+02	n	
7.00E-02	I	6.00E-06	I	4.00E-03	I	1.00E-01	I	V		1		4.58E+02	Carbon Tetrachloride	56-23-5	6.50E+01	c**	2.90E+02	c**	4.70E+01	c**	2.00E+02	c**	4.60E+01	c**	5.00E+00
						1.00E-01	P	V		1		5.89E+03	Carbonyl Sulfide	463-58-1	6.70E+01	n	2.80E+02	n	1.00E+02	n	4.40E+02	n	2.10E+02	n	
				1.00E-02	I					1	0.1		Carbosulfan	55285-14-8	6.30E+02	n	8.20E+03	n					5.10E+01	n	
				1.00E-01	I					1	0.1		Carboxin	5234-68-4	6.30E+03	n	8.20E+04	n					1.90E+03	n	
						9.00E-04	I			1			Ceric oxide	1306-38-3	1.30E+06	nm	5.40E+06	nm	9.40E-01	n	3.90E+00	n			
				1.00E-01	I			V		1			Chloral Hydrate	302-17-0	7.80E+03	n	1.20E+05	nm					2.00E+03	n	
				1.50E-02	I					1	0.1		Chloramben	133-90-4	9.50E+02	n	1.20E+04	n					2.90E+02	n	
										1	0.1		Chloramines, Organic	E701235											4.0E+03(G)
4.00E-01	H									1	0.1		Chloranil	118-75-2	1.30E+02	c	5.70E+02	c					1.80E+01	c	
				5.00E-04	G			V		1	0.04		Chlordane (alpha)	5103-71-9	3.60E+01	n	5.00E+02	n					3.60E+00	n	
				5.00E-04	G			V		1	0.04		Chlordane (gamma)	5103-74-2	3.60E+01	n	5.00E+02	n					1.00E+01	n	
3.50E-01	I	1.00E-04	I	5.00E-04	I	7.00E-04	I	V		1	0.04		Chlordane (technical mixture)	12789-03-6	3.50E+01	n	4.50E+02	n	7.30E-01	n	3.10E+00	n	7.40E-01	n	2.00E+00
1.00E+01	I	4.60E-03	C	3.00E-04	I					1	0.1		Chlordecone (Kepone)	143-50-0	5.40E+00	c**	2.30E+01	c*	6.10E-02	c	2.70E-01	c	3.50E-01	c**	
				7.00E-04	A					1	0.1		Chlorfenvinphos	470-90-6	4.40E+01	n	5.70E+02	n					1.10E+01	n	
				9.00E-02	O					1	0.1		Chlorimuron, Ethyl-	90982-32-4	5.70E+03	n	7.40E+04	n					1.80E+03	n	
				1.00E-01	I	1.50E-04	A	V		1		2.78E+03	Chlorine	7782-50-5	1.80E-01	n	7.80E-01	n	1.50E-01	n	6.40E-01	n	3.00E-01	n	4.0E+03(G)
				3.00E-02	I	2.00E-04	I	V		1			Chlorine Dioxide	10049-04-4	2.30E+03	n	3.40E+04	n	2.10E-01	n	8.80E-01	n	4.20E-01	n	8.0E+02(G)
				3.00E-02	I					1			Chlorite (Sodium Salt)	7758-19-2	2.30E+03	n	3.50E+04	n					6.00E+02	n	1.00E+03
						5.00E+01	I	V		1		1.15E+03	Chloro-1,1-difluoroethane, 1-	75-68-3	5.40E+04	ns	2.30E+05	nms	5.20E+04	n	2.20E+05	n	1.00E+05	n	
		3.00E-04	I	2.00E-02	H	2.00E-02	I	V		1		7.86E+02	Chloro-1,3-butadiene, 2-	126-99-8	1.00E+00	c*	4.40E+00	c*	9.40E-01	c*	4.10E+00	c*	1.90E+00	c*	
4.60E-01	H									1	0.1		Chloro-2-methylaniline HCl, 4-	3165-93-3	1.20E+02	c	5.00E+02	c					1.70E+01	c	
1.00E-01	P	7.70E-05	C	3.00E-03	X					1	0.1		Chloro-2-methylaniline, 4-	95-69-2	1.90E+02	n	2.30E+03	c**	3.60E+00	c	1.60E+01	c	5.40E+01	n	
2.70E-01	X							V		1		1.18E+04	Chloroacetaldehyde, 2-	107-20-0	2.60E+02	c	1.20E+03	c					2.90E+01	c	
										1	0.1		Chloroacetic Acid	79-11-8											6.0E+01(G)
						3.00E-05	I			1	0.1		Chloroacetophenone, 2-	532-27-4	4.30E+04	n	1.80E+05	nm	3.10E-02	n	1.30E-01	n			
2.00E-01	P			5.00E-04	P					1	0.1		Chloroaniline, p-	106-47-8	3.20E+01	n	4.10E+02	n					9.50E+00	n	
				2.00E-02	I	5.00E-02	P	V		1		7.61E+02	Chlorobenzene	108-90-7	2.80E+02	n	1.30E+03	ns	5.20E+01	n	2.20E+02	n	7.80E+01	n	1.00E+02
				1.00E-01	X					1	0.1		Chlorobenzene sulfonic acid, p-	98-66-8	6.30E+03	n	8.20E+04	n					2.00E+03	n	
1.10E-01	C	3.10E-05	C	2.00E-02	I					1	0.1		Chlorobenzilate	510-15-6	4.90E+02	c**	2.10E+03	c**	9.10E+00	c	4.00E+01	c	3.10E+01	c**	
				3.00E-02	X					1	0.1		Chlorobenzoic Acid, p-	74-11-3	1.90E+03	n	2.50E+04	n					5.10E+02	n	
		8.60E-06	C	3.00E-03	P	3.00E-01	P	V		1		2.90E+02	Chlorobenzotrifluoride, 4-	98-56-6	2.10E+02	n	9.60E+02	c**s	3.30E+01	c**	1.40E+02	c**	3.50E+01	n	
				4.00E-02	P			V		1		7.28E+02	Chlorobutane, 1-	109-69-3	3.10E+03	ns	4.70E+04	ns					6.40E+02	n	
						5.00E+01	I	V		1		1.68E+03	Chlorodifluoromethane	75-45-6	4.90E+04	ns	2.10E+05	nms	5.20E+04	n	2.20E+05	n	1.00E+05	n	
				2.00E-02	P			V		1		1.11E+05	Chloroethanol, 2-	107-07-3	1.60E+03	n	2.30E+04	n					4.00E+02	n	
3.10E-02	C	2.30E-05	I	1.00E-02	I	9.80E-02	A	V		1		2.54E+03	Chloroform	67-66-3	3.20E+01	c**	1.40E+02	c**	1.20E+01	c**	5.30E+01	c**	2.20E+01	c**	8.0E+01(G)

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	GIABS	ABS_d	C_{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
						9.00E-02	I	V		1		1.32E+03	Chloromethane	74-87-3	1.10E+02	n	4.60E+02	n	9.40E+01	n	3.90E+02	n	1.90E+02	n	
2.40E+00	C	6.90E-04	C					V		1		9.32E+03	Chloromethyl Methyl Ether	107-30-2	2.00E+00	c	8.90E+00	c	4.10E-01	c	1.80E+00	c	6.50E-01	c	
3.00E-01	P			3.00E-03	P	1.00E-05	X			1	0.1		Chloronitrobenzene, o-	88-73-3	1.80E+02	c**	7.70E+02	c**	1.00E-02	n	4.40E-02	n	2.40E+01	c**	
6.00E-02	P			7.00E-04	P	2.00E-03	P			1	0.1		Chloronitrobenzene, p-	100-00-5	4.40E+01	n	5.70E+02	n	2.10E+00	n	8.80E+00	n	1.30E+01	n	
				5.00E-03	I			V		1		2.74E+04	Chlorophenol, 2-	95-57-8	3.90E+02	n	5.80E+03	n					9.10E+01	n	
						4.00E-04	C	V		1		6.17E+02	Chloropicrin	76-06-2	2.00E+00	n	8.20E+00	n	4.20E-01	n	1.80E+00	n	8.30E-01	n	
1.70E-02	C			1.50E-02	I					1	0.1		Chlorothalonil	1897-45-6	9.50E+02	n	1.20E+04	n					2.60E+02	n	
				2.00E-02	I			V		1		9.07E+02	Chlorotoluene, o-	95-49-8	1.60E+03	ns	2.30E+04	ns					2.40E+02	n	
				2.00E-02	X			V		1		2.53E+02	Chlorotoluene, p-	106-43-4	1.60E+03	ns	2.30E+04	ns					2.50E+02	n	
2.40E+02	C	6.90E-02	C							1	0.1		Chlorozotocin	54749-90-5	2.30E-01	c	9.60E-01	c	4.10E-03	c	1.80E-02	c	3.20E-02	c	
				5.00E-03	O					1	0.1		Chlorpropham	101-21-3	3.20E+02	n	4.10E+03	n					7.10E+01	n	
				1.00E-03	A					1	0.1		Chlorpyrifos	2921-88-2	6.30E+01	n	8.20E+02	n					8.40E+00	n	
				1.00E-02	H					1	0.1		Chlorpyrifos Methyl	5598-13-0	6.30E+02	n	8.20E+03	n					1.20E+02	n	
				5.00E-02	O					1	0.1		Chlorsulfuron	64902-72-3	3.20E+03	n	4.10E+04	n					9.90E+02	n	
				1.00E-02	I					1	0.1		Chlorthal-dimethyl	1861-32-1	6.30E+02	n	8.20E+03	n					1.20E+02	n	
				8.00E-04	H					1	0.1		Chlorthiophos	60238-56-4	5.10E+01	n	6.60E+02	n					2.80E+00	n	
				1.50E+00	I					0.013			Chromium(III), Insoluble Salts	16065-83-1	1.20E+05	nm	1.80E+06	nm					2.20E+04	n	
5.00E-01	C	8.40E-02	G	3.00E-03	I	1.00E-04	I		M	0.025			Chromium(VI)	18540-29-9	3.00E+01	c**	6.30E+02	c**	1.20E-03	c*	1.50E-02	c*	3.50E+00	c*	
										0.013			Chromium, Total	7440-47-3											1.00E+02
				1.30E-02	I					1	0.1		Clofentezine	74115-24-5	8.20E+02	n	1.10E+04	n					2.30E+02	n	
		9.00E-03	P	3.00E-04	P	6.00E-06	P			1			Cobalt	7440-48-4	2.30E+01	n	3.50E+02	n	6.30E-03	n	2.60E-02	n	6.00E+00	n	
		6.20E-04	I					V	M	1			Coke Oven Emissions	E649830					1.60E-01	c	2.00E+00	c			
				4.00E-02	H					1			Copper	7440-50-8	3.10E+03	n	4.70E+04	n					8.00E+02	n	1.30E+03
				5.00E-02	I	6.00E-01	C			1	0.1		Cresol, m-	108-39-4	3.20E+03	n	4.10E+04	n	6.30E+02	n	2.60E+03	n	9.30E+02	n	
				5.00E-02	I	6.00E-01	C			1	0.1		Cresol, o-	95-48-7	3.20E+03	n	4.10E+04	n	6.30E+02	n	2.60E+03	n	9.30E+02	n	
				2.00E-02	P	6.00E-01	C			1	0.1		Cresol, p-	106-44-5	1.30E+03	n	1.60E+04	n	6.30E+02	n	2.60E+03	n	3.70E+02	n	
				1.00E-01	A					1	0.1		Cresol, p-chloro-m-	59-50-7	6.30E+03	n	8.20E+04	n					1.40E+03	n	
				1.00E-01	A	6.00E-01	C			1	0.1		Cresols	1319-77-3	6.30E+03	n	8.20E+04	n	6.30E+02	n	2.60E+03	n	1.50E+03	n	
1.90E+00	H			1.00E-03	P			V		1		1.66E+04	Crotonaldehyde, trans-	123-73-9	3.70E+01	c**	1.70E+02	c**					4.00E+00	c**	
				1.00E-01	I	4.00E-01	I	V		1		2.68E+02	Cumene	98-82-8	1.90E+03	ns	9.90E+03	ns	4.20E+02	n	1.80E+03	n	4.50E+02	n	
2.20E-01	C	6.30E-05	C							1	0.1		Cupferron	135-20-6	2.50E+02	c	1.00E+03	c	4.50E+00	c	1.90E+01	c	3.50E+01	c	
8.40E-01	H			2.00E-03	H					1	0.1		Cyanazine	21725-46-2	6.50E+01	c**	2.70E+02	c**					8.80E+00	c**	
													Cyanides												
				1.00E-03	I					1			~Calcium Cyanide	592-01-8	7.80E+01	n	1.20E+03	n					2.00E+01	n	
				5.00E-03	I					1			~Copper Cyanide	544-92-3	3.90E+02	n	5.80E+03	n					1.00E+02	n	
				6.00E-04	I	8.00E-04	G	V		1		9.54E+05	~Cyanide (CN-)	57-12-5	2.30E+01	n	1.50E+02	n	8.30E-01	n	3.50E+00	n	1.50E+00	n	2.00E+02
				1.00E-03	I			V		1			~Cyanogen	460-19-5	7.80E+01	n	1.20E+03	n					2.00E+01	n	
				9.00E-02	I			V		1			~Cyanogen Bromide	506-68-3	7.00E+03	n	1.10E+05	nm					1.80E+03	n	
				5.00E-02	I			V		1			~Cyanogen Chloride	506-77-4	3.90E+03	n	5.80E+04	n					1.00E+03	n	
				6.00E-04	I	8.00E-04	I	V		1		1.00E+07	~Hydrogen Cyanide	74-90-8	2.30E+01	n	1.50E+02	n	8.30E-01	n	3.50E+00	n	1.50E+00	n	
				2.00E-03	I					1			~Potassium Cyanide	151-50-8	1.60E+02	n	2.30E+03	n					4.00E+01	n	
				5.00E-03	I					0.04			~Potassium Silver Cyanide	506-61-6	3.90E+02	n	5.80E+03	n					8.20E+01	n	
				1.00E-01	I					0.04			~Silver Cyanide	506-64-9	7.80E+03	n	1.20E+05	nm					1.80E+03	n	
				1.00E-03	I					1			~Sodium Cyanide	143-33-9	7.80E+01	n	1.20E+03	n					2.00E+01	n	2.00E+02
				2.00E-04	P					1			~Thiocyanates	E1790665	1.60E+01	n	2.30E+02	n					4.00E+00	n	
				2.00E-04	X			V		1			~Thiocyanic Acid	463-56-9	1.60E+01	n	2.30E+02	n					4.00E+00	n	
				5.00E-02	I					1			~Zinc Cyanide	557-21-1	3.90E+03	n	5.80E+04	n					1.00E+03	n	
						6.00E+00	I	V		1		1.17E+02	Cyclohexane	110-82-7	6.50E+03	ns	2.70E+04	ns	6.30E+03	n	2.60E+04	n	1.30E+04	n	
2.00E-02	X			2.00E-02	X					1	0.1		Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	1.30E+03	n	1.10E+04	c**					2.80E+02	c**	
				5.00E+00	I	7.00E-01	P	V		1		5.11E+03	Cyclohexanone	108-94-1	2.80E+04	ns	1.30E+05	nms	7.30E+02	n	3.10E+03	n	1.40E+03	n	
				5.00E-03	P	1.00E+00	X	V		1		2.83E+02	Cyclohexene	110-83-8	3.10E+02	ns	3.10E+03	ns	1.00E+03	n	4.40E+03	n	7.00E+01	n	
				2.00E-01	I			V		1		2.93E+05	Cyclohexylamine	108-91-8	1.60E+04	n	2.30E+05	nm					3.80E+03	n	
				2.50E-02	I					1	0.1		Cyfluthrin	68359-37-5	1.60E+03	n	2.10E+04	n					1.20E+02	n	
				7.20E-02	O					1	0.1		Cypermethrin	52315-07-8	4.50E+03	n	5.90E+04	n					1.40E+03	n	
				5.00E-01	O					1	0.1		Cyromazine	66215-27-8	3.20E+04	n	4.10E+05	nm					9.90E+03	n	
2.40E-01	I	6.90E-05	C	3.00E-05	X					1	0.1		DDD, p,p`- (DDD)	72-54-8	1.90E+00	n	2.50E+01	n	4.10E+00	c	1.80E+01	c	6.30E-02	n	
3.40E-01	I	9.70E-05	C	3.00E-04	X			V		1			DDE, p,p'-	72-55-9	2.30E+01	n	3.50E+02	n	2.90E+00	c	1.30E+01	c	4.60E+00	c**	
3.40E-01	I	9.70E-05	I	5.00E-04	I					1	0.03		DDT	50-29-3	3.70E+01	n	5.20E+02	n	2.90E+00	c	1.30E+01	c	1.00E+01	n	
				3.00E-02	I					1	0.1		Dalapon	75-99-0	1.90E+03	n	2.50E+04	n					6.00E+02	n	2.00E+02
1.80E-02	C	5.10E-06	C	1.50E-01	I					1	0.1		Daminozide	1596-84-5	3.00E+03	c**	1.30E+04	c**	5.50E+01	c	2.40E+02	c	4.30E+02	c**	
7.00E-04	I			7.00E-03	I					1	0.1		Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5	4.40E+02	n	5.70E+03	n					1.40E+02	n	
				4.00E-05	I					1	0.1		Demeton	8065-48-3	2.50E+00	n	3.30E+01	n					4.20E-01	n	
1.20E-03	I			6.00E-01	I					1	0.1		Di(2-ethylhexyl)adipate	103-23-1	3.80E+04	n	1.90E+05	c**m					6.50E+03	c**	4.00E+02
6.10E-02	H									1	0.1		Diallate	2303-16-4	8.90E+02	c	3.80E+03	c					5.40E+01	c	
				7.00E-04	A					1	0.1		Diazinon	333-41-5	4.40E+01	n	5.70E+02	n					1.00E+01	n	
				1.00E-02	X			V		1			Dibenzothiophene	132-65-0	7.80E+02	n	1.20E+04	n					6.50E+01	n	
8.00E-01	P	6.00E-03	P	2.00E-04	P	2.00E-04	I	V	M	1		9.79E+02	Dibromo-3-chloropropane, 1,2-	96-12-8	5.30E-01	c**	6.40E+00	c**	1.70E-02	c*	2.00E-01	c**	3.30E-02	c*	2.00E-01
										1	0.1		Dibromoacetic acid	631-64-1											6.0E+01(G)
				4.00E-04	X			V		1		1.59E+02	Dibromobenzene, 1,3-	108-36-1	3.10E+01	n	4.70E+02	ns					5.30E+00	n	
				1.00E-02	I			V		1			Dibromobenzene, 1,4-	106-37-6	7.80E+02	n	1.20E+04	n					1.30E+02	n	
8.40E-02	I			2.00E-02	I			V		1		8.02E+02	Dibromochloromethane	124-48-1	8.30E+02	c**s	3.90E+03	c**s					8.70E+01	c**	8.0E+01(G)

RMLs TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	GIABS	ABS_d	C_{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
2.00E+00	I	6.00E-04	I	9.00E-03	I	9.00E-03	I	V		1		1.34E+03	Dibromoethane, 1,2-	106-93-4	3.60E+00	c*	1.60E+01	c*	4.70E-01	c*	2.00E+00	c*	7.50E-01	c*	5.00E-02
						4.00E-03	X	V		1		2.82E+03	Dibromomethane (Methylene Bromide)	74-95-3	2.40E+01	n	9.90E+01	n	4.20E+00	n	1.80E+01	n	8.30E+00	n	
				3.00E-04	P					1	0.1		Dibutyltin Compounds	E1790661	1.90E+01	n	2.50E+02	n					6.00E+00	n	
				3.00E-02	I					1	0.1		Dicamba	1918-00-9	1.90E+03	n	2.50E+04	n					5.70E+02	n	
										1	0.1		Dichloramine	3400-09-7											4.0E+03(G)
		4.20E-03	P					V		1		5.54E+02	Dichloro-2-butene, 1,4-	764-41-0	2.10E-01	c	9.40E-01	c	6.70E-02	c	2.90E-01	c	1.30E-01	c	
		4.20E-03	P					V		1		5.19E+02	Dichloro-2-butene, cis-1,4-	1476-11-5	7.40E-01	c	3.20E+00	c	6.70E-02	c	2.90E-01	c	1.30E-01	c	
		4.20E-03	P					V		1		7.60E+02	Dichloro-2-butene, trans-1,4-	110-57-6	7.40E-01	c	3.20E+00	c	6.70E-02	c	2.90E-01	c	1.30E-01	c	
5.00E-02	I			4.00E-03	I					1	0.1		Dichloroacetic Acid	79-43-6	2.50E+02	n	3.30E+03	n					7.90E+01	n	6.0E+01(G)
				9.00E-02	I	2.00E-01	H	V		1		3.76E+02	Dichlorobenzene, 1,2-	95-50-1	1.80E+03	ns	9.30E+03	ns	2.10E+02	n	8.80E+02	n	3.00E+02	n	6.00E+02
5.40E-03	C	1.10E-05	C	7.00E-02	A	8.00E-01	I	V		1			Dichlorobenzene, 1,4-	106-46-7	2.60E+02	c*	1.10E+03	c*	2.60E+01	c*	1.10E+02	c*	4.80E+01	c*	7.50E+01
4.50E-01	I	3.40E-04	C							1	0.1		Dichlorobenzidine, 3,3'-	91-94-1	1.20E+02	c	5.10E+02	c	8.30E-01	c	3.60E+00	c	1.30E+01	c	
				9.00E-03	X					1	0.1		Dichlorobenzophenone, 4,4'-	90-98-2	5.70E+02	n	7.40E+03	n					7.80E+01	n	
				2.00E-01	I	1.00E-01	X	V		1		8.45E+02	Dichlorodifluoromethane	75-71-8	8.70E+01	n	3.70E+02	n	1.00E+02	n	4.40E+02	n	2.00E+02	n	
5.70E-03	C	1.60E-06	C	2.00E-01	P			V		1		1.69E+03	Dichloroethane, 1,1-	75-34-3	3.60E+02	c*	1.60E+03	c	1.80E+02	c	7.70E+02	c	2.80E+02	c*	
9.10E-02	I	2.60E-05	I	6.00E-03	X	7.00E-03	P	V		1		2.98E+03	Dichloroethane, 1,2-	107-06-2	3.10E+01	n	1.40E+02	n	7.30E+00	n	3.10E+01	n	1.30E+01	n	5.00E+00
				5.00E-02	I	2.00E-01	I	V		1		1.19E+03	Dichloroethylene, 1,1-	75-35-4	2.30E+02	n	1.00E+03	n	2.10E+02	n	8.80E+02	n	2.80E+02	n	7.00E+00
				2.00E-03	I			V		1		2.37E+03	Dichloroethylene, cis-1,2-	156-59-2	1.60E+02	n	2.30E+03	n					3.60E+01	n	7.00E+01
				2.00E-02	I	4.00E-02	X	V		1		1.85E+03	Dichloroethylene, trans-1,2-	156-60-5	7.00E+01	n	3.00E+02	n	4.20E+01	n	1.80E+02	n	6.80E+01	n	1.00E+02
				3.00E-03	I					1	0.1		Dichlorophenol, 2,4-	120-83-2	1.90E+02	n	2.50E+03	n					4.60E+01	n	
				1.00E-02	I					1	0.05		Dichlorophenoxy Acetic Acid, 2,4-	94-75-7	7.00E+02	n	9.60E+03	n					1.70E+02	n	7.00E+01
3.70E-02	P	3.70E-06	P	4.00E-02	P	4.00E-03	I	V		1		1.36E+03	Dichloropropane, 1,2-	78-87-5	1.60E+01	n	6.60E+01	n	4.20E+00	n	1.80E+01	n	8.20E+00	n	5.00E+00
				2.00E-02	P			V		1		1.49E+03	Dichloropropane, 1,3-	142-28-9	1.60E+03	ns	2.30E+04	ns					3.70E+02	n	
				3.00E-03	I					1	0.1		Dichloropropanol, 2,3-	616-23-9	1.90E+02	n	2.50E+03	n					5.90E+01	n	
1.00E-01	I	4.00E-06	I	3.00E-02	I	2.00E-02	I	V		1		1.57E+03	Dichloropropene, 1,3-	542-75-6	7.20E+01	n	3.10E+02	n	2.10E+01	n	8.80E+01	n	3.90E+01	n	
2.90E-01	I	8.30E-05	C	5.00E-04	I	5.00E-04	I			1	0.1		Dichlorvos	62-73-7	3.20E+01	n	4.10E+02	n	5.20E-01	n	2.20E+00	n	9.90E+00	n	
				3.00E-05	O					1	0.1		Dicrotophos	141-66-2	1.90E+00	n	2.50E+01	n					6.00E-01	n	
				8.00E-02	P	3.00E-04	X	V		1		2.56E+02	Dicyclopentadiene	77-73-6	1.30E+00	n	5.40E+00	n	3.10E-01	n	1.30E+00	n	6.30E-01	n	
1.60E+01	I	4.60E-03	I	5.00E-05	I					1	0.1		Dieldrin	60-57-1	3.20E+00	n	1.40E+01	c**	6.10E-02	c	2.70E-01	c	1.80E-01	c**	
		3.00E-04	C			5.00E-03	I			1	0.1		Diesel Engine Exhaust	E17136615					9.40E-01	c**	4.10E+00	c**			
				2.00E-03	P	2.00E-04	P			1	0.1		Diethanolamine	111-42-2	1.30E+02	n	1.60E+03	n	2.10E-01	n	8.80E-01	n	4.00E+01	n	
				3.00E-02	P	1.00E-04	P			1	0.1		Diethylene Glycol Monobutyl Ether	112-34-5	1.90E+03	n	2.40E+04	n	1.00E-01	n	4.40E-01	n	6.00E+02	n	
				6.00E-02	P	3.00E-04	P			1	0.1		Diethylene Glycol Monoethyl Ether	111-90-0	3.80E+03	n	4.80E+04	n	3.10E-01	n	1.30E+00	n	1.20E+03	n	
				1.00E-03	P			V		1		1.12E+05	Diethylformamide	617-84-5	7.80E+01	n	1.20E+03	n					2.00E+01	n	
3.50E+02	C	1.00E-01	C							1	0.1		Diethylstilbestrol	56-53-1	1.60E-01	c	6.60E-01	c	2.80E-03	c	1.20E-02	c	5.10E-03	c	
				8.30E-02	O					1	0.1		Difenzoquat	43222-48-6	5.20E+03	n	6.80E+04	n					1.70E+03	n	
				2.00E-02	I					1	0.1		Diflubenzuron	35367-38-5	1.30E+03	n	1.60E+04	n					2.90E+02	n	
						4.00E+01	I	V		1		1.43E+03	Difluoroethane, 1,1-	75-37-6	4.80E+04	ns	2.00E+05	nms	4.20E+04	n	1.80E+05	n	8.30E+04	n	
						3.00E+01	X	V		1		6.91E+02	Difluoropropane, 2,2-	420-45-1	2.40E+04	ns	1.00E+05	ns	3.10E+04	n	1.30E+05	n	6.30E+04	n	
4.40E-02	C	1.30E-05	C					V		1			Dihydrosafrole	94-58-6	9.90E+02	c	4.50E+03	c	2.20E+01	c	9.40E+01	c	3.00E+01	c	
						7.00E-01	P	V		1		2.26E+03	Diisopropyl Ether	108-20-3	2.20E+03	n	9.40E+03	ns	7.30E+02	n	3.10E+03	n	1.50E+03	n	
				8.00E-02	I			V		1		5.30E+02	Diisopropyl Methylphosphonate	1445-75-6	6.30E+03	ns	9.30E+04	ns					1.60E+03	n	
				2.20E-02	O					1	0.1		Dimethipin	55290-64-7	1.40E+03	n	1.80E+04	n					4.40E+02	n	
				2.20E-03	O					1	0.1		Dimethoate	60-51-5	1.40E+02	n	1.80E+03	n					4.40E+01	n	
1.60E+00	P									1	0.1		Dimethoxybenzidine, 3,3'-	119-90-4	3.40E+01	c	1.40E+02	c					4.70E+00	c	
1.70E-03	P			6.00E-02	P					1	0.1		Dimethyl methylphosphonate	756-79-6	3.80E+03	n	4.90E+04	n					1.20E+03	n	
4.60E+00	C	1.30E-03	C							1	0.1		Dimethylamino azobenzene [p-]	60-11-7	1.20E+01	c	5.00E+01	c	2.20E-01	c	9.40E-01	c	5.00E-01	c	
5.80E-01	H									1	0.1		Dimethylaniline HCl, 2,4-	21436-96-4	9.40E+01	c	4.00E+02	c					1.30E+01	c	
2.00E-01	P			2.00E-03	X					1	0.1		Dimethylaniline, 2,4-	95-68-1	1.30E+02	n	1.10E+03	c**					3.70E+01	c**	
2.70E-02	P			2.00E-03	I			V		1		8.30E+02	Dimethylaniline, N,N-	121-69-7	1.60E+02	n	2.30E+03	ns					3.50E+01	n	
1.10E+01	P									1	0.1		Dimethylbenzidine, 3,3'-	119-93-7	4.90E+00	c	2.10E+01	c					6.50E-01	c	
				1.00E-01	P	3.00E-02	I	V		1		1.06E+05	Dimethylformamide	68-12-2	2.60E+03	n	1.50E+04	n	3.10E+01	n	1.30E+02	n	6.10E+01	n	
				1.00E-04	X	2.00E-06	X	V		1		1.72E+05	Dimethylhydrazine, 1,1-	57-14-7	5.70E-02	n	2.40E-01	n	2.10E-03	n	8.80E-03	n	4.20E-03	n	
5.50E+02	C	1.60E-01	C					V		1		1.89E+05	Dimethylhydrazine, 1,2-	540-73-8	8.80E-02	c	4.10E-01	c	1.80E-03	c	7.70E-03	c	2.80E-03	c	
				2.00E-02	I					1	0.1		Dimethylphenol, 2,4-	105-67-9	1.30E+03	n	1.60E+04	n					3.60E+02	n	
				6.00E-04	I					1	0.1		Dimethylphenol, 2,6-	576-26-1	3.80E+01	n	4.90E+02	n					1.10E+01	n	
				1.00E-03	I					1	0.1		Dimethylphenol, 3,4-	95-65-8	6.30E+01	n	8.20E+02	n					1.80E+01	n	
4.50E-02	C	1.30E-05	C					V		1		4.73E+02	Dimethylvinylchloride	513-37-1	1.10E+02	c	4.80E+02	cs	2.20E+01	c	9.40E+01	c	3.30E+01	c	
				8.00E-05	X					1	0.1		Dinitro-o-cresol, 4,6-	534-52-1	5.10E+00	n	6.60E+01	n					1.50E+00	n	
				2.00E-03	I					1	0.1		Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5	1.30E+02	n	1.60E+03	n					2.30E+01	n	
				4.00E-04	X	2.00E-03	X			1	0.1		Dinitroaniline, 3,5-	618-87-1	2.50E+01	n	3.30E+02	n	2.10E+00	n	8.80E+00	n	7.70E+00	n	
				1.00E-04	P					1	0.1		Dinitrobenzene, 1,2-	528-29-0	6.30E+00	n	8.20E+01	n					1.90E+00	n	
				1.00E-04	I					1	0.1		Dinitrobenzene, 1,3-	99-65-0	6.30E+00	n	8.20E+01	n					2.00E+00	n	
				1.00E-04	P					1	0.1		Dinitrobenzene, 1,4-	100-25-4	6.30E+00	n	8.20E+01	n					2.00E+00	n	
				2.00E-03	I					1	0.1		Dinitrophenol, 2,4-	51-28-5	1.30E+02	n	1.60E+03	n					3.90E+01	n	
6.80E-01	I									1	0.1		Dinitrotoluene Mixture, 2,4/2,6-	E1615210	8.00E+01	c	3.40E+02	c					1.10E+01	c	
3.10E-01	C	8.90E-05	C	2.00E-03	I					1	0.102		Dinitrotoluene, 2,4-	121-14-2	1.30E+02	n	7.40E+02	c**	3.20E+00	c	1.40E+01	c	2.40E+01	c**	
1.50E+00	P			3.00E-04	X					1	0.099		Dinitrotoluene, 2,6-	606-20-2	1.90E+01	n	1.50E+02	c**					4.90E+00	c**	
				1.00E-04	X					1	0.006		Dinitrotoluene, 2-Amino-4,6-	35572-78-2	7.70E+00	n	1.10E+02	n					1.90E+00	n	
				1.00E-04	X					1	0.009		Dinitrotoluene, 4-Amino-2,6-	19406-51-0	7.70E+00	n	1.10E+02	n					1.90E+00	n	
4.50E-01	X			9.00E-04	X					1	0.1		Dinitrotoluene, Technical grade	25321-14-6	5.70E+01	n	5.10E+02	c**					1.00E+01	c**	

RMLs TR=1E-04 THQ=1.0

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Toxicity and Chemical-specific Information														Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	GIABS	ABS_d	C_{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
				1.00E-03	I					1	0.1		Dinoseb	88-85-7	6.30E+01	n	8.20E+02	n					1.50E+01	n	7.00E+00
1.00E-01	I	5.00E-06	I	3.00E-02	I	3.00E-02	I	V		1		1.16E+05	Dioxane, 1,4-	123-91-1	5.30E+02	c**	2.40E+03	c**	3.10E+01	n	1.30E+02	n	4.60E+01	c**	
													Dioxins												
6.20E+03	I	1.30E+00	I							1	0.03		~Hexachlorodibenzo-p-dioxin, Mixture	34465-46-8	1.00E-02	c	4.70E-02	c	2.20E-04	c	9.40E-04	c	1.30E-03	c	
1.30E+05	C	3.80E+01	C	7.00E-10	I	4.00E-08	C	V		1	0.03		~TCDD, 2,3,7,8-	1746-01-6	5.10E-05	n	7.20E-04	n	7.40E-06	c**	3.20E-05	c**	1.20E-05	c**	3.00E-05
				3.00E-02	I					1	0.1		Diphenamid	957-51-7	1.90E+03	n	2.50E+04	n					5.30E+02	n	
						4.00E-04	X	V		1			Diphenyl Ether	101-84-8	3.40E+01	n	1.40E+02	n	4.20E-01	n	1.80E+00	n	8.30E-01	n	
				8.00E-04	X					1	0.1		Diphenyl Sulfone	127-63-9	5.10E+01	n	6.60E+02	n					1.50E+01	n	
				1.00E-01	O					1	0.1		Diphenylamine	122-39-4	6.30E+03	n	8.20E+04	n					1.30E+03	n	
8.00E-01	I	2.20E-04	I							1	0.1		Diphenylhydrazine, 1,2-	122-66-7	6.80E+01	c	2.90E+02	c	1.30E+00	c	5.60E+00	c	7.80E+00	c	
				2.20E-03	I					1	0.1		Diquat	2764-72-9	1.40E+02	n	1.80E+03	n					4.00E+01	n	2.00E+01
7.40E+00	C	2.10E-03	C							1	0.1		Direct Black 38	1937-37-7	7.30E+00	c	3.10E+01	c	1.30E-01	c	5.80E-01	c	1.10E+00	c	
7.40E+00	C	2.10E-03	C							1	0.1		Direct Blue 6	2602-46-2	7.30E+00	c	3.10E+01	c	1.30E-01	c	5.80E-01	c	1.10E+00	c	
6.70E+00	C	1.90E-03	C							1	0.1		Direct Brown 95	16071-86-6	8.10E+00	c	3.40E+01	c	1.50E-01	c	6.50E-01	c	1.20E+00	c	
				4.00E-05	I					1	0.1		Disulfoton	298-04-4	2.50E+00	n	3.30E+01	n					5.00E-01	n	
				1.00E-02	I			V		1			Dithiane, 1,4-	505-29-3	7.80E+02	n	1.20E+04	n					2.00E+02	n	
				2.00E-03	I					1	0.1		Diuron	330-54-1	1.30E+02	n	1.60E+03	n					3.60E+01	n	
				2.00E-02	O					1	0.1		Dodine	2439-10-3	1.30E+03	n	1.60E+04	n					4.00E+02	n	
				5.00E-02	O			V		1			EPTC	759-94-4	3.90E+03	n	5.80E+04	n					7.50E+02	n	
				6.00E-03	I			V		1			Endosulfan	115-29-7	4.70E+02	n	7.00E+03	n					1.00E+02	n	
				6.00E-03	P					1	0.1		Endosulfan Sulfate	1031-07-8	3.80E+02	n	4.90E+03	n					1.10E+02	n	
				2.00E-02	I					1	0.1		Endothall	145-73-3	1.30E+03	n	1.60E+04	n					3.80E+02	n	1.00E+02
				3.00E-04	I					1	0.1		Endrin	72-20-8	1.90E+01	n	2.50E+02	n					2.30E+00	n	2.00E+00
9.90E-03	I	1.20E-06	I	6.00E-03	P	1.00E-03	I	V		1		1.05E+04	Epichlorohydrin	106-89-8	1.90E+01	n	8.20E+01	n	1.00E+00	n	4.40E+00	n	2.00E+00	n	
						2.00E-02	I	V		1		1.53E+04	Epoxybutane, 1,2-	106-88-7	1.60E+02	n	6.70E+02	n	2.10E+01	n	8.80E+01	n	4.20E+01	n	
				4.00E-02	P					1	0.1		Ethanol, 2-(2-methoxyethoxy)-	111-77-3	2.50E+03	n	3.30E+04	n					8.00E+02	n	
				5.00E-03	I					1	0.1		Ethephon	16672-87-0	3.20E+02	n	4.10E+03	n					1.00E+02	n	
				5.00E-04	I					1	0.1		Ethion	563-12-2	3.20E+01	n	4.10E+02	n					4.30E+00	n	
				1.00E-01	P	6.00E-02	P	V		1		2.38E+04	Ethoxyethanol Acetate, 2-	111-15-9	2.60E+03	n	1.40E+04	n	6.30E+01	n	2.60E+02	n	1.20E+02	n	
				9.00E-02	P	4.00E-02	P	V		1		1.06E+05	Ethoxyethanol, 2-	110-80-5	2.60E+03	n	1.50E+04	n	4.20E+01	n	1.80E+02	n	8.00E+01	n	
				7.00E-01	P	7.00E-02	P	V		1		1.08E+04	Ethyl Acetate	141-78-6	6.20E+02	n	2.60E+03	n	7.30E+01	n	3.10E+02	n	1.40E+02	n	
				5.00E-03	P	8.00E-03	P	V		1		2.50E+03	Ethyl Acrylate	140-88-5	4.70E+01	n	2.10E+02	n	8.30E+00	n	3.50E+01	n	1.40E+01	n	
						4.00E+00	P	V		1		2.12E+03	Ethyl Chloride (Chloroethane)	75-00-3	5.40E+03	ns	2.30E+04	ns	4.20E+03	n	1.80E+04	n	8.30E+03	n	
				2.00E-01	I			V		1		1.01E+04	Ethyl Ether	60-29-7	1.60E+04	ns	2.30E+05	nms					3.90E+03	n	
						3.00E-01	P	V		1		1.10E+03	Ethyl Methacrylate	97-63-2	1.80E+03	ns	7.60E+03	ns	3.10E+02	n	1.30E+03	n	6.30E+02	n	
		8.00E-08	I	1.00E+00	I	4.00E+01	I	V		1		2.87E+03	Ethyl Tertiary Butyl Ether (ETBE)	637-92-3	1.30E+04	c**s	5.60E+04	c**s	3.50E+03	c*	1.50E+04	c*	7.00E+03	c**	
				1.00E-05	I					1	0.1		Ethyl-p-nitrophenyl Phosphonate	2104-64-5	6.30E-01	n	8.20E+00	n					8.90E-02	n	
1.10E-02	C	2.50E-06	C	5.00E-02	P	1.00E+00	I	V		1		4.80E+02	Ethylbenzene	100-41-4	5.80E+02	c**s	2.50E+03	c**s	1.10E+02	c**	4.90E+02	c**	1.50E+02	c**	7.00E+02
				7.00E-02	P					1	0.1		Ethylene Cyanohydrin	109-78-4	4.40E+03	n	5.70E+04	n					1.40E+03	n	
				9.00E-02	P			V		1		1.89E+05	Ethylene Diamine	107-15-3	7.00E+03	n	1.10E+05	nm					1.80E+03	n	
				8.00E-01	A	4.00E-01	C			1	0.1		Ethylene Glycol	107-21-1	5.10E+04	n	6.60E+05	nm	4.20E+02	n	1.80E+03	n	1.60E+04	n	
				1.00E-01	I	1.60E+00	I			1	0.1		Ethylene Glycol Monobutyl Ether	111-76-2	6.30E+03	n	8.20E+04	n	1.70E+03	n	7.00E+03	n	2.00E+03	n	
3.10E-01	C	3.00E-03	I			3.00E-02	C	V	M	1		1.21E+05	Ethylene Oxide	75-21-8	2.00E-01	c	2.50E+00	c	3.40E-02	c	4.10E-01	c	6.70E-02	c	
4.50E-02	C	1.30E-05	C	8.00E-05	I					1	0.1		Ethylene Thiourea	96-45-7	5.10E+00	n	6.60E+01	n	2.20E+01	c	9.40E+01	c	1.60E+00	n	
6.50E+01	C	1.90E-02	C					V		1		1.54E+05	Ethyleneimine	151-56-4	2.70E-01	c	1.20E+00	c	1.50E-02	c	6.50E-02	c	2.40E-02	c	
				3.00E+00	I					1	0.1		Ethylphthalyl Ethyl Glycolate	84-72-0	1.90E+05	nm	2.50E+06	nm					5.80E+04	n	
				2.50E-04	I					1	0.1		Fenamiphos	22224-92-6	1.60E+01	n	2.10E+02	n					4.40E+00	n	
				2.50E-02	I					1	0.1		Fenpropathrin	39515-41-8	1.60E+03	n	2.10E+04	n					6.40E+01	n	
				2.50E-02	I					1	0.1		Fenvalerate	51630-58-1	1.60E+03	n	2.10E+04	n					5.00E+02	n	
				1.30E-02	I					1	0.1		Fluometuron	2164-17-2	8.20E+02	n	1.10E+04	n					2.40E+02	n	
				4.00E-02	C	1.30E-02	C			1			Fluoride	16984-48-8	3.10E+03	n	4.70E+04	n	1.40E+01	n	5.70E+01	n	8.00E+02	n	4.00E+03
				6.00E-02	I	1.30E-02	C			1			Fluorine (Soluble Fluoride)	7782-41-4	4.70E+03	n	7.00E+04	n	1.40E+01	n	5.70E+01	n	1.20E+03	n	4.00E+03
				8.00E-02	I					1	0.1		Fluridone	59756-60-4	5.10E+03	n	6.60E+04	n					1.40E+03	n	
				4.00E-02	O					1	0.1		Flurprimidol	56425-91-3	2.50E+03	n	3.30E+04	n					6.90E+02	n	
				2.00E-03	O					1	0.1		Flusilazole	85509-19-9	1.30E+02	n	1.60E+03	n					3.10E+01	n	
				5.00E-01	O					1	0.1		Flutolanil	66332-96-5	3.20E+04	n	4.10E+05	nm					7.90E+03	n	
				1.00E-02	I					1	0.1		Fluvalinate	69409-94-5	6.30E+02	n	8.20E+03	n					2.00E+02	n	
				9.00E-02	O					1	0.1		Folpet	133-07-3	5.70E+03	n	7.40E+04	n					1.60E+03	n	
				1.00E-02	O					1	0.1		Fomesafen	72178-02-0	6.30E+02	n	8.20E+03	n					1.90E+02	n	
				2.00E-03	I					1	0.1		Fonofos	944-22-9	1.30E+02	n	1.60E+03	n					2.40E+01	n	
2.10E-02	C	1.30E-05	I	2.00E-01	I	9.80E-03	A	V		1		4.24E+04	Formaldehyde	50-00-0	7.60E+02	n	3.30E+03	n	1.00E+01	n	4.30E+01	n	2.00E+01	n	
				9.00E-01	P	3.00E-04	X	V		1		1.06E+05	Formic Acid	64-18-6	2.90E+01	n	1.20E+02	n	3.10E-01	n	1.30E+00	n	6.30E-01	n	
				2.50E+00	O					1	0.1		Fosetyl-AL	39148-24-8	1.60E+05	nm	2.10E+06	nm					5.00E+04	n	
													Furans												
				1.00E-03	X			V		1			~Dibenzofuran	132-64-9	7.80E+01	n	1.20E+03	n					7.90E+00	n	
				1.00E-03	I			V		1		6.22E+03	~Furan	110-00-9	7.80E+01	n	1.20E+03	n					1.90E+01	n	
				9.00E-01	I	2.00E+00	I	V		1		1.65E+05	~Tetrahydrofuran	109-99-9	1.80E+04	n	9.50E+04	n	2.10E+03	n	8.80E+03	n	3.40E+03	n	
3.80E+00	H									1	0.1		Furazolidone	67-45-8	1.40E+01	c	6.00E+01	c					2.00E+00	c	
				3.00E-03	I	5.00E-02	H	V		1		1.01E+04	Furfural	98-01-1	2.10E+02	n	2.60E+03	n	5.20E+01	n	2.20E+02	n	3.80E+01	n	
1.50E+00	C	4.30E-04	C							1	0.1		Furium	531-82-8	3.60E+01	c	1.50E+02	c	6.50E-01	c	2.90E+00	c	5.10E+00	c	
3.00E-02	I	8.60E-06	C							1	0.1		Furmecyclox	60568-05-0	1.80E+03	c	7.70E+03	c	3.30E+01	c	1.40E+02	c	1.10E+02	c	

RMLs
TR=1E-04
THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	GIABS	ABS_d	C_{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
				6.00E-03	O					1	0.1		Glufosinate, Ammonium	77182-82-2	3.80E+02	n	4.90E+03	n					1.20E+02	n	
				1.00E-01	A	8.00E-05	C			1	0.1		Glutaraldehyde	111-30-8	6.00E+03	n	7.00E+04	n	8.30E-02	n	3.50E-01	n	2.00E+03	n	
				4.00E-04	I	1.00E-03	X	V		1		1.06E+05	Glycidaldehyde	765-34-4	2.30E+01	n	2.10E+02	n	1.00E+00	n	4.40E+00	n	1.70E+00	n	
				1.00E-01	I					1	0.1		Glyphosate	1071-83-6	6.30E+03	n	8.20E+04	n					2.00E+03	n	7.00E+02
				1.00E-02	X			V		1			Guanidine	113-00-8	7.80E+02	n	1.20E+04	n					2.00E+02	n	
				2.00E-02	P					1	0.1		Guanidine Chloride	50-01-1	1.30E+03	n	1.60E+04	n					4.00E+02	n	
				3.00E-02	X					1	0.1		Guanidine Nitrate	506-93-4	1.90E+03	n	2.50E+04	n					6.00E+02	n	
				5.00E-05	I					1	0.1		Haloxyfop, Methyl	69806-40-2	3.20E+00	n	4.10E+01	n					7.60E-01	n	
4.50E+00	I	1.30E-03	I	1.00E-04	A			V		1			Heptachlor	76-44-8	7.80E+00	n	6.30E+01	c**	2.20E-01	c	9.40E-01	c	1.40E-01	c**	4.00E-01
9.10E+00	I	2.60E-03	I	1.30E-05	I			V		1			Heptachlor Epoxide	1024-57-3	1.00E+00	n	1.50E+01	n	1.10E-01	c	4.70E-01	c	1.20E-01	n	2.00E-01
						3.00E-03	X	V		1		2.09E+02	Heptanal, n-	111-71-7	2.40E+01	n	1.00E+02	n	3.10E+00	n	1.30E+01	n	6.30E+00	n	
				3.00E-04	X	4.00E-01	P	V		1		5.79E+01	Heptane, N-	142-82-5	2.20E+01	n	2.90E+02	ns	4.20E+02	n	1.80E+03	n	6.00E+00	n	
				2.00E-03	I			V		1			Hexabromobenzene	87-82-1	1.60E+02	n	2.30E+03	n					4.00E+01	n	
				2.00E-04	I					1	0.1		Hexabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-153)	68631-49-2	1.30E+01	n	1.60E+02	n					4.00E+00	n	
1.60E+00	I	4.60E-04	I	1.00E-05	P			V		1			Hexachlorobenzene	118-74-1	7.80E-01	n	1.20E+01	n	6.10E-01	c	2.70E+00	c	2.00E-01	n	1.00E+00
7.80E-02	I	2.20E-05	I	1.00E-03	P			V		1		1.68E+01	Hexachlorobutadiene	87-68-3	7.80E+01	ns	5.30E+02	c**s	1.30E+01	c	5.60E+01	c	6.50E+00	n	
6.30E+00	I	1.80E-03	I	8.00E-03	A					1	0.1		Hexachlorocyclohexane, Alpha-	319-84-6	8.60E+00	c*	3.60E+01	c	1.60E-01	c	6.80E-01	c	7.20E-01	c	
1.80E+00	I	5.30E-04	I							1	0.1		Hexachlorocyclohexane, Beta-	319-85-7	3.00E+01	c	1.30E+02	c	5.30E-01	c	2.30E+00	c	2.50E+00	c	
1.10E+00	C	3.10E-04	C	1.00E-05	A					1	0.04		Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	7.10E-01	n	1.00E+01	n	9.10E-01	c	4.00E+00	c	1.20E-01	n	2.00E-01
1.80E+00	I	5.10E-04	I							1	0.1		Hexachlorocyclohexane, Technical	608-73-1	3.00E+01	c	1.30E+02	c	5.50E-01	c	2.40E+00	c	2.50E+00	c	
				6.00E-03	I	2.00E-04	I	V		1		1.57E+01	Hexachlorocyclopentadiene	77-47-4	1.80E+00	n	7.50E+00	n	2.10E-01	n	8.80E-01	n	4.10E-01	n	5.00E+01
4.00E-02	I	1.10E-05	C	7.00E-04	I	3.00E-02	I	V		1			Hexachloroethane	67-72-1	4.50E+01	n	4.60E+02	n	2.60E+01	c**	1.10E+02	c**	6.20E+00	n	
				3.00E-04	I					1	0.1		Hexachlorophene	70-30-4	1.90E+01	n	2.50E+02	n					6.00E+00	n	
8.00E-02	I			4.00E-03	I					1	0.015		Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	3.00E+02	n	3.80E+03	c**					8.00E+01	n	
						1.00E-05	I	V		1		3.39E+03	Hexamethylene Diisocyanate, 1,6-	822-06-0	3.10E+00	n	1.30E+01	n	1.00E-02	n	4.40E-02	n	2.10E-02	n	
						4.00E-04	C			1	0.1		Hexamethylene diisocyanate biuret	4035-89-6	5.70E+05	nm	2.40E+06	nm	4.20E-01	n	1.80E+00	n			
						4.00E-04	C			1	0.1		Hexamethylene diisocyanate isocyanurate	3779-63-3	5.70E+05	nm	2.40E+06	nm	4.20E-01	n	1.80E+00	n			
				4.00E-04	P					1	0.1		Hexamethylphosphoramide	680-31-9	2.50E+01	n	3.30E+02	n					8.00E+00	n	
		2.00E-07	X			6.00E-01	P	V		1		1.41E+02	Hexane, Commercial	E5241997	5.20E+02	ns	2.20E+03	ns	6.30E+02	n	2.60E+03	n	1.30E+03	n	
						7.00E-01	I	V		1		1.41E+02	Hexane, N-	110-54-3	6.10E+02	ns	2.50E+03	ns	7.30E+02	n	3.10E+03	n	1.50E+03	n	
				2.00E+00	P					1	0.1		Hexanedioic Acid	124-04-9	1.30E+05	nm	1.60E+06	nm					4.00E+04	n	
9.50E-03	P			7.00E-02	P	4.00E-04	P	V		1		2.74E+02	Hexanol, 1-,2-ethyl- (2-Ethyl-1-hexanol)	104-76-7	1.50E+01	n	6.30E+01	n	4.20E-01	n	1.80E+00	n	8.30E-01	n	
				5.00E-03	I	3.00E-02	I	V		1		3.28E+03	Hexanone, 2-	591-78-6	2.00E+02	n	1.30E+03	n	3.10E+01	n	1.30E+02	n	3.80E+01	n	
				3.30E-02	I					1	0.1		Hexazinone	51235-04-2	2.10E+03	n	2.70E+04	n					6.40E+02	n	
				2.50E-02	I					1	0.1		Hexythiazox	78587-05-0	1.60E+03	n	2.10E+04	n					1.10E+02	n	
				1.70E-02	O					1	0.1		Hydramethylnon	67485-29-4	1.10E+03	n	1.40E+04	n					3.40E+02	n	
3.00E+00	I	4.90E-03	I			3.00E-05	P	V		1		1.12E+05	Hydrazine	302-01-2	2.00E+00	n	8.60E+00	n	3.10E-02	n	1.30E-01	n	6.30E-02	n	
3.00E+00	I	4.90E-03	I							1			Hydrazine Sulfate	10034-93-2	2.30E+01	c	1.10E+02	c	5.70E-02	c	2.50E-01	c	2.60E+00	c	
						2.00E-02	I	V		1			Hydrogen Chloride	7647-01-0	2.80E+07	nm	1.20E+08	nm	2.10E+01	n	8.80E+01	n	4.20E+01	n	
				4.00E-02	C	1.40E-02	C	V		1			Hydrogen Fluoride	7664-39-3	3.10E+03	n	4.70E+04	n	1.50E+01	n	6.10E+01	n	2.80E+01	n	
						2.00E-03	I	V		1			Hydrogen Sulfide	7783-06-4	2.80E+06	nm	1.20E+07	nm	2.10E+00	n	8.80E+00	n	4.20E+00	n	
6.00E-02	P			4.00E-02	P					1	0.1		Hydroquinone	123-31-9	9.00E+02	c**	3.80E+03	c**					1.30E+02	c**	
6.10E-02	O			1.10E-01	O					1	0.1		Imazalil	35554-44-0	8.90E+02	c**	3.80E+03	c*					9.00E+01	c*	
				2.50E-01	I					1	0.1		Imazaquin	81335-37-7	1.60E+04	n	2.10E+05	nm					4.90E+03	n	
				2.50E+00	O					1	0.1		Imazethapyr	81335-77-5	1.60E+05	nm	2.10E+06	nm					4.70E+04	n	
				1.00E-02	A					1			Iodine	7553-56-2	7.80E+02	n	1.20E+04	n					2.00E+02	n	
				4.00E-02	I					1	0.1		Iprodione	36734-19-7	2.50E+03	n	3.30E+04	n					7.40E+02	n	
				7.00E-01	P					1			Iron	7439-89-6	5.50E+04	n	8.20E+05	nm					1.40E+04	n	
				3.00E-01	I			V		1		1.00E+04	Isobutyl Alcohol	78-83-1	2.30E+04	ns	3.50E+05	nms					5.90E+03	n	
9.50E-04	I			2.00E-01	I	2.00E+00	C			1	0.1		Isophorone	78-59-1	1.30E+04	n	1.60E+05	nm	2.10E+03	n	8.80E+03	n	3.80E+03	n	
				1.50E-02	I			V		1			Isopropalin	33820-53-0	1.20E+03	n	1.80E+04	n					4.00E+01	n	
				2.00E+00	P	2.00E-01	P	V		1		1.09E+05	Isopropanol	67-63-0	5.60E+03	n	2.40E+04	n	2.10E+02	n	8.80E+02	n	4.10E+02	n	
				1.00E-01	I					1	0.1		Isopropyl Methyl Phosphonic Acid	1832-54-8	6.30E+03	n	8.20E+04	n					2.00E+03	n	
				5.00E-02	I					1	0.1		Isoxaben	82558-50-7	3.20E+03	n	4.10E+04	n					7.30E+02	n	
						3.00E-01	A	V		1			JP-7	E1737665	4.30E+08	nm	1.80E+09	nm	3.10E+02	n	1.30E+03	n	6.30E+02	n	
				8.00E-03	O					1	0.1		Lactofen	77501-63-4	5.10E+02	n	6.60E+03	n					1.00E+02	n	
				2.00E-04	X					1	0.1		Lactonitrile	78-97-7	1.30E+01	n	1.60E+02	n					4.00E+00	n	
				5.00E-05	P					1			Lanthanum	7439-91-0	3.90E+00	n	5.80E+01	n					1.00E+00	n	
				2.10E-05	P					1	0.1		Lanthanum Acetate Hydrate	100587-90-4	1.30E+00	n	1.70E+01	n					4.20E-01	n	
				1.90E-05	P					1			Lanthanum Chloride Heptahydrate	10025-84-0	1.50E+00	n	2.20E+01	n					3.70E-01	n	
				2.80E-05	P					1			Lanthanum Chloride, Anhydrous	10099-58-8	2.20E+00	n	3.30E+01	n					5.70E-01	n	
				1.60E-05	P					1			Lanthanum Nitrate Hexahydrate	10277-43-7	1.30E+00	n	1.90E+01	n					3.20E-01	n	
													Lead Compounds												
8.50E-03	C	1.20E-05	C							1			~Lead Phosphate	7446-27-7	8.20E+03	c	3.80E+04	c	2.30E+01	c	1.00E+02	c	9.10E+02	c	
2.10E-01	C	8.00E-05	C							1	0.1		~Lead acetate	301-04-2	2.60E+02	c	1.10E+03	c	3.50E+00	c	1.50E+01	c	3.70E+01	c	
										1			~Lead and Compounds	7439-92-1	4.00E+02	G	8.00E+02	G	1.50E-01	G			1.50E+01	G	1.50E+01
3.80E-02	C	1.10E-05	C							1	0.1		~Lead subacetate	1335-32-6	1.40E+03	c	6.00E+03	c	2.60E+01	c	1.10E+02	c	2.10E+02	c	
				1.00E-07	I			V		1		2.43E+00	~Tetraethyl Lead	78-00-2	7.80E-03	n	1.20E-01	n					1.30E-03	n	
				5.00E-06	P			V		1		3.83E+02	Lewisite	541-25-3	3.90E-01	n	5.80E+00	n					9.00E-02	n	
				7.70E-03	O					1	0.1		Linuron	330-55-2	4.90E+02	n	6.30E+03	n					1.30E+02	n	
				2.00E-03	P					1			Lithium	7439-93-2	1.60E+02	n	2.30E+03	n					4.00E+01	n	

RMLs TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	GIABS	ABS_d	C_{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
				5.00E-04	I					1	0.1		MCPA	94-74-6	3.20E+01	n	4.10E+02	n					7.50E+00	n	
				4.40E-03	O					1	0.1		MCPB	94-81-5	2.80E+02	n	3.60E+03	n					6.50E+01	n	
				1.00E-03	I					1	0.1		MCPP	93-65-2	6.30E+01	n	8.20E+02	n					1.60E+01	n	
				2.00E-02	I					1	0.1		Malathion	121-75-5	1.30E+03	n	1.60E+04	n					3.90E+02	n	
				1.00E-01	I	7.00E-04	C			1	0.1		Maleic Anhydride	108-31-6	6.30E+03	n	8.00E+04	n	7.30E-01	n	3.10E+00	n	1.90E+03	n	
				5.00E-01	I					1	0.1		Maleic Hydrazide	123-33-1	3.20E+04	n	4.10E+05	nm					1.00E+04	n	
				1.00E-04	P					1	0.1		Malononitrile	109-77-3	6.30E+00	n	8.20E+01	n					2.00E+00	n	
				3.00E-02	H					1	0.1		Mancozeb	8018-01-7	1.90E+03	n	2.50E+04	n					5.40E+02	n	
				5.00E-03	I					1	0.1		Maneb	12427-38-2	3.20E+02	n	4.10E+03	n					9.80E+01	n	
				1.40E-01	I	5.00E-05	I			1			Manganese (Diet)	7439-96-5					5.20E-02	n	2.20E-01	n			
				2.40E-02	G	5.00E-05	I			0.04			Manganese (Non-diet)	7439-96-5	1.80E+03	n	2.60E+04	n	5.20E-02	n	2.20E-01	n	4.30E+02	n	
				9.00E-05	H					1	0.1		Mephosfolan	950-10-7	5.70E+00	n	7.40E+01	n					1.80E+00	n	
				3.00E-02	I					1	0.1		Mepiquat Chloride	24307-26-4	1.90E+03	n	2.50E+04	n					6.00E+02	n	
1.10E-02	P			4.00E-03	P					1	0.1		Mercaptobenzothiazole, 2-	149-30-4	2.50E+02	n	3.30E+03	n					7.20E+01	n	
													Mercury Compounds												
				3.00E-04	I	3.00E-04	G			0.07			~Mercuric Chloride (and other Mercury salts)	7487-94-7	2.30E+01	n	3.50E+02	n	3.10E-01	n	1.30E+00	n	5.70E+00	n	2.00E+00
						3.00E-04	I	V		1		3.13E+00	~Mercury (elemental)	7439-97-6	1.10E+01	ns	4.60E+01	ns	3.10E-01	n	1.30E+00	n	6.30E-01	n	2.00E+00
				1.00E-04	I					1			~Methyl Mercury	22967-92-6	7.80E+00	n	1.20E+02	n					2.00E+00	n	
				8.00E-05	I					1	0.1		~Phenylmercuric Acetate	62-38-4	5.10E+00	n	6.60E+01	n					1.60E+00	n	
				3.00E-05	I			V		1			Merphos	150-50-5	2.30E+00	n	3.50E+01	n					6.00E-01	n	
				6.00E-02	I					1	0.1		Metalaxyl	57837-19-1	3.80E+03	n	4.90E+04	n					1.20E+03	n	
				1.00E-04	I	3.00E-02	P	V		1		4.58E+03	Methacrylonitrile	126-98-7	7.50E+00	n	1.00E+02	n	3.10E+01	n	1.30E+02	n	1.90E+00	n	
				5.00E-05	I					1	0.1		Methamidophos	10265-92-6	3.20E+00	n	4.10E+01	n					1.00E+00	n	
				2.00E+00	I	2.00E+01	I	V		1		1.06E+05	Methanol	67-56-1	1.20E+05	nms	1.20E+06	nms	2.10E+04	n	8.80E+04	n	2.00E+04	n	
				1.50E-03	O					1	0.1		Methidathion	950-37-8	9.50E+01	n	1.20E+03	n					2.90E+01	n	
				2.50E-02	I					1	0.1		Methomyl	16752-77-5	1.60E+03	n	2.10E+04	n					5.00E+02	n	
4.90E-02	C									1	0.1		Methoxy-5-nitroaniline, 2-	99-59-2	1.10E+03	c	4.70E+03	c					1.50E+02	c	
				5.00E-03	I					1	0.1		Methoxychlor	72-43-5	3.20E+02	n	4.10E+03	n					3.70E+01	n	4.00E+01
				8.00E-03	P	1.00E-03	P	V		1		1.15E+05	Methoxyethanol Acetate, 2-	110-49-6	1.10E+02	n	5.10E+02	n	1.00E+00	n	4.40E+00	n	2.10E+00	n	
				5.00E-03	P	7.00E-03	P	V		1		1.06E+05	Methoxyethanol, 2-	109-86-4	2.60E+02	n	2.00E+03	n	7.30E+00	n	3.10E+01	n	1.30E+01	n	
				1.00E+00	X			V		1		2.90E+04	Methyl Acetate	79-20-9	7.80E+04	ns	1.20E+06	nms					2.00E+04	n	
						2.00E-02	P	V		1		6.75E+03	Methyl Acrylate	96-33-3	1.50E+02	n	6.10E+02	n	2.10E+01	n	8.80E+01	n	4.20E+01	n	
				6.00E-01	I	5.00E+00	I	V		1		2.84E+04	Methyl Ethyl Ketone (2-Butanone)	78-93-3	2.70E+04	n	1.90E+05	nms	5.20E+03	n	2.20E+04	n	5.60E+03	n	
		1.00E-03	X	1.00E-03	P	2.00E-05	X	V		1		1.80E+05	Methyl Hydrazine	60-34-4	1.00E+00	n	4.40E+00	n	2.10E-02	n	8.80E-02	n	4.20E-02	n	
						3.00E+00	I	V		1		3.36E+03	Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1	3.30E+04	ns	1.40E+05	nms	3.10E+03	n	1.30E+04	n	6.30E+03	n	
						1.00E-03	C	V		1		1.01E+04	Methyl Isocyanate	624-83-9	4.60E+00	n	1.90E+01	n	1.00E+00	n	4.40E+00	n	2.10E+00	n	
				1.40E+00	I	7.00E-01	I	V		1		2.36E+03	Methyl Methacrylate	80-62-6	4.40E+03	ns	1.90E+04	ns	7.30E+02	n	3.10E+03	n	1.40E+03	n	
				2.50E-04	I					1	0.1		Methyl Parathion	298-00-0	1.60E+01	n	2.10E+02	n					4.50E+00	n	
				6.00E-02	X					1	0.1		Methyl Phosphonic Acid	993-13-5	3.80E+03	n	4.90E+04	n					1.20E+03	n	
				6.00E-03	H	4.00E-02	H	V		1		3.93E+02	Methyl Styrene (Mixed Isomers)	25013-15-4	3.20E+02	n	2.60E+03	ns	4.20E+01	n	1.80E+02	n	2.30E+01	n	
9.90E-02	C	2.80E-05	C							1	0.1		Methyl methanesulfonate	66-27-3	5.50E+02	c	2.30E+03	c	1.00E+01	c	4.40E+01	c	7.90E+01	c	
1.80E-03	C	2.60E-07	C			3.00E+00	I	V		1		8.87E+03	Methyl tert-Butyl Ether (MTBE)	1634-04-4	4.70E+03	c**	2.10E+04	c**s	1.10E+03	c**	4.70E+03	c**	1.40E+03	c**	
				3.00E-04	X					1	0.1		Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2	1.90E+01	n	2.50E+02	n					6.00E+00	n	
						3.00E+00	X	V		1		2.45E+03	Methyl-2-Pentanol, 4-	108-11-2	5.40E+04	ns	2.30E+05	nms	3.10E+03	n	1.30E+04	n	6.30E+03	n	
9.00E-03	P			2.00E-02	X					1	0.1		Methyl-5-Nitroaniline, 2-	99-55-8	1.30E+03	n	1.60E+04	n					3.80E+02	n	
8.30E+00	C	2.40E-03	C							1	0.1		Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	6.50E+00	c	2.80E+01	c	1.20E-01	c	5.10E-01	c	9.40E-01	c	
1.30E-01	C	3.70E-05	C							1	0.1		Methylaniline Hydrochloride, 2-	636-21-5	4.20E+02	c	1.80E+03	c	7.60E+00	c	3.30E+01	c	6.00E+01	c	
				1.00E-02	A					1	0.1		Methylarsonic acid	124-58-3	6.30E+02	n	8.20E+03	n					2.00E+02	n	
				2.00E-04	X					1	0.1		Methylbenzene,1-4-diamine monohydrochloride, 2-	74612-12-7	1.30E+01	n	1.60E+02	n					4.00E+00	n	
1.00E-01	X			3.00E-04	X					1	0.1		Methylbenzene-1,4-diamine sulfate, 2-	615-50-9	1.90E+01	n	2.50E+02	n					6.00E+00	n	
2.20E+01	C	6.30E-03	C						M	1	0.1		Methylcholanthrene, 3-	56-49-5	5.50E-01	c	1.00E+01	c	1.60E-02	c	1.90E-01	c	1.10E-01	c	
2.00E-03	I	1.00E-08	I	6.00E-03	I	6.00E-01	I	V	M	1		3.32E+03	Methylene Chloride	75-09-2	3.50E+02	n	3.20E+03	n	6.30E+02	n	2.60E+03	n	1.10E+02	n	5.00E+00
1.00E-01	P	4.30E-04	C	2.00E-03	P				M	1	0.1		Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	1.20E+02	c**	1.60E+03	n	2.40E-01	c	2.90E+00	c	1.60E+01	c**	
4.60E-02	I	1.30E-05	C							1	0.1		Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	1.20E+03	c	5.00E+03	c	2.20E+01	c	9.40E+01	c	7.00E+01	c	
1.60E+00	C	4.60E-04	C			2.00E-02	C			1	0.1		Methylenebisbenzenamine, 4,4'-	101-77-9	3.40E+01	c	1.40E+02	c	6.10E-01	c*	2.70E+00	c*	4.70E+00	c	
						6.00E-04	I			1	0.1		Methylenediphenyl Diisocyanate	101-68-8	8.50E+05	nm	3.60E+06	nm	6.30E-01	n	2.60E+00	n			
				7.00E-02	H			V		1		5.00E+02	Methylstyrene, Alpha-	98-83-9	5.50E+03	ns	8.20E+04	ns					7.80E+02	n	
				1.50E-01	I					1	0.1		Metolachlor	51218-45-2	9.50E+03	n	1.20E+05	nm					2.70E+03	n	
				2.50E-02	I					1	0.1		Metribuzin	21087-64-9	1.60E+03	n	2.10E+04	n					4.90E+02	n	
				2.50E-01	I					1	0.1		Metsulfuron-methyl	74223-64-6	1.60E+04	n	2.10E+05	nm					4.90E+03	n	
		4.50E-06	X	1.00E-02	X	1.00E-01	P	V		1		6.86E+00	Midrange Aliphatic Hydrocarbon Streams	E1790669	6.50E+01	c**s	2.80E+02	c**s	6.20E+01	c**	2.70E+02	c**	1.00E+02	n	
				3.00E+00	P			V		1		3.42E-01	Mineral oils	8012-95-1	2.30E+05	nms	3.50E+06	nms					6.00E+04	n	
1.80E+01	C	5.10E-03	C	2.00E-04	I			V		1			Mirex	2385-85-5	3.60E+00	c**	1.70E+01	c*	5.50E-02	c	2.40E-01	c	8.80E-02	c*	
				2.00E-03	I					1	0.1		Molinate	2212-67-1	1.30E+02	n	1.60E+03	n					3.00E+01	n	
				5.00E-03	I	2.00E-03	A			1			Molybdenum	7439-98-7	3.90E+02	n	5.80E+03	n	2.10E+00	n	8.80E+00	n	1.00E+02	n	
				1.00E-01	I					1			Monochloramine	10599-90-3	7.80E+03	n	1.20E+05	nm					2.00E+03	n	4.0E+03(G)
				2.00E-03	P					1	0.1		Monomethylaniline	100-61-8	1.30E+02	n	1.60E+03	n					3.80E+01	n	
				2.50E-02	I					1	0.1		Myclobutanil	88671-89-0	1.60E+03	n	2.10E+04	n					4.50E+02	n	
				3.00E-04	X					1	0.1		N,N'-Diphenyl-1,4-benzenediamine	74-31-7	1.90E+01	n	2.50E+02	n					3.60E+00	n	
				2.00E-03	I			V		1			Naled	300-76-5	1.60E+02	n	2.30E+03	n					4.00E+01	n	
				3.00E-02	X	1.00E-01	P	V		1			Naphtha, High Flash Aromatic (HFAN)	64742-95-6	2.30E+03	n	3.50E+04	n	1.00E+02	n	4.40E+02	n	1.50E+02	n	

RMLs
R=1E-04
THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD$_o$ (mg/kg-day)	key	RfC$_i$ (mg/m^3)	key	vol	mutagen	GIABS	ABS$_d$	C$_{sat}$ (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
1.80E+00	C	0.00E+00	C							1	0.1		Naphthylamine, 2-	91-59-8	3.00E+01	c	1.30E+02	c					3.90E+00	c	
				1.20E-01	O					1	0.1		Napropamide	15299-99-7	7.60E+03	n	9.80E+04	n					2.00E+03	n	
		2.60E-04	C	1.10E-02	C	1.40E-05	C			1	0.1		Nickel Acetate	373-02-4	6.70E+02	n	8.10E+03	n	1.50E-02	n	6.10E-02	n	2.20E+02	n	
		2.60E-04	C	1.10E-02	C	1.40E-05	C			1	0.1		Nickel Carbonate	3333-67-3	6.70E+02	n	8.10E+03	n	1.50E-02	n	6.10E-02	n	2.20E+02	n	
		2.60E-04	C	1.10E-02	C	1.40E-05	C	V		1			Nickel Carbonyl	13463-39-3	8.20E+02	n	1.10E+04	n	1.50E-02	n	6.10E-02	n	2.90E-02	n	
		2.60E-04	C	1.10E-02	C	1.40E-05	C			0.04			Nickel Hydroxide	12054-48-7	8.20E+02	n	1.10E+04	n	1.50E-02	n	6.10E-02	n	2.00E+02	n	
		2.60E-04	C	1.10E-02	C	2.00E-05	C			0.04			Nickel Oxide	1313-99-1	8.40E+02	n	1.20E+04	n	2.10E-02	n	8.80E-02	n	2.00E+02	n	
		2.40E-04	I	1.10E-02	C	1.40E-05	C			0.04			Nickel Refinery Dust	E715532	8.20E+02	n	1.10E+04	n	1.50E-02	n	6.10E-02	n	2.20E+02	n	
		2.60E-04	C	2.00E-02	I	9.00E-05	A			0.04			Nickel Soluble Salts	7440-02-0	1.50E+03	n	2.20E+04	n	9.40E-02	n	3.90E-01	n	3.90E+02	n	
1.70E+00	C	4.80E-04	I	1.10E-02	C	1.40E-05	C			0.04			Nickel Subsulfide	12035-72-2	4.10E+01	c*	1.90E+02	c*	1.50E-02	n	6.10E-02	n	4.50E+00	c*	
9.10E-01	C	2.60E-04	C	1.10E-02	C	1.40E-05	C			1	0.1		Nickelocene	1271-28-9	6.00E+01	c*	2.50E+02	c*	1.50E-02	n	6.10E-02	n	8.60E+00	c*	
				1.60E+00	I					1			Nitrate (measured as nitrogen)	14797-55-8	1.30E+05	nm	1.90E+06	nm					3.20E+04	n	1.00E+04
										1			Nitrate + Nitrite (measured as nitrogen)	E701177											1.00E+04
				1.00E-01	I					1			Nitrite (measured as nitrogen)	14797-65-0	7.80E+03	n	1.20E+05	nm					2.00E+03	n	1.00E+03
				1.00E-02	X	5.00E-05	X			1	0.1		Nitroaniline, 2-	88-74-4	6.30E+02	n	8.00E+03	n	5.20E-02	n	2.20E-01	n	1.90E+02	n	
2.00E-02	P			4.00E-03	P	6.00E-03	P			1	0.1		Nitroaniline, 4-	100-01-6	2.50E+02	n	3.30E+03	n	6.30E+00	n	2.60E+01	n	7.80E+01	n	
		4.00E-05	I	2.00E-03	I	9.00E-03	I	V		1		3.05E+03	Nitrobenzene	98-95-3	1.30E+02	n	1.30E+03	n	7.00E+00	c**	3.10E+01	c**	1.30E+01	n	
				3.00E+03	P					1	0.1		Nitrocellulose	9004-70-0	1.90E+08	nm	2.50E+09	nm					6.00E+07	n	
				7.00E-02	H					1	0.1		Nitrofurantoin	67-20-9	4.40E+03	n	5.70E+04	n					1.40E+03	n	
1.30E+00	C	3.70E-04	C							1	0.1		Nitrofurazone	59-87-0	4.20E+01	c	1.80E+02	c	7.60E-01	c	3.30E+00	c	6.00E+00	c	
1.70E-02	P			1.00E-04	P					1	0.1		Nitroglycerin	55-63-0	6.30E+00	n	8.20E+01	n					2.00E+00	n	
				1.00E-01	I					1	0.1		Nitroguanidine	556-88-7	6.30E+03	n	8.20E+04	n					2.00E+03	n	
		8.80E-06	P			5.00E-03	P	V		1		1.80E+04	Nitromethane	75-52-5	8.80E+01	n	3.70E+02	n	5.20E+00	n	2.20E+01	n	1.00E+01	n	
		5.80E-04	X			2.00E-02	I	V		1		4.86E+03	Nitropropane, 2-	79-46-9	6.40E+00	c*	2.80E+01	c*	4.80E-01	c*	2.10E+00	c*	9.70E-01	c*	
2.70E+01	C	7.70E-03	C						M	1	0.1		Nitroso-N-ethylurea, N-	759-73-9	4.50E-01	c	8.50E+00	c	1.30E-02	c	1.60E-01	c	9.20E-02	c	
1.20E+02	C	3.40E-02	C						M	1	0.1		Nitroso-N-methylurea, N-	684-93-5	1.00E-01	c	1.90E+00	c	3.00E-03	c	3.60E-02	c	2.10E-02	c	
5.40E+00	I	1.60E-03	I					V		1			Nitroso-di-N-butylamine, N-	924-16-3	9.90E+00	c	4.60E+01	c	1.80E-01	c	7.70E-01	c	2.70E-01	c	
7.00E+00	I	2.00E-03	C							1	0.1		Nitroso-di-N-propylamine, N-	621-64-7	7.80E+00	c	3.30E+01	c	1.40E-01	c	6.10E-01	c	1.10E+00	c	
2.80E+00	I	8.00E-04	C							1	0.1		Nitrosodiethanolamine, N-	1116-54-7	1.90E+01	c	8.20E+01	c	3.50E-01	c	1.50E+00	c	2.80E+00	c	
1.50E+02	I	4.30E-02	I						M	1	0.1		Nitrosodiethylamine, N-	55-18-5	8.10E-02	c	1.50E+00	c	2.40E-03	c	2.90E-02	c	1.70E-02	c	
5.10E+01	I	1.40E-02	I	8.00E-06	P	4.00E-05	X	V	M	1		2.37E+05	Nitrosodimethylamine, N-	62-75-9	2.00E-01	c**	3.40E+00	c**	7.20E-03	c**	8.80E-02	c**	1.10E-02	c**	
4.90E-03	I	2.60E-06	C							1	0.1		Nitrosodiphenylamine, N-	86-30-6	1.10E+04	c	4.70E+04	c	1.10E+02	c	4.70E+02	c	1.20E+03	c	
2.20E+01	I	6.30E-03	C					V		1		1.08E+05	Nitrosomethylethylamine, N-	10595-95-6	2.00E+00	c	9.10E+00	c	4.50E-02	c	1.90E-01	c	7.10E-02	c	
6.70E+00	C	1.90E-03	C							1	0.1		Nitrosomorpholine [N-]	59-89-2	8.10E+00	c	3.40E+01	c	1.50E-01	c	6.50E-01	c	1.20E+00	c	
9.40E+00	C	2.70E-03	C							1	0.1		Nitrosopiperidine [N-]	100-75-4	5.80E+00	c	2.40E+01	c	1.00E-01	c	4.50E-01	c	8.20E-01	c	
2.10E+00	I	6.10E-04	I							1	0.1		Nitrosopyrrolidine, N-	930-55-2	2.60E+01	c	1.10E+02	c	4.60E-01	c	2.00E+00	c	3.70E+00	c	
				1.00E-04	X					1	0.1		Nitrotoluene, m-	99-08-1	6.30E+00	n	8.20E+01	n					1.70E+00	n	
2.20E-01	P			9.00E-04	P			V		1		1.51E+03	Nitrotoluene, o-	88-72-2	7.00E+01	n	1.10E+03	n					1.60E+01	n	
1.60E-02	P			4.00E-03	P					1	0.1		Nitrotoluene, p-	99-99-0	2.50E+02	n	3.30E+03	n					7.10E+01	n	
				3.00E-04	X	2.00E-02	P	V		1		6.86E+00	Nonane, n-	111-84-2	1.10E+01	ns	7.20E+01	ns	2.10E+01	n	8.80E+01	n	5.30E+00	n	
				1.50E-03	O					1	0.1		Norflurazon	27314-13-2	9.50E+01	n	1.20E+03	n					2.90E+01	n	
				3.00E-03	I					1	0.1		Octabromodiphenyl Ether	32536-52-0	1.90E+02	n	2.50E+03	n					6.00E+01	n	
				5.00E-02	I					1	0.006		Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0	3.90E+03	n	5.70E+04	n					1.00E+03	n	
				2.00E-03	H					1	0.1		Octamethylpyrophosphoramide	152-16-9	1.30E+02	n	1.60E+03	n					4.00E+01	n	
7.80E-03	O			1.90E-01	O					1	0.1		Oryzalin	19044-88-3	7.00E+03	c**	2.90E+04	c**					7.90E+02	c**	
				5.00E-03	I					1	0.1		Oxadiazon	19666-30-9	3.20E+02	n	4.10E+03	n					4.70E+01	n	
				2.50E-02	I					1	0.1		Oxamyl	23135-22-0	1.60E+03	n	2.10E+04	n					5.00E+02	n	2.00E+02
7.30E-02	O			4.00E-02	O					1	0.1		Oxyfluorfen	42874-03-3	7.40E+02	c**	3.10E+03	c*					5.40E+01	c**	
				1.30E-02	I					1	0.1		Paclobutrazol	76738-62-0	8.20E+02	n	1.10E+04	n					2.30E+02	n	
				4.50E-03	I					1	0.1		Paraquat Dichloride	1910-42-5	2.80E+02	n	3.70E+03	n					9.00E+01	n	
				6.00E-03	H					1	0.1		Parathion	56-38-2	3.80E+02	n	4.90E+03	n					8.60E+01	n	
				5.00E-02	H			V		1			Pebulate	1114-71-2	3.90E+03	n	5.80E+04	n					5.60E+02	n	
				3.00E-01	O					1	0.1		Pendimethalin	40487-42-1	1.90E+04	n	2.50E+05	nm					1.40E+03	n	
				2.00E-03	I			V		1		3.12E-01	Pentabromodiphenyl Ether	32534-81-9	1.60E+02	ns	2.30E+03	ns					4.00E+01	n	
				1.00E-04	I					1	0.1		Pentabromodiphenyl ether, 2,2',4,4',5- (BDE-99)	60348-60-9	6.30E+00	n	8.20E+01	n					2.00E+00	n	
				8.00E-04	I			V		1			Pentachlorobenzene	608-93-5	6.30E+01	n	9.30E+02	n					3.20E+00	n	
9.00E-02	P							V		1		4.57E+02	Pentachloroethane	76-01-7	7.70E+02	cs	3.60E+03	cs					6.50E+01	c	
2.60E-01	H			3.00E-03	I			V		1			Pentachloronitrobenzene	82-68-8	2.30E+02	n	1.30E+03	c**					1.20E+01	c**	
4.00E-01	I	5.10E-06	C	5.00E-03	I					1	0.25		Pentachlorophenol	87-86-5	1.00E+02	c**	4.00E+02	c**	5.50E+01	c	2.40E+02	c	4.10E+00	c**	1.00E+00
4.30E-03	X			9.00E-03	P					1	0.1		Pentaerythritol tetranitrate (PETN)	78-11-5	5.70E+02	n	7.40E+03	n					1.70E+02	n	
				1.00E-04	X					1	0.1		Pentamethylphosphoramide (PMPA)	10159-46-3	6.30E+00	n	8.20E+01	n					2.00E+00	n	
						1.00E+00	P	V		1		3.88E+02	Pentane, n-	109-66-0	8.10E+02	ns	3.40E+03	ns	1.00E+03	n	4.40E+03	n	2.10E+03	n	
													Per- and Polyfluoroalkyl Substances (PFAS)												
				3.00E-06	D					1	0.1		~Ammonium perfluoro-2-methyl-3-oxahexanoate	62037-80-3	1.90E-01	n	2.50E+00	n					6.00E-02	n	
				3.00E-06	D			V		1			~Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6	2.30E-01	n	3.50E+00	n					6.00E-02	n	
				3.00E-04	P					1	0.1		~Perfluorobutanesulfonate	45187-15-3	1.90E+01	n	2.50E+02	n					6.00E+00	n	
				3.00E-04	P					1	0.1		~Perfluorobutanesulfonic acid (PFBS)	375-73-5	1.90E+01	n	2.50E+02	n					6.00E+00	n	
				2.00E-05	A					1	0.1		~Perfluorohexanesulfonate	108427-53-8	1.30E+00	n	1.60E+01	n					3.90E-01	n	
				2.00E-05	A					1	0.1		~Perfluorohexanesulfonic acid (PFHxS)	355-46-4	1.30E+00	n	1.60E+01	n					3.90E-01	n	
				3.00E-06	A					1	0.1		~Perfluorononanoate	72007-68-2	1.90E-01	n	2.50E+00	n					5.90E-02	n	
				3.00E-06	A					1	0.1		~Perfluorononanoic acid (PFNA)	375-95-1	1.90E-01	n	2.50E+00	n					5.90E-02	n	

RMLS TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	GIABS	ABS_d	C_{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
				2.00E-06	A					1	0.1		~Perfluorooctanesulfonate	45298-90-6	1.30E-01	n	1.60E+00	n					4.00E-02	n	
				2.00E-06	A					1	0.1		~Perfluorooctanesulfonic acid (PFOS)	1763-23-1	1.30E-01	n	1.60E+00	n					4.00E-02	n	
7.00E-02	D			3.00E-06	A					1	0.1		~Perfluorooctanoate	45285-51-6	1.90E-01	n	2.50E+00	n					6.00E-02	n	
7.00E-02	D			3.00E-06	A					1	0.1		~Perfluorooctanoic acid (PFOA)	335-67-1	1.90E-01	n	2.50E+00	n					6.00E-02	n	
				3.00E-04	P					1	0.1		~Potassium perfluorobutanesulfonate	29420-49-3	1.90E+01	n	2.50E+02	n					6.00E+00	n	
				2.00E-06	A					1	0.1		~Potassium perfluorooctanesulfonate	2795-39-3	1.30E-01	n	1.60E+00	n					4.00E-02	n	
													Perchlorates													
				7.00E-04	I					1			~Ammonium Perchlorate	7790-98-9	5.50E+01	n	8.20E+02	n					1.40E+01	n	
				7.00E-04	I					1			~Lithium Perchlorate	7791-03-9	5.50E+01	n	8.20E+02	n					1.40E+01	n	
				7.00E-04	I					1			~Perchlorate and Perchlorate Salts	14797-73-0	5.50E+01	n	8.20E+02	n					1.40E+01	n	1.5E+01(G)
				7.00E-04	I					1			~Potassium Perchlorate	7778-74-7	5.50E+01	n	8.20E+02	n					1.40E+01	n	
				7.00E-04	I					1			~Sodium Perchlorate	7601-89-0	5.50E+01	n	8.20E+02	n					1.40E+01	n	
				5.00E-02	I					1	0.1		Permethrin	52645-53-1	3.20E+03	n	4.10E+04	n					1.00E+03	n	
2.20E-03	C	6.30E-07	C							1	0.1		Phenacetin	62-44-2	2.50E+04	c	1.00E+05	cm	4.50E+02	c	1.90E+03	c	3.40E+03	c	
				2.40E-01	O					1	0.1		Phenmedipham	13684-63-4	1.50E+04	n	2.00E+05	nm					3.80E+03	n	
				3.00E-01	I	2.00E-01	C			1	0.1		Phenol	108-95-2	1.90E+04	n	2.50E+05	nm	2.10E+02	n	8.80E+02	n	5.80E+03	n	
				4.00E-03	I					1	0.1		Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1	2.50E+02	n	3.30E+03	n					7.80E+01	n	
				5.00E-04	X					1	0.1		Phenothiazine	92-84-2	3.20E+01	n	4.10E+02	n					4.30E+00	n	
				2.00E-04	X			V		1		1.29E+02	Phenyl Isothiocyanate	103-72-0	1.60E+01	n	2.30E+02	ns					2.60E+00	n	
				6.00E-03	I					1	0.1		Phenylenediamine, m-	108-45-2	3.80E+02	n	4.90E+03	n					1.20E+02	n	
1.20E-01	P			4.00E-03	P					1	0.1		Phenylenediamine, o-	95-54-5	2.50E+02	n	1.90E+03	c**					6.50E+01	c**	
				1.00E-03	X					1	0.1		Phenylenediamine, p-	106-50-3	6.30E+01	n	8.20E+02	n					2.00E+01	n	
1.90E-03	H									1	0.1		Phenylphenol, 2-	90-43-7	2.80E+04	c	1.20E+05	cm					3.00E+03	c	
				2.00E-04	H					1	0.1		Phorate	298-02-2	1.30E+01	n	1.60E+02	n					3.00E+00	n	
						3.00E-04	I	V		1		1.61E+03	Phosgene	75-44-5	3.10E-01	n	1.30E+00	n	3.10E-01	n	1.30E+00	n	6.30E-01	n	
				2.00E-02	I					1	0.1		Phosmet	732-11-6	1.30E+03	n	1.60E+04	n					3.70E+02	n	
				3.00E-04	I	3.00E-04	I	V		1			Phosphine	7803-51-2	2.30E+01	n	3.50E+02	n	3.10E-01	n	1.30E+00	n	5.70E-01	n	
						1.00E-02	I			1			Phosphoric Acid	7664-38-2	1.40E+07	nm	6.00E+07	nm	1.00E+01	n	4.40E+01	n			
				2.00E-05	I			V		1			Phosphorus, White	7723-14-0	1.60E+00	n	2.30E+01	n					4.00E-01	n	
													Phthalates													
1.40E-02	I	2.40E-06	C	2.00E-02	I					1	0.1		~Bis(2-ethylhexyl)phthalate	117-81-7	1.30E+03	n	1.60E+04	c**	1.20E+02	c	5.10E+02	c	4.00E+02	n	6.00E+00
1.90E-03	P			2.00E-01	I					1	0.1		~Butyl Benzyl Phthalate	85-68-7	1.30E+04	n	1.20E+05	c**m					1.60E+03	c**	
				1.00E+00	I					1	0.1		~Butylphthalyl Butylglycolate	85-70-1	6.30E+04	n	8.20E+05	nm					1.30E+04	n	
				1.00E-01	I					1	0.1		~Dibutyl Phthalate	84-74-2	6.30E+03	n	8.20E+04	n					9.00E+02	n	
				8.00E-01	I					1	0.1		~Diethyl Phthalate	84-66-2	5.10E+04	n	6.60E+05	nm					1.50E+04	n	
				1.00E-01	I			V		1			~Dimethylterephthalate	120-61-6	7.80E+03	n	1.20E+05	nm					1.90E+03	n	
				1.00E-02	P					1	0.1		~Octyl Phthalate, di-N-	117-84-0	6.30E+02	n	8.20E+03	n					2.00E+02	n	
				5.00E-01	X					1	0.1		~Phthalic Acid, p-	100-21-0	3.20E+04	n	4.10E+05	nm					9.40E+03	n	
				2.00E+00	I	2.00E-02	C			1	0.1		~Phthalic Anhydride	85-44-9	1.30E+05	nm	1.60E+06	nm	2.10E+01	n	8.80E+01	n	3.90E+04	n	
				7.00E-02	I					1	0.1		Picloram	1918-02-1	4.40E+03	n	5.70E+04	n					1.40E+03	n	5.00E+02
				1.00E-04	X					1	0.1		Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3	6.30E+00	n	8.20E+01	n					2.00E+00	n	
				2.00E-03	X					1	0.1		Picric Acid (2,4,6-Trinitrophenol)	88-89-1	1.30E+02	n	1.60E+03	n					4.00E+01	n	
				7.30E-04	O					1	0.1		Pirimiphos, Methyl	29232-93-7	4.60E+01	n	6.00E+02	n					8.90E+00	n	
3.00E+01	C	8.60E-03	C	7.00E-06	H					1	0.1		Polybrominated Biphenyls	36355-01-8	4.40E-01	n	5.70E+00	n	3.30E-02	c	1.40E-01	c	1.40E-01	n	
													Polychlorinated Biphenyls (PCBs)													
7.00E-02	G	2.00E-05	G	7.00E-05	I			V		1	0.14		~Aroclor 1016	12674-11-2	4.10E+00	n	5.10E+01	n	1.40E+01	c	6.10E+01	c	1.40E+00	n	
2.00E+00	G	5.70E-04	G					V		1	0.14		~Aroclor 1221	11104-28-2	2.00E+01	c	8.30E+01	c	4.90E-01	c	2.10E+00	c	4.70E-01	c	
2.00E+00	G	5.70E-04	G					V		1	0.14		~Aroclor 1232	11141-16-5	1.70E+01	c	7.20E+01	c	4.90E-01	c	2.10E+00	c	4.70E-01	c	
2.00E+00	G	5.70E-04	G					V		1	0.14		~Aroclor 1242	53469-21-9	2.30E+01	c	9.50E+01	c	4.90E-01	c	2.10E+00	c	7.80E-01	c	
2.00E+00	G	5.70E-04	G					V		1	0.14		~Aroclor 1248	12672-29-6	2.30E+01	c	9.40E+01	c	4.90E-01	c	2.10E+00	c	7.80E-01	c	
2.00E+00	G	5.70E-04	G	2.00E-05	I			V		1	0.14		~Aroclor 1254	11097-69-1	1.20E+00	n	1.50E+01	n	4.90E-01	c	2.10E+00	c	4.00E-01	n	
2.00E+00	G	5.70E-04	G					V		1	0.14		~Aroclor 1260	11096-82-5	2.40E+01	c	9.90E+01	c	4.90E-01	c	2.10E+00	c	7.80E-01	c	
				6.00E-04	X			V		1	0.14		~Aroclor 5460	11126-42-4	3.50E+01	n	4.40E+02	n					1.20E+01	n	
3.90E+00	W	1.10E-03	W	2.30E-05	W	1.30E-03	W	V		1	0.14		~Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	1.40E+00	n	1.70E+01	n	2.50E-01	c**	1.10E+00	c**	4.00E-01	c**	
3.90E+00	W	1.10E-03	W	2.30E-05	W	1.30E-03	W	V		1	0.14		~Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167)	52663-72-6	1.40E+00	n	1.70E+01	n	2.50E-01	c**	1.10E+00	c**	4.00E-01	c**	
3.90E+00	W	1.10E-03	W	2.30E-05	W	1.30E-03	W	V		1	0.14		~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	1.40E+00	n	1.70E+01	n	2.50E-01	c**	1.10E+00	c**	4.00E-01	c**	
3.90E+00	W	1.10E-03	W	2.30E-05	W	1.30E-03	W	V		1	0.14		~Hexachlorobiphenyl, 2,3,3',4,4',5- (PCB 156)	38380-08-4	1.40E+00	n	1.70E+01	n	2.50E-01	c**	1.10E+00	c**	4.00E-01	c**	
3.90E+03	W	1.10E+00	W	2.30E-08	W	1.30E-06	W	V		1	0.14		~Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	1.40E-03	n	1.70E-02	n	2.50E-04	c**	1.10E-03	c**	4.00E-04	c**	
3.90E+00	W	1.10E-03	W	2.30E-05	W	1.30E-03	W	V		1	0.14		~Pentachlorobiphenyl, 2',3,4,4',5- (PCB 123)	65510-44-3	1.40E+00	n	1.70E+01	n	2.50E-01	c**	1.10E+00	c**	4.00E-01	c**	
3.90E+00	W	1.10E-03	W	2.30E-05	W	1.30E-03	W	V		1	0.14		~Pentachlorobiphenyl, 2,3',4,4',5- (PCB 118)	31508-00-6	1.40E+00	n	1.70E+01	n	2.50E-01	c**	1.10E+00	c**	4.00E-01	c**	
3.90E+00	W	1.10E-03	W	2.30E-05	W	1.30E-03	W	V		1	0.14		~Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	1.40E+00	n	1.70E+01	n	2.50E-01	c**	1.10E+00	c**	4.00E-01	c**	
3.90E+00	W	1.10E-03	W	2.30E-05	W	1.30E-03	W	V		1	0.14		~Pentachlorobiphenyl, 2,3,4,4',5- (PCB 114)	74472-37-0	1.40E+00	n	1.70E+01	n	2.50E-01	c**	1.10E+00	c**	4.00E-01	c**	
1.30E+04	W	3.80E+00	W	7.00E-09	W	4.00E-07	W	V		1	0.14		~Pentachlorobiphenyl, 3,3',4,4',5- (PCB 126)	57465-28-8	4.10E-04	n	5.10E-03	n	7.40E-05	c**	3.20E-04	c**	1.20E-04	c**	
2.00E+00	I	5.70E-04	I					V		1	0.14		~Polychlorinated Biphenyls (high risk)	1336-36-3	2.30E+01	c	9.40E+01	c	4.90E-01	c	2.10E+00	c			5.00E-01
4.00E-01	I	1.00E-04	I					V		1	0.14		~Polychlorinated Biphenyls (low risk)	1336-36-3					2.80E+00	c	1.20E+01	c	4.40E+00	c	5.00E-01
7.00E-02	I	2.00E-05	I					V		1	0.14		~Polychlorinated Biphenyls (lowest risk)	1336-36-3					1.40E+01	c	6.10E+01	c			5.00E-01
1.30E+01	W	3.80E-03	W	7.00E-06	W	4.00E-04	W			1	0.14		~Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	4.10E-01	n	5.10E+00	n	7.40E-02	c**	3.20E-01	c**	1.40E-01	n	
3.90E+01	W	1.10E-02	W	2.30E-06	W	1.30E-04	W	V		1	0.14		~Tetrachlorobiphenyl, 3,4,4',5- (PCB 81)	70362-50-4	1.40E-01	n	1.70E+00	n	2.50E-02	c**	1.10E-01	c**	4.00E-02	c**	
						6.00E-04	I			1	0.1		Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9	8.50E+05	nm	3.60E+06	nm	6.30E-01	n	2.60E+00	n			
													Polynuclear Aromatic Hydrocarbons (PAHs)													
				6.00E-02	I			V		1	0.13		~Acenaphthene	83-32-9	3.60E+03	n	4.50E+04	n					5.30E+02	n	

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	GIABS	ABS_d	C_{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
				3.00E-01	I			V		1	0.13		~Anthracene	120-12-7	1.80E+04	n	2.30E+05	nm					1.80E+03	n	
1.00E-01	E	6.00E-05	E					V	M	1	0.13		~Benz[a]anthracene	56-55-3	1.10E+02	c	2.10E+03	c	1.70E+00	c	2.00E+01	c	3.00E+00	c	
				9.00E-05	X	2.00E-06	X			1	0.1		~Benzo(e)pyrene	192-97-2	5.70E+00	n	7.30E+01	n	2.10E-03	n	8.80E-03	n	1.80E+00	n	
1.20E+00	C	1.10E-04	C							1	0.13		~Benzo(j)fluoranthene	205-82-3	4.20E+01	c	1.80E+02	c	2.60E+00	c	1.10E+01	c	6.50E+00	c	
1.00E+00	I	6.00E-04	I	3.00E-04	I	2.00E-06	I		M	1	0.13		~Benzo[a]pyrene	50-32-8	1.10E+01	c**	2.10E+02	c**	2.10E-03	n	8.80E-03	n	2.50E+00	c**	2.00E-01
1.00E-01	E	6.00E-05	E						M	1	0.13		~Benzo[b]fluoranthene	205-99-2	1.10E+02	c	2.10E+03	c	1.70E+00	c	2.00E+01	c	2.50E+01	c	
1.00E-02	E	6.00E-06	E						M	1	0.13		~Benzo[k]fluoranthene	207-08-9	1.10E+03	c	2.10E+04	c	1.70E+01	c	2.00E+02	c	2.50E+02	c	
				8.00E-02	I			V		1	0.13		~Chloronaphthalene, Beta-	91-58-7	4.80E+03	n	6.00E+04	n					7.50E+02	n	
1.00E-03	E	6.00E-07	E						M	1	0.13		~Chrysene	218-01-9	1.10E+04	c	2.10E+05	cm	1.70E+02	c	2.00E+03	c	2.50E+03	c	
1.00E+00	E	6.00E-04	E						M	1	0.13		~Dibenz[a,h]anthracene	53-70-3	1.10E+01	c	2.10E+02	c	1.70E-01	c	2.00E+00	c	2.50E+00	c	
1.20E+01	C	1.10E-03	C							1	0.13		~Dibenzo(a,e)pyrene	192-65-4	4.20E+00	c	1.80E+01	c	2.60E-01	c	1.10E+00	c	6.50E-01	c	
2.50E+02	C	7.10E-02	C						M	1	0.13		~Dimethylbenz(a)anthracene, 7,12-	57-97-6	4.60E-02	c	8.40E-01	c	1.40E-03	c	1.70E-02	c	1.00E-02	c	
				4.00E-02	I					1	0.13		~Fluoranthene	206-44-0	2.40E+03	n	3.00E+04	n					8.00E+02	n	
				4.00E-02	I			V		1	0.13		~Fluorene	86-73-7	2.40E+03	n	3.00E+04	n					2.90E+02	n	
1.00E-01	E	6.00E-05	E						M	1	0.13		~Indeno[1,2,3-cd]pyrene	193-39-5	1.10E+02	c	2.10E+03	c	1.70E+00	c	2.00E+01	c	2.50E+01	c	
2.90E-02	P			7.00E-02	A			V		1	0.13	3.94E+02	~Methylnaphthalene, 1-	90-12-0	1.80E+03	c**s	7.30E+03	c**s					1.10E+02	c**	
				4.00E-03	I			V		1	0.13		~Methylnaphthalene, 2-	91-57-6	2.40E+02	n	3.00E+03	n					3.60E+01	n	
1.20E-01	C	3.40E-05	C	2.00E-02	I	3.00E-03	I	V		1	0.13		~Naphthalene	91-20-3	1.30E+02	n	5.90E+02	n	3.10E+00	n	1.30E+01	n	6.10E+00	n	
1.20E+00	C	1.10E-04	C							1	0.13		~Nitropyrene, 4-	57835-92-4	4.20E+01	c	1.80E+02	c	2.60E+00	c	1.10E+01	c	1.90E+00	c	
				3.00E-02	I			V		1	0.13		~Pyrene	129-00-0	1.80E+03	n	2.30E+04	n					1.20E+02	n	
1.50E-01	I			9.00E-03	I					1	0.1		Prochloraz	67747-09-5	3.60E+02	c**	1.50E+03	c**					3.80E+01	c**	
				6.00E-03	H			V		1			Profluralin	26399-36-0	4.70E+02	n	7.00E+03	n					2.60E+01	n	
				1.50E-02	I					1	0.1		Prometon	1610-18-0	9.50E+02	n	1.20E+04	n					2.50E+02	n	
				4.00E-02	O					1	0.1		Prometryn	7287-19-6	2.50E+03	n	3.30E+04	n					6.00E+02	n	
				7.50E-02	I					1	0.1		Pronamide	23950-58-5	4.70E+03	n	6.20E+04	n					1.20E+03	n	
				1.30E-02	I					1	0.1		Propachlor	1918-16-7	8.20E+02	n	1.10E+04	n					2.50E+02	n	
				5.00E-03	I					1	0.1		Propanil	709-98-8	3.20E+02	n	4.10E+03	n					8.20E+01	n	
1.90E-01	O			4.00E-02	O					1	0.1		Propargite	2312-35-8	2.80E+02	c**	1.20E+03	c*					1.60E+01	c*	
				2.00E-03	I			V		1		1.11E+05	Propargyl Alcohol	107-19-7	1.60E+02	n	2.30E+03	n					4.00E+01	n	
				2.00E-02	I					1	0.1		Propazine	139-40-2	1.30E+03	n	1.60E+04	n					3.40E+02	n	
				2.00E-02	I					1	0.1		Propham	122-42-9	1.30E+03	n	1.60E+04	n					3.50E+02	n	
				1.00E-01	O					1	0.1		Propiconazole	60207-90-1	6.30E+03	n	8.20E+04	n					1.60E+03	n	
						8.00E-03	I	V		1		3.26E+04	Propionaldehyde	123-38-6	7.50E+01	n	3.10E+02	n	8.30E+00	n	3.50E+01	n	1.70E+01	n	
				1.00E-01	X	1.00E+00	X	V		1		2.64E+02	Propyl benzene	103-65-1	3.80E+03	ns	2.40E+04	ns	1.00E+03	n	4.40E+03	n	6.60E+02	n	
						3.00E+00	C	V		1		3.49E+02	Propylene	115-07-1	2.20E+03	ns	9.30E+03	ns	3.10E+03	n	1.30E+04	n	6.30E+03	n	
				2.00E+01	P					1	0.1		Propylene Glycol	57-55-6	1.30E+06	nm	1.60E+07	nm					4.00E+05	n	
						2.70E-04	A			1	0.1		Propylene Glycol Dinitrate	6423-43-4	3.90E+05	nm	1.60E+06	nm	2.80E-01	n	1.20E+00	n			
				7.00E-01	H	2.00E+00	I	V		1		1.06E+05	Propylene Glycol Monomethyl Ether	107-98-2	4.10E+04	n	3.70E+05	nms	2.10E+03	n	8.80E+03	n	3.20E+03	n	
2.40E-01	I	3.70E-06	I			3.00E-02	I	V		1		7.77E+04	Propylene Oxide	75-56-9	2.10E+02	c**	9.70E+02	c**	3.10E+01	n	1.30E+02	n	2.70E+01	c**	
				1.00E-03	I			V		1		5.30E+05	Pyridine	110-86-1	7.80E+01	n	1.20E+03	n					2.00E+01	n	
				5.00E-04	I					1	0.1		Quinalphos	13593-03-8	3.20E+01	n	4.10E+02	n					5.10E+00	n	
3.00E+00	I									1	0.1		Quinoline	91-22-5	1.80E+01	c	7.70E+01	c					2.40E+00	c	
				9.00E-03	I					1	0.1		Quizalofop-ethyl	76578-14-8	5.70E+02	n	7.40E+03	n					1.20E+02	n	
						3.00E+04	A			1			Refractory Ceramic Fibers (units in fibers)	E715557					3.10E+04	G	1.30E+05	G			
				3.00E-02	I					1	0.1		Resmethrin	10453-86-8	1.90E+03	n	2.50E+04	n					6.70E+01	n	
				5.00E-02	H			V		1			Ronnel	299-84-3	3.90E+03	n	5.80E+04	n					4.10E+02	n	
				4.00E-03	I					1	0.1		Rotenone	83-79-4	2.50E+02	n	3.30E+03	n					6.10E+01	n	
2.20E-01	C	6.30E-05	C						M	1	0.1		Safrole	94-59-7	5.50E+01	c	1.00E+03	c	1.60E+00	c	1.90E+01	c	9.60E+00	c	
				5.00E-03	I					1			Selenious Acid	7783-00-8	3.90E+02	n	5.80E+03	n					1.00E+02	n	
				5.00E-03	I	2.00E-02	C			1			Selenium	7782-49-2	3.90E+02	n	5.80E+03	n	2.10E+01	n	8.80E+01	n	1.00E+02	n	5.00E+01
				5.00E-03	C	2.00E-02	C			1			Selenium Sulfide	7446-34-6	3.90E+02	n	5.80E+03	n	2.10E+01	n	8.80E+01	n	1.00E+02	n	
				1.40E-01	O					1	0.1		Sethoxydim	74051-80-2	8.80E+03	n	1.10E+05	nm					1.60E+03	n	
						3.00E-03	C			1			Silica (crystalline, respirable)	7631-86-9	4.30E+06	nm	1.80E+07	nm	3.10E+00	n	1.30E+01	n			
				5.00E-03	I					0.04			Silver	7440-22-4	3.90E+02	n	5.80E+03	n					9.40E+01	n	
1.20E-01	H			5.00E-03	I					1	0.1		Simazine	122-34-9	3.20E+02	n	1.90E+03	c**					6.10E+01	c**	4.00E+00
				1.30E-02	I					1	0.1		Sodium Acifluorfen	62476-59-9	8.20E+02	n	1.10E+04	n					2.60E+02	n	
				4.00E-03	I					1			Sodium Azide	26628-22-8	3.10E+02	n	4.70E+03	n					8.00E+01	n	
2.70E-01	H			3.00E-02	I					1	0.1		Sodium Diethyldithiocarbamate	148-18-5	2.00E+02	c**	8.50E+02	c*					2.90E+01	c*	
				5.00E-02	A	1.30E-02	C			1			Sodium Fluoride	7681-49-4	3.90E+03	n	5.80E+04	n	1.40E+01	n	5.70E+01	n	1.00E+03	n	4.00E+03
				2.00E-05	I					1	0.1		Sodium Fluoroacetate	62-74-8	1.30E+00	n	1.60E+01	n					4.00E-01	n	
				1.00E-03	H					1			Sodium Metavanadate	13718-26-8	7.80E+01	n	1.20E+03	n					2.00E+01	n	
				8.00E-04	P					1			Sodium Tungstate	13472-45-2	6.30E+01	n	9.30E+02	n					1.60E+01	n	
				8.00E-04	P					1			Sodium Tungstate Dihydrate	10213-10-2	6.30E+01	n	9.30E+02	n					1.60E+01	n	
2.40E-02	H			3.00E-02	I					1	0.1		Stirofos (Tetrachlorovinphos)	961-11-5	1.90E+03	n	9.60E+03	c**					2.80E+02	c**	
				6.00E-01	I					1			Strontium, Stable	7440-24-6	4.70E+04	n	7.00E+05	nm					1.20E+04	n	
				3.00E-04	I					1	0.1		Strychnine	57-24-9	1.90E+01	n	2.50E+02	n					5.90E+00	n	
				2.00E-01	I	1.00E+00	I	V		1		8.67E+02	Styrene	100-42-5	6.00E+03	ns	3.50E+04	ns	1.00E+03	n	4.40E+03	n	1.20E+03	n	1.00E+02
				3.00E-03	P					1	0.1		Styrene-Acrylonitrile (SAN) Trimer (THNA isomer)	57964-39-3	1.90E+02	n	2.50E+03	n					4.80E+01	n	
				3.00E-03	P					1	0.1		Styrene-Acrylonitrile (SAN) Trimer (THNP isomer)	57964-40-6	1.90E+02	n	2.50E+03	n					4.80E+01	n	
				1.00E-03	P	2.00E-03	X			1	0.1		Sulfolane	126-33-0	6.30E+01	n	8.20E+02	n	2.10E+00	n	8.80E+00	n	2.00E+01	n	
				8.00E-04	P					1	0.1		Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9	5.10E+01	n	6.60E+02	n					1.10E+01	n	

RMLS
TR=1E-04
THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	GIABS	ABS_d	C_{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
						1.00E-03	C	V		1			Sulfur Trioxide	7446-11-9	1.40E+06	nm	6.00E+06	nm	1.00E+00	n	4.40E+00	n	2.10E+00	n	
						1.00E-03	C			1			Sulfuric Acid	7664-93-9	1.40E+06	nm	6.00E+06	nm	1.00E+00	n	4.40E+00	n			
2.50E-02	I	7.10E-06	I	5.00E-02	H					1	0.1		Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl ester	140-57-8	2.20E+03	c**	9.20E+03	c**	4.00E+01	c	1.70E+02	c	1.30E+02	c**	
				3.00E-02	H					1	0.1		TCMTB	21564-17-0	1.90E+03	n	2.50E+04	n					4.80E+02	n	
				7.00E-02	I					1	0.1		Tebuthiuron	34014-18-1	4.40E+03	n	5.70E+04	n					1.40E+03	n	
				2.00E-02	H					1	0.1		Temephos	3383-96-8	1.30E+03	n	1.60E+04	n					4.00E+02	n	
				1.30E-02	I					1	0.1		Terbacil	5902-51-2	8.20E+02	n	1.10E+04	n					2.50E+02	n	
				2.50E-05	H			V		1		3.09E+01	Terbufos	13071-79-9	2.00E+00	n	2.90E+01	n					2.40E-01	n	
				1.00E-03	I					1	0.1		Terbutryn	886-50-0	6.30E+01	n	8.20E+02	n					1.30E+01	n	
5.00E-03	C	1.30E-06	C					V		1			Tert-Butyl Acetate	540-88-5	8.10E+02	c	3.60E+03	c	2.20E+02	c	9.40E+02	c	3.30E+02	c	
				1.00E-04	I					1	0.1		Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1	6.30E+00	n	8.20E+01	n					2.00E+00	n	
				3.00E-05	P			V		1			Tetrachlorobenzene, 1,2,4,5-	95-94-3	2.30E+00	n	3.50E+01	n					1.70E-01	n	
2.60E-02	I	7.40E-06	I	3.00E-02	I			V		1		6.80E+02	Tetrachloroethane, 1,1,1,2-	630-20-6	2.00E+02	c*	8.80E+02	c*s	3.80E+01	c	1.70E+02	c	5.70E+01	c**	
2.00E-01	I	5.80E-05	C	2.00E-02	I			V		1		1.90E+03	Tetrachloroethane, 1,1,2,2-	79-34-5	6.00E+01	c*	2.70E+02	c*	4.80E+00	c	2.10E+01	c	7.60E+00	c*	
2.10E-03	I	2.60E-07	I	6.00E-03	I	4.00E-02	I	V		1		1.66E+02	Tetrachloroethylene	127-18-4	8.10E+01	n	3.90E+02	ns	4.20E+01	n	1.80E+02	n	4.10E+01	n	5.00E+00
				3.00E-02	I					1	0.1		Tetrachlorophenol, 2,3,4,6-	58-90-2	1.90E+03	n	2.50E+04	n					2.40E+02	n	
1.60E+01	X			6.00E-05	X			V		1			Tetrachlorotoluene, p- alpha, alpha, alpha-	5216-25-1	4.30E+00	c**	2.00E+01	c**					1.70E-01	c**	
				5.00E-04	I					1	0.1		Tetraethyl Dithiopyrophosphate	3689-24-5	3.20E+01	n	4.10E+02	n					7.10E+00	n	
						8.00E+01	I	V		1		2.05E+03	Tetrafluoroethane, 1,1,1,2-	811-97-2	1.00E+05	nms	4.30E+05	nms	8.30E+04	n	3.50E+05	n	1.70E+05	n	
				1.00E-04	X					1	0.1		Tetramethylphosphoramide, -N,N,N',N" (TMPA)	16853-36-4	6.30E+00	n	8.20E+01	n					2.00E+00	n	
				2.00E-03	P					1	0.00065		Tetryl (Trinitrophenylmethylnitramine)	479-45-8	1.60E+02	n	2.30E+03	n					3.90E+01	n	
				2.00E-05	G					1			Thallic Oxide	1314-32-5	1.60E+00	n	2.30E+01	n					4.00E-01	n	
				1.00E-05	X					1			Thallium (I) Nitrate	10102-45-1	7.80E-01	n	1.20E+01	n					2.00E-01	n	
				1.00E-05	X					1			Thallium (Soluble Salts)	7440-28-0	7.80E-01	n	1.20E+01	n					2.00E-01	n	2.00E+00
				1.00E-05	X			V		1			Thallium Acetate	563-68-8	7.80E-01	n	1.20E+01	n					2.00E-01	n	
				2.00E-05	X					1	0.1		Thallium Carbonate	6533-73-9	1.30E+00	n	1.60E+01	n					4.00E-01	n	
				1.00E-05	X					1			Thallium Chloride	7791-12-0	7.80E-01	n	1.20E+01	n					2.00E-01	n	
				1.00E-05	G					1			Thallium Selenite	12039-52-0	7.80E-01	n	1.20E+01	n					2.00E-01	n	
				2.00E-05	X					1			Thallium Sulfate	7446-18-6	1.60E+00	n	2.30E+01	n					4.00E-01	n	
				4.30E-02	O					1	0.1		Thifensulfuron-methyl	79277-27-3	2.70E+03	n	3.50E+04	n					8.60E+02	n	
				1.00E-02	I					1	0.1		Thiobencarb	28249-77-6	6.30E+02	n	8.20E+03	n					1.60E+02	n	
				7.00E-02	X					1	0.0075		Thiodiglycol	111-48-8	5.40E+03	n	7.90E+04	n					1.40E+03	n	
				3.00E-04	H					1	0.1		Thiofanox	39196-18-4	1.90E+01	n	2.50E+02	n					5.30E+00	n	
1.20E-02	O			1.60E-01	O					1	0.1		Thiophanate, Methyl	23564-05-8	4.70E+03	c**	2.00E+04	c**					6.70E+02	c**	
				1.50E-02	O					1	0.1		Thiram	137-26-8	9.50E+02	n	1.20E+04	n					2.90E+02	n	
				6.00E-01	H					1			Tin	7440-31-5	4.70E+04	n	7.00E+05	nm					1.20E+04	n	
						1.00E-04	A	V		1			Titanium Tetrachloride	7550-45-0	1.40E+05	nm	6.00E+05	nm	1.00E-01	n	4.40E-01	n	2.10E-01	n	
				8.00E-02	I	5.00E+00	I	V		1		8.18E+02	Toluene	108-88-3	4.90E+03	ns	4.70E+04	ns	5.20E+03	n	2.20E+04	n	1.10E+03	n	1.00E+03
3.90E-02	C	1.10E-05	C			8.00E-06	C	V		1			Toluene-2,4-diisocyanate	584-84-9	6.40E+00	n	2.70E+01	n	8.30E-03	n	3.50E-02	n	1.70E-02	n	
1.80E-01	X			2.00E-04	X					1	0.1		Toluene-2,5-diamine	95-70-5	1.30E+01	n	1.60E+02	n					4.00E+00	n	
3.90E-02	C	1.10E-05	C			8.00E-06	C	V		1		1.71E+03	Toluene-2,6-diisocyanate	91-08-7	5.30E+00	n	2.20E+01	n	8.30E-03	n	3.50E-02	n	1.70E-02	n	
				1.00E-04	X					1	0.1		Toluenediamine, 2,3-	2687-25-4	6.30E+00	n	8.20E+01	n					2.00E+00	n	
				1.00E-04	X					1	0.1		Toluenediamine, 3,4-	496-72-0	6.30E+00	n	8.20E+01	n					2.00E+00	n	
				5.00E-03	P					1	0.1		Toluic Acid, p-	99-94-5	3.20E+02	n	4.10E+03	n					9.00E+01	n	
1.60E-02	P	5.10E-05	C							1	0.1		Toluidine, o- (Methylaniline, 2-)	95-53-4	3.40E+03	c	1.40E+04	c	5.50E+00	c	2.40E+01	c	4.70E+02	c	
3.00E-02	P			4.00E-03	X					1	0.1		Toluidine, p-	106-49-0	2.50E+02	n	3.30E+03	n					7.70E+01	n	
				3.00E+00	P			V		1		3.42E-01	Total Petroleum Hydrocarbons (Aliphatic High)	E1790670	2.30E+05	nms	3.50E+06	nms					6.00E+04	n	
						6.00E-01	P	V		1		1.41E+02	Total Petroleum Hydrocarbons (Aliphatic Low)	E1790666	5.20E+02	ns	2.20E+03	ns	6.30E+02	n	2.60E+03	n	1.30E+03	n	
				1.00E-02	X	1.00E-01	P	V		1		6.86E+00	Total Petroleum Hydrocarbons (Aliphatic Medium)	E1790668	9.60E+01	ns	4.40E+02	ns	1.00E+02	n	4.40E+02	n	1.00E+02	n	
				4.00E-02	P					1	0.13		Total Petroleum Hydrocarbons (Aromatic High)	E1790676	2.40E+03	n	3.00E+04	n					8.00E+02	n	
				4.00E-03	P	3.00E-02	P	V		1		1.82E+03	Total Petroleum Hydrocarbons (Aromatic Low)	E1790672	8.20E+01	n	4.20E+02	n	3.10E+01	n	1.30E+02	n	3.30E+01	n	
				4.00E-03	P	3.00E-03	P	V		1	0.13		Total Petroleum Hydrocarbons (Aromatic Medium)	E1790674	9.70E+01	n	5.60E+02	n	3.10E+00	n	1.30E+01	n	5.50E+00	n	
1.10E+00	I	3.20E-04	I	9.00E-05	P					1	0.1		Toxaphene	8001-35-2	5.70E+00	n	7.40E+01	n	8.80E-01	c	3.80E+00	c	1.80E+00	n	3.00E+00
				3.00E-05	X					1	0.1		Toxaphene, Weathered	E1841606	1.90E+00	n	2.50E+01	n					6.00E-01	n	
				7.50E-03	I					1	0.1		Tralomethrin	66841-25-6	4.70E+02	n	6.20E+03	n					1.50E+02	n	
				3.00E-04	A			V		1			Tri-n-butyltin	688-73-3	2.30E+01	n	3.50E+02	n					3.70E+00	n	
				8.00E+01	X					1	0.1		Triacetin	102-76-1	5.10E+06	nm	6.60E+07	nm					1.60E+06	n	
				3.40E-02	O					1	0.1		Triadimefon	43121-43-3	2.10E+03	n	2.80E+04	n					6.30E+02	n	
7.20E-02	O			2.50E-02	O			V		1			Triallate	2303-17-5	9.70E+02	c**	4.60E+03	c**					4.70E+01	c**	
				1.00E-02	I					1	0.1		Triasulfuron	82097-50-5	6.30E+02	n	8.20E+03	n					2.00E+02	n	
				8.00E-03	I					1	0.1		Tribenuron-methyl	101200-48-0	5.10E+02	n	6.60E+03	n					1.60E+02	n	
				5.00E-03	I			V		1			Tribromobenzene, 1,2,4-	615-54-3	3.90E+02	n	5.80E+03	n					4.50E+01	n	
				9.00E-03	X					1	0.1		Tribromophenol, 2,4,6-	118-79-6	5.70E+02	n	7.40E+03	n					1.20E+02	n	
				2.00E-04	O					1	0.1		Tribufos	78-48-8	1.30E+01	n	1.60E+02	n					5.70E-01	n	
9.00E-03	P			1.00E-02	P					1	0.1		Tributyl Phosphate	126-73-8	6.30E+02	n	8.20E+03	n					1.20E+02	n	
				3.00E-04	P					1	0.1		Tributyltin Compounds	E1790679	1.90E+01	n	2.50E+02	n					6.00E+00	n	
				3.00E-04	I					1	0.1		Tributyltin Oxide	56-35-9	1.90E+01	n	2.50E+02	n					5.70E+00	n	
										1	0.1		Trichloramine	10025-85-1											4.0E+03(G)
				3.00E+01	I	5.00E+00	P	V		1		9.10E+02	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1	6.70E+03	ns	2.80E+04	ns	5.20E+03	n	2.20E+04	n	1.00E+04	n	
7.00E-02	I			2.00E-02	I					1	0.1		Trichloroacetic Acid	76-03-9	7.80E+02	c**	3.30E+03	c**					1.10E+02	c**	6.0E+01(G)
2.90E-02	H									1	0.1		Trichloroaniline HCl, 2,4,6-	33663-50-2	1.90E+03	c	7.90E+03	c					2.70E+02	c	

RMLs TR=1E-04 HQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information													Contaminant		Removal Levels										
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	GIABS	ABS_d	C_{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Resident Air (ug/m^3)	key	Industrial Air (ug/m^3)	key	Tapwater (ug/L)	key	MCL (ug/L)
7.00E-03	X			3.00E-05	X					1	0.1		Trichloroaniline, 2,4,6-	634-93-5	1.90E+00	n	2.50E+01	n					4.00E-01	n	
				8.00E-04	X			V		1			Trichlorobenzene, 1,2,3-	87-61-6	6.30E+01	n	9.30E+02	n					7.00E+00	n	
2.90E-02	P			1.00E-02	I	2.00E-03	P	V		1		4.04E+02	Trichlorobenzene, 1,2,4-	120-82-1	5.80E+01	n	2.60E+02	n	2.10E+00	n	8.80E+00	n	4.00E+00	n	7.00E+01
				2.00E+00	I	5.00E+00	I	V		1		6.40E+02	Trichloroethane, 1,1,1-	71-55-6	8.10E+03	ns	3.60E+04	ns	5.20E+03	n	2.20E+04	n	8.00E+03	n	2.00E+02
5.70E-02	I	1.60E-05	I	4.00E-03	I	2.00E-04	X	V		1		2.16E+03	Trichloroethane, 1,1,2-	79-00-5	1.50E+00	n	6.30E+00	n	2.10E-01	n	8.80E-01	n	4.10E-01	n	5.00E+00
4.60E-02	I	4.10E-06	I	5.00E-04	I	2.00E-03	I	V	M	1		6.92E+02	Trichloroethylene	79-01-6	4.10E+00	n	1.90E+01	n	2.10E+00	n	8.80E+00	n	2.80E+00	n	5.00E+00
				3.00E-01	I			V		1		1.23E+03	Trichlorofluoromethane	75-69-4	2.30E+04	ns	3.50E+05	nms					5.20E+03	n	
				1.00E-01	I					1	0.1		Trichlorophenol, 2,4,5-	95-95-4	6.30E+03	n	8.20E+04	n					1.20E+03	n	
1.10E-02	I	3.10E-06	I	1.00E-03	P					1	0.1		Trichlorophenol, 2,4,6-	88-06-2	6.30E+01	n	8.20E+02	n	9.10E+01	c	4.00E+02	c	1.20E+01	n	
				1.00E-02	I					1	0.1		Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5	6.30E+02	n	8.20E+03	n					1.60E+02	n	
				8.00E-03	I					1	0.1		Trichlorophenoxypropionic acid, -2,4,5	93-72-1	5.10E+02	n	6.60E+03	n					1.10E+02	n	5.00E+01
				5.00E-03	I			V		1		1.28E+03	Trichloropropane, 1,1,2-	598-77-6	3.90E+02	n	5.80E+03	ns					8.80E+01	n	
3.00E+01	I			4.00E-03	I	3.00E-04	I	V	M	1		1.40E+03	Trichloropropane, 1,2,3-	96-18-4	5.10E-01	c**	1.10E+01	c**	3.10E-01	n	1.30E+00	n	7.50E-02	c**	
				3.00E-03	X	3.00E-04	P	V		1		3.11E+02	Trichloropropene, 1,2,3-	96-19-5	7.30E-01	n	3.10E+00	n	3.10E-01	n	1.30E+00	n	6.20E-01	n	
				2.00E-02	A					1	0.1		Tricresyl Phosphate (TCP)	1330-78-5	1.30E+03	n	1.60E+04	n					1.60E+02	n	
				3.00E-03	I					1	0.1		Tridiphane	58138-08-2	1.90E+02	n	2.50E+03	n					1.80E+01	n	
						7.00E-03	I	V		1		2.79E+04	Triethylamine	121-44-8	1.20E+02	n	4.80E+02	n	7.30E+00	n	3.10E+01	n	1.50E+01	n	
				2.00E+00	P					1	0.1		Triethylene Glycol	112-27-6	1.30E+05	nm	1.60E+06	nm					4.00E+04	n	
						2.00E+01	P	V		1		4.81E+03	Trifluoroethane, 1,1,1-	420-46-2	1.50E+04	ns	6.20E+04	ns	2.10E+04	n	8.80E+04	n	4.20E+04	n	
7.70E-03	I			7.50E-03	I			V		1			Trifluralin	1582-09-8	5.90E+02	n	8.80E+03	n					4.00E+01	n	
2.00E-02	P			1.00E-02	P					1	0.1		Trimethyl Phosphate	512-56-1	6.30E+02	n	8.20E+03	n					2.00E+02	n	
				1.00E-02	I	6.00E-02	I	V		1		2.93E+02	Trimethylbenzene, 1,2,3-	526-73-8	3.40E+02	ns	2.00E+03	ns	6.30E+01	n	2.60E+02	n	5.50E+01	n	
				1.00E-02	I	6.00E-02	I	V		1		2.19E+02	Trimethylbenzene, 1,2,4-	95-63-6	3.00E+02	ns	1.80E+03	ns	6.30E+01	n	2.60E+02	n	5.60E+01	n	
				1.00E-02	I	6.00E-02	I	V		1		1.82E+02	Trimethylbenzene, 1,3,5-	108-67-8	2.70E+02	ns	1.50E+03	ns	6.30E+01	n	2.60E+02	n	6.00E+01	n	
				1.00E-02	X			V		1		2.96E+01	Trimethylpentene, 2,4,4-	25167-70-8	7.80E+02	ns	1.20E+04	ns					3.80E+01	n	
				3.00E-02	I					1	0.019		Trinitrobenzene, 1,3,5-	99-35-4	2.20E+03	n	3.20E+04	n					5.90E+02	n	
3.00E-02	I			5.00E-04	I					1	0.032		Trinitrotoluene, 2,4,6-	118-96-7	3.60E+01	n	5.10E+02	n					9.80E+00	n	
				2.00E-02	P					1	0.1		Triphenylphosphine Oxide	791-28-6	1.30E+03	n	1.60E+04	n					3.60E+02	n	
				2.00E-02	A					1	0.1		Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8	1.30E+03	n	1.60E+04	n					3.60E+02	n	
				1.00E-02	X					1	0.1		Tris(1-chloro-2-propyl)phosphate	13674-84-5	6.30E+02	n	8.20E+03	n					1.90E+02	n	
2.30E+00	C	6.60E-04	C					V		1		4.67E+02	Tris(2,3-dibromopropyl)phosphate	126-72-7	2.80E+01	c	1.30E+02	c	4.30E-01	c	1.90E+00	c	6.80E-01	c	
2.00E-02	P			7.00E-03	P					1	0.1		Tris(2-chloroethyl)phosphate	115-96-8	4.40E+02	n	5.70E+03	n					1.40E+02	n	
3.20E-03	P			1.00E-01	P					1	0.1		Tris(2-ethylhexyl)phosphate	78-42-2	6.30E+03	n	7.20E+04	c**					2.00E+03	n	
				8.00E-04	P					1			Tungsten	7440-33-7	6.30E+01	n	9.30E+02	n					1.60E+01	n	
				2.00E-04	A	4.00E-05	A			1			Uranium	7440-61-1	1.60E+01	n	2.30E+02	n	4.20E-02	n	1.80E-01	n	4.00E+00	n	3.00E+01
1.00E+00	C	2.90E-04	C						M	1	0.1		Urethane	51-79-6	1.20E+01	c	2.30E+02	c	3.50E-01	c	4.20E+00	c	2.50E+00	c	
		8.30E-03	P	9.00E-03	I	7.00E-06	P			0.026			Vanadium Pentoxide	1314-62-1	6.60E+02	n	8.40E+03	n	7.30E-03	n	3.10E-02	n	1.50E+02	n	
				5.00E-03	G	1.00E-04	A			0.026			Vanadium and Compounds	7440-62-2	3.90E+02	n	5.80E+03	n	1.00E-01	n	4.40E-01	n	8.60E+01	n	
				1.00E-03	I			V		1			Vernolate	1929-77-7	7.80E+01	n	1.20E+03	n					1.10E+01	n	
				1.20E-03	O					1	0.1		Vinclozolin	50471-44-8	7.60E+01	n	9.80E+02	n					2.10E+01	n	
				1.00E+00	H	2.00E-01	I	V		1		2.75E+03	Vinyl Acetate	108-05-4	9.10E+02	n	3.80E+03	ns	2.10E+02	n	8.80E+02	n	4.10E+02	n	
		1.50E-05	P			3.00E-03	I	V		1		2.47E+03	Vinyl Bromide	593-60-2	4.30E+00	n	1.80E+01	n	3.10E+00	n	1.30E+01	n	6.30E+00	n	
7.20E-01	I	4.40E-06	I	3.00E-03	I	8.00E-02	A	V	M	1		3.92E+03	Vinyl Chloride	75-01-4	5.90E+00	c*	1.70E+02	c**	1.70E+01	c**	2.80E+02	c**	1.90E+00	c*	2.00E+00
				3.00E-04	I					1	0.1		Warfarin	81-81-2	1.90E+01	n	2.50E+02	n					5.60E+00	n	
				2.00E-01	G	1.00E-01	G	V		1		3.88E+02	Xylene, m-	108-38-3	5.50E+02	ns	2.40E+03	ns	1.00E+02	n	4.40E+02	n	1.90E+02	n	
				2.00E-01	G	1.00E-01	G	V		1		4.34E+02	Xylene, o-	95-47-6	6.40E+02	ns	2.80E+03	ns	1.00E+02	n	4.40E+02	n	1.90E+02	n	
				2.00E-01	G	1.00E-01	G	V		1		3.90E+02	Xylene, p-	106-42-3	5.60E+02	ns	2.40E+03	ns	1.00E+02	n	4.40E+02	n	1.90E+02	n	
				2.00E-01	I	1.00E-01	I	V		1		2.60E+02	Xylenes	1330-20-7	5.80E+02	ns	2.50E+03	ns	1.00E+02	n	4.40E+02	n	1.90E+02	n	1.00E+04
				3.00E-04	I					1			Zinc Phosphide	1314-84-7	2.30E+01	n	3.50E+02	n					6.00E+00	n	
				3.00E-01	I					1			Zinc and Compounds	7440-66-6	2.30E+04	n	3.50E+05	nm					6.00E+03	n	
				5.00E-02	I					1	0.1		Zineb	12122-67-7	3.20E+03	n	4.10E+04	n					9.90E+02	n	
				8.00E-05	X					1			Zirconium	7440-67-7	6.30E+00	n	9.30E+01	n					1.60E+00	n	

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
				3.0E-04	O						1.36E+09		1	0.1	Acephate	30560-19-1					2.3E+01	9.9E+01		1.9E+01
		2.2E-06	I			9.0E-03	I	V		1.07E+05	1.36E+09	8.72E+03	1		Acetaldehyde	75-07-0			1.1E+03	1.1E+03			8.2E+01	8.2E+01
				2.0E-02	I						1.36E+09		1	0.1	Acetochlor	34256-82-1					1.6E+03	6.6E+03		1.3E+03
				9.0E-01	I			V		1.14E+05	1.36E+09	1.37E+04	1		Acetone	67-64-1					7.0E+04			7.0E+04
						2.0E-03	X				1.36E+09		1	0.1	Acetone Cyanohydrin	75-86-5							2.8E+06	2.8E+06
						6.0E-02	I	V		1.28E+05	1.36E+09	1.30E+04	1		Acetonitrile	75-05-8							8.1E+02	8.1E+02
				1.0E-01	I			V		2.52E+03	1.36E+09	5.97E+04	1		Acetophenone	98-86-2					7.8E+03			7.8E+03
3.8E+00	C	1.3E-03	C								1.36E+09		1	0.1	Acetylaminofluorene, 2-	53-96-3	1.8E+01	6.5E+01	2.9E+05	1.4E+01				
				5.0E-04	I	2.0E-05	I	V		2.27E+04	1.36E+09	6.91E+03	1		Acrolein	107-02-8					3.9E+01		1.4E-01	1.4E-01
5.0E-01	I	1.0E-04	I	2.0E-03	I	6.0E-03	I		M		1.36E+09		1	0.1	Acrylamide	79-06-1	3.1E+01	1.2E+02	1.4E+06	2.4E+01	1.6E+02	6.6E+02	8.5E+06	1.3E+02
				5.0E-01	I	2.0E-04	P	V		1.09E+05	1.36E+09	9.53E+04	1		Acrylic Acid	79-10-7					3.9E+04		2.0E+01	2.0E+01
5.4E-01	I	6.8E-05	I	1.0E-02	A	2.0E-03	I	V		1.13E+04	1.36E+09	7.69E+03	1		Acrylonitrile	107-13-1	1.3E+02		3.2E+01	2.5E+01	7.8E+02		1.6E+01	1.6E+01
						6.0E-03	P				1.36E+09		1	0.1	Adiponitrile	111-69-3							8.5E+06	8.5E+06
5.6E-02	C			1.0E-02	I						1.36E+09		1	0.1	Alachlor	15972-60-8	1.2E+03	4.4E+03		9.7E+02	7.8E+02	3.3E+03		6.3E+02
				1.0E-03	I						1.36E+09		1	0.1	Aldicarb	116-06-3					7.8E+01	3.3E+02		6.3E+01
				1.0E-03	I						1.36E+09		1	0.1	Aldicarb Sulfone	1646-88-4					7.8E+01	3.3E+02		6.3E+01
											1.36E+09		1	0.1	Aldicarb sulfoxide	1646-87-3								
1.7E+01	I	4.9E-03	I	3.0E-05	I			V			1.36E+09	1.72E+06	1		Aldrin	309-00-2	4.1E+00		9.8E+01	3.9E+00	2.3E+00			2.3E+00
				4.0E-03	P	1.0E-04	X	V		1.11E+05	1.36E+09	3.42E+04	1		Allyl Alcohol	107-18-6					3.1E+02		3.6E+00	3.5E+00
2.1E-02	C	6.0E-06	C			1.0E-03	I	V		1.42E+03	1.36E+09	1.58E+03	1		Allyl Chloride	107-05-1	3.3E+03		7.4E+01	7.2E+01			1.7E+00	1.7E+00
				1.0E+00	P	5.0E-03	P				1.36E+09		1		Aluminum	7429-90-5					7.8E+04		7.1E+06	7.7E+04
				4.0E-04	I						1.36E+09		1		Aluminum Phosphide	20859-73-8					3.1E+01			3.1E+01
				9.0E-03	I						1.36E+09		1	0.1	Ametryn	834-12-8					7.0E+02	3.0E+03		5.7E+02
2.1E+01	C	6.0E-03	C								1.36E+09		1	0.1	Aminobiphenyl, 4-	92-67-1	3.3E+00	1.2E+01	6.4E+04	2.6E+00				
				8.0E-02	P						1.36E+09		1	0.1	Aminophenol, m-	591-27-5					6.3E+03	2.6E+04		5.1E+03
				4.0E-03	X						1.36E+09		1	0.1	Aminophenol, o-	95-55-6					3.1E+02	1.3E+03		2.5E+02
				2.0E-02	P						1.36E+09		1	0.1	Aminophenol, p-	123-30-8					1.6E+03	6.6E+03		1.3E+03
				2.5E-03	I						1.36E+09		1	0.1	Amitraz	33089-61-1					2.0E+02	8.2E+02		1.6E+02
						5.0E-01	I	V					1		Ammonia	7664-41-7								
				2.0E-03	X						1.36E+09		1	0.1	Ammonium Picrate	131-74-8					1.6E+02	6.6E+02		1.3E+02
				2.0E-01	I						1.36E+09		1		Ammonium Sulfamate	7773-06-0					1.6E+04			1.6E+04
						3.0E-03	X	V		1.37E+04	1.36E+09	2.62E+04	1		Amyl Alcohol, tert-	75-85-4							8.2E+01	8.2E+01
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I				1.36E+09		1	0.1	Aniline	62-53-3	1.2E+04	4.3E+04	2.4E+08	9.5E+03	5.5E+02	2.3E+03	1.4E+06	4.4E+02
4.0E-02	P			2.0E-03	X						1.36E+09		1	0.1	Anthraquinone, 9,10-	84-65-1	1.7E+03	6.2E+03		1.4E+03	1.6E+02	6.6E+02		1.3E+02
				4.0E-04	I	3.0E-04	A				1.36E+09		0.15		Antimony (metallic)	7440-36-0					3.1E+01		4.3E+05	3.1E+01
				5.0E-04	H						1.36E+09		0.15		Antimony Pentoxide	1314-60-9					3.9E+01			3.9E+01
				4.0E-04	H						1.36E+09		0.15		Antimony Tetroxide	1332-81-6					3.1E+01			3.1E+01
						2.0E-04	I				1.36E+09		0.15		Antimony Trioxide	1309-64-4							2.8E+05	2.8E+05
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C				1.36E+09		1	0.03	Arsenic, Inorganic	7440-38-2	7.7E+01	5.5E+02	8.9E+04	6.8E+01	3.9E+01	3.3E+02	2.1E+04	3.5E+01
				3.5E-06	C	5.0E-05	I				1.36E+09		1		Arsine	7784-42-1					2.7E-01		7.1E+04	2.7E-01
													1		Asbestos (units in fibers)	1332-21-4								
				3.6E-02	O						1.36E+09		1	0.1	Asulam	3337-71-1					2.8E+03	1.2E+04		2.3E+03
2.3E-01	C			3.0E-03	A						1.36E+09		1	0.1	Atrazine	1912-24-9	3.0E+02	1.1E+03		2.4E+02	2.3E+02	9.9E+02		1.9E+02
8.8E-01	C	2.5E-04	C								1.36E+09		1	0.1	Auramine	492-80-8	7.9E+01	2.8E+02	1.5E+06	6.2E+01				
				4.0E-04	I						1.36E+09		1	0.1	Avermectin B1	65195-55-3					3.1E+01	1.3E+02		2.5E+01
				3.0E-03	A	1.0E-02	A				1.36E+09		1	0.1	Azinphos-methyl	86-50-0					2.3E+02	9.9E+02	1.4E+07	1.9E+02
1.1E-01	I	3.1E-05	I					V			1.36E+09	5.23E+05	1		Azobenzene	103-33-3	6.3E+02		4.7E+03	5.6E+02				
				1.0E+00	P	7.0E-06	P				1.36E+09		1	0.1	Azodicarbonamide	123-77-3					7.8E+04	3.3E+05	9.9E+03	8.6E+03
				2.0E-01	I	5.0E-04	H				1.36E+09		0.07		Barium	7440-39-3					1.6E+04		7.1E+05	1.5E+04
				5.0E-03	O			V			1.36E+09	3.07E+05	1		Benfluralin	1861-40-1					3.9E+02			3.9E+02
				5.0E-02	I						1.36E+09		1	0.1	Benomyl	17804-35-2					3.9E+03	1.6E+04		3.2E+03
				2.0E-01	I						1.36E+09		1	0.1	Bensulfuron-methyl	83055-99-6					1.6E+04	6.6E+04		1.3E+04
				3.0E-02	I						1.36E+09		1	0.1	Bentazon	25057-89-0					2.3E+03	9.9E+03		1.9E+03
4.0E-03	P			1.0E-01	I			V		1.16E+03	1.36E+09	2.25E+04	1		Benzaldehyde	100-52-7	1.7E+04			1.7E+04	7.8E+03			7.8E+03
5.5E-02	I	7.8E-06	I	4.0E-03	I	3.0E-02	I	V		1.82E+03	1.36E+09	3.54E+03	1		Benzene	71-43-2	1.3E+03		1.3E+02	1.2E+02	3.1E+02		1.1E+02	8.2E+01
1.0E-01	X			3.0E-04	X						1.36E+09		1	0.1	Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1	7.0E+02	2.5E+03		5.4E+02	2.3E+01	9.9E+01		1.9E+01
				1.0E-03	P			V		1.26E+03	1.36E+09	1.94E+04	1		Benzenethiol	108-98-5					7.8E+01			7.8E+01
2.3E+02	I	6.7E-02	I	3.0E-03	I				M		1.36E+09		1	0.1	Benzidine	92-87-5	6.7E-02	2.6E-01	2.1E+03	5.3E-02	2.3E+02	9.9E+02		1.9E+02
				4.0E+00	I						1.36E+09		1	0.1	Benzoic Acid	65-85-0					3.1E+05	1.3E+06		2.5E+05
1.3E+01	I							V		3.24E+02	1.36E+09	6.76E+04	1		Benzotrichloride	98-07-7	5.3E+00			5.3E+00				
				1.0E-01	P						1.36E+09		1	0.1	Benzyl Alcohol	100-51-6					7.8E+03	3.3E+04		6.3E+03
1.7E-01	I	4.9E-05	C	2.0E-03	P	1.0E-03	P	V		1.46E+03	1.36E+09	2.55E+04	1		Benzyl Chloride	100-44-7	4.1E+02		1.5E+02	1.1E+02	1.6E+02		2.7E+01	2.3E+01
		2.4E-03	I	2.0E-03	I	2.0E-05	I				1.36E+09		0.007		Beryllium and compounds	7440-41-7			1.6E+05	1.6E+05	1.6E+02		2.8E+04	1.6E+02
				9.0E-03	P						1.36E+09		1	0.1	Bifenox	42576-02-3					7.0E+02	3.0E+03		5.7E+02
				1.5E-02	I						1.36E+09		1	0.1	Biphenthrin	82657-04-3					1.2E+03	4.9E+03		9.5E+02
8.0E-03	I			5.0E-01	I	4.0E-04	X	V			1.36E+09	1.14E+05	1		Biphenyl, 1,1'-	92-52-4	8.7E+03			8.7E+03	3.9E+04		4.8E+01	4.7E+01
				4.0E-02	I			V		1.02E+03	1.36E+09	3.50E+04	1		Bis(2-chloro-1-methylethyl) ether	108-60-1					3.1E+03			3.1E+03
				3.0E-03	P						1.36E+09		1	0.1	Bis(2-chloroethoxy)methane	111-91-1					2.3E+02	9.9E+02		1.9E+02
1.1E+00	I	3.3E-04	I					V		5.05E+03	1.36E+09	4.25E+04	1		Bis(2-chloroethyl)ether	111-44-4	6.3E+01		3.6E+01	2.3E+01				
2.2E+02	I	6.2E-02	I					V		4.22E+03	1.36E+09	1.88E+03	1		Bis(chloromethyl)ether	542-88-1	3.2E-01		8.5E-03	8.3E-03				
				5.0E-02	I						1.36E+09		1	0.1	Bisphenol A	80-05-7					3.9E+03	1.6E+04		3.2E+03
				2.0E-01	I	2.0E-02	H				1.36E+09		1		Boron And Borates Only	7440-42-8					1.6E+04		2.8E+07	1.6E+04
				2.0E+00	P	2.0E-02	P	V			1.36E+09		1		Boron Trichloride	10294-34-5					1.6E+05		2.8E+07	1.6E+05
				4.0E-02	C	1.3E-02	C	V			1.36E+09		1		Boron Trifluoride	7637-07-2					3.1E+03		1.8E+07	3.1E+03
7.0E-01	I			4.0E-03	I						1.36E+09		1		Bromate	15541-45-4	9.9E+01			9.9E+01	3.1E+02			3.1E+02
				1.0E-04	X	6.0E-05	X	V		2.38E+03	1.36E+09	5.92E+03	1		Bromo-2-chloroethane, 1-	107-04-0					7.8E+00		3.7E-01	3.5E-01
				3.0E-04	X			V		8.96E+02	1.36E+09	1.12E+04	1		Bromo-3-fluorobenzene, 1-	1073-06-9					2.3E+01			2.3E+01

RMLs TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
				3.0E-04	X			V		3.23E+02	1.36E+09	1.14E+04	1		Bromo-4-fluorobenzene, 1-	460-00-4					2.3E+01			2.3E+01
											1.36E+09		1	0.1	Bromoacetic acid	79-08-3								
				8.0E-03	I	6.0E-02	I	V		6.79E+02	1.36E+09	8.37E+03	1		Bromobenzene	108-86-1					6.3E+02		5.2E+02	2.9E+02
						4.0E-02	X	V		4.04E+03	1.36E+09	3.58E+03	1		Bromochloromethane	74-97-5							1.5E+02	1.5E+02
6.2E-02	I	3.7E-05	C	8.0E-03	P			V		9.32E+02	1.36E+09	3.97E+03	1		Bromodichloromethane	75-27-4	1.1E+03		3.0E+01	2.9E+01	6.3E+02			6.3E+02
7.9E-03	I	1.1E-06	I	2.0E-02	I			V		9.15E+02	1.36E+09	9.70E+03	1		Bromoform	75-25-2	8.8E+03		2.5E+03	1.9E+03	1.6E+03			1.6E+03
				1.4E-03	I	5.0E-03	I	V		3.59E+03	1.36E+09	1.40E+03	1		Bromomethane	74-83-9					1.1E+02		7.3E+00	6.8E+00
				5.0E-03	H			V			1.36E+09	1.24E+05	1		Bromophos	2104-96-3					3.9E+02			3.9E+02
						1.0E-01	A	V		9.66E+02	1.36E+09	2.14E+03	1		Bromopropane, 1-	106-94-5							2.2E+02	2.2E+02
1.0E-01	O			1.5E-02	O						1.36E+09		1	0.1	Bromoxynil	1689-84-5	6.7E+02	2.4E+03		5.3E+02	1.2E+03	4.9E+03		9.5E+02
1.0E-01	O			1.5E-02	O			V			1.36E+09	4.74E+05	1		Bromoxynil Octanoate	1689-99-2	6.7E+02			6.7E+02	1.2E+03			1.2E+03
6.0E-01	C	3.0E-05	I			2.0E-03	I	V		6.67E+02	1.36E+09	8.66E+02	1		Butadiene, 1,3-	106-99-0	1.2E+02		8.1E+00	7.6E+00			1.8E+00	1.8E+00
				1.0E-01	I			V		7.64E+03	1.36E+09	3.00E+04	1		Butanol, N-	71-36-3					7.8E+03			7.8E+03
5.0E-04	I			4.0E-01	I	5.0E+00	I	V			1.36E+09	2.87E+04	1		Butyl Alcohol, t-	75-65-0	1.4E+05			1.4E+05	3.1E+04		1.5E+05	2.6E+04
				2.0E+00	P	3.0E+01	P	V		2.13E+04	1.36E+09	2.92E+04	1		Butyl alcohol, sec-	78-92-2					1.6E+05		9.1E+05	1.3E+05
				5.0E-02	I			V			1.36E+09	8.63E+04	1		Butylate	2008-41-5					3.9E+03			3.9E+03
2.0E-04	C	5.7E-08	C								1.36E+09		1	0.1	Butylated hydroxyanisole	25013-16-5	3.5E+05	1.2E+06	6.7E+09	2.7E+05				
3.6E-03	P			3.0E-01	P						1.36E+09		1	0.1	Butylated hydroxytoluene	128-37-0	1.9E+04	6.9E+04		1.5E+04	2.3E+04	9.9E+04		1.9E+04
				5.0E-02	P			V		1.08E+02	1.36E+09	8.14E+03	1		Butylbenzene, n-	104-51-8					3.9E+03			3.9E+03
				1.0E-01	X			V		1.45E+02	1.36E+09	7.35E+03	1		Butylbenzene, sec-	135-98-8					7.8E+03			7.8E+03
				1.0E-01	X			V		1.83E+02	1.36E+09	7.36E+03	1		Butylbenzene, tert-	98-06-6					7.8E+03			7.8E+03
				2.0E-02	A						1.36E+09		1	0.1	Cacodylic Acid	75-60-5					1.6E+03	6.6E+03		1.3E+03
		1.8E-03	I	1.0E-04	A	1.0E-05	A				1.36E+09		0.025	0.001	Cadmium (Diet)	7440-43-9			2.1E+05	2.1E+05	7.8E+00	8.2E+01	1.4E+04	7.1E+00
		1.8E-03	I	1.0E-04	A	1.0E-05	A						0.05	0.001	Cadmium (Water)	7440-43-9								
				5.0E-01	I	2.2E-03	C				1.36E+09		1	0.1	Caprolactam	105-60-2					3.9E+04	1.6E+05	3.1E+06	3.1E+04
1.5E-01	C	4.3E-05	C	2.0E-03	I						1.36E+09		1	0.1	Captafol	2425-06-1	4.6E+02	1.6E+03	8.9E+06	3.6E+02	1.6E+02	6.6E+02		1.3E+02
2.3E-03	C	6.6E-07	C	1.3E-01	I						1.36E+09		1	0.1	Captan	133-06-2	3.0E+04	1.1E+05	5.8E+08	2.4E+04	1.0E+04	4.3E+04		8.2E+03
				1.0E-01	I						1.36E+09		1	0.1	Carbaryl	63-25-2					7.8E+03	3.3E+04		6.3E+03
				5.0E-03	I						1.36E+09		1	0.1	Carbofuran	1563-66-2					3.9E+02	1.6E+03		3.2E+02
				1.0E-01	I	7.0E-01	I	V		7.38E+02	1.36E+09	1.17E+03	1		Carbon Disulfide	75-15-0					7.8E+03		8.5E+02	7.7E+02
7.0E-02	I	6.0E-06	I	4.0E-03	I	1.0E-01	I	V		4.58E+02	1.36E+09	1.49E+03	1		Carbon Tetrachloride	56-23-5	9.9E+02		7.0E+01	6.5E+01	3.1E+02		1.6E+02	1.0E+02
						1.0E-01	P	V		5.89E+03	1.36E+09	6.46E+02	1		Carbonyl Sulfide	463-58-1							6.7E+01	6.7E+01
				1.0E-02	I						1.36E+09		1	0.1	Carbosulfan	55285-14-8					7.8E+02	3.3E+03		6.3E+02
				1.0E-01	I						1.36E+09		1	0.1	Carboxin	5234-68-4					7.8E+03	3.3E+04		6.3E+03
						9.0E-04	I				1.36E+09		1		Ceric oxide	1306-38-3							1.3E+06	1.3E+06
				1.0E-01	I			V			1.36E+09	1.45E+05	1		Chloral Hydrate	302-17-0					7.8E+03			7.8E+03
				1.5E-02	I						1.36E+09		1	0.1	Chloramben	133-90-4					1.2E+03	4.9E+03		9.5E+02
											1.36E+09		1	0.1	Chloramines, Organic	E701235								
4.0E-01	H										1.36E+09		1	0.1	Chloranil	118-75-2	1.7E+02	6.1E+02		1.3E+02				
				5.0E-04	G			V			1.36E+09	1.49E+06	1	0.04	Chlordane (alpha)	5103-71-9					3.9E+01	4.1E+02		3.6E+01
				5.0E-04	G			V			1.36E+09	1.49E+06	1	0.04	Chlordane (gamma)	5103-74-2					3.9E+01	4.1E+02		3.6E+01
3.5E-01	I	1.0E-04	I	5.0E-04	I	7.0E-04	I	V			1.36E+09	1.53E+06	1	0.04	Chlordane (technical mixture)	12789-03-6	2.0E+02	1.8E+03	4.3E+03	1.7E+02	3.9E+01	4.1E+02	1.1E+03	3.5E+01
1.0E+01	I	4.6E-03	C	3.0E-04	I						1.36E+09		1	0.1	Chlordecone (Kepone)	143-50-0	7.0E+00	2.5E+01	8.3E+04	5.4E+00	2.3E+01	9.9E+01		1.9E+01
				7.0E-04	A						1.36E+09		1	0.1	Chlorfenvinphos	470-90-6					5.5E+01	2.3E+02		4.4E+01
				9.0E-02	O						1.36E+09		1	0.1	Chlorimuron, Ethyl-	90982-32-4					7.0E+03	3.0E+04		5.7E+03
				1.0E-01	I	1.5E-04	A	V		2.78E+03	1.36E+09	1.22E+03	1		Chlorine	7782-50-5					7.8E+03		1.8E-01	1.8E-01
				3.0E-02	I	2.0E-04	I	V			1.36E+09		1		Chlorine Dioxide	10049-04-4					2.3E+03		2.8E+05	2.3E+03
				3.0E-02	I						1.36E+09		1		Chlorite (Sodium Salt)	7758-19-2					2.3E+03			2.3E+03
						5.0E+01	I	V		1.15E+03	1.36E+09	1.03E+03	1		Chloro-1,1-difluoroethane, 1-	75-68-3							5.4E+04	5.4E+04
		3.0E-04	I	2.0E-02	H	2.0E-02	I	V		7.86E+02	1.36E+09	1.08E+03	1		Chloro-1,3-butadiene, 2-	126-99-8			1.0E+00	1.0E+00	1.6E+03		2.2E+01	2.2E+01
4.6E-01	H										1.36E+09		1	0.1	Chloro-2-methylaniline HCl, 4-	3165-93-3	1.5E+02	5.4E+02		1.2E+02				
1.0E-01	P	7.7E-05	C	3.0E-03	X						1.36E+09		1	0.1	Chloro-2-methylaniline, 4-	95-69-2	7.0E+02	2.5E+03	5.0E+06	5.4E+02	2.3E+02	9.9E+02		1.9E+02
2.7E-01	X							V		1.18E+04	1.36E+09	1.62E+04	1		Chloroacetaldehyde, 2-	107-20-0	2.6E+02			2.6E+02				
											1.36E+09		1	0.1	Chloroacetic Acid	79-11-8								
						3.0E-05	I				1.36E+09		1	0.1	Chloroacetophenone, 2-	532-27-4							4.3E+04	4.3E+04
2.0E-01	P			5.0E-04	P						1.36E+09		1	0.1	Chloroaniline, p-	106-47-8	3.5E+02	1.2E+03		2.7E+02	3.9E+01	1.6E+02		3.2E+01
				2.0E-02	I	5.0E-02	P	V		7.61E+02	1.36E+09	6.45E+03	1		Chlorobenzene	108-90-7					1.6E+03		3.4E+02	2.8E+02
				1.0E-01	X						1.36E+09		1	0.1	Chlorobenzene sulfonic acid, p-	98-66-8					7.8E+03	3.3E+04		6.3E+03
1.1E-01	C	3.1E-05	C	2.0E-02	I						1.36E+09		1	0.1	Chlorobenzilate	510-15-6	6.3E+02	2.2E+03	1.2E+07	4.9E+02	1.6E+03	6.6E+03		1.3E+03
				3.0E-02	X						1.36E+09		1	0.1	Chlorobenzoic Acid, p-	74-11-3					2.3E+03	9.9E+03		1.9E+03
		8.6E-06	C	3.0E-03	P	3.0E-01	P	V		2.90E+02	1.36E+09	6.76E+03	1		Chlorobenzotrifluoride, 4-	98-56-6			2.2E+02	2.2E+02	2.3E+02		2.1E+03	2.1E+02
				4.0E-02	P			V		7.28E+02	1.36E+09	1.76E+03	1		Chlorobutane, 1-	109-69-3					3.1E+03			3.1E+03
						5.0E+01	I	V		1.68E+03	1.36E+09	9.38E+02	1		Chlorodifluoromethane	75-45-6							4.9E+04	4.9E+04
				2.0E-02	P			V		1.11E+05	1.36E+09	7.81E+04	1		Chloroethanol, 2-	107-07-3					1.6E+03			1.6E+03
3.1E-02	C	2.3E-05	I	1.0E-02	I	9.8E-02	A	V		2.54E+03	1.36E+09	2.63E+03	1		Chloroform	67-66-3	2.2E+03		3.2E+01	3.2E+01	7.8E+02		2.7E+02	2.0E+02
						9.0E-02	I	V		1.32E+03	1.36E+09	1.18E+03	1		Chloromethane	74-87-3							1.1E+02	1.1E+02
2.4E+00	C	6.9E-04	C					V		9.32E+03	1.36E+09	5.33E+03	1		Chloromethyl Methyl Ether	107-30-2	2.9E+01		2.2E+00	2.0E+00				
3.0E-01	P			3.0E-03	P	1.0E-05	X				1.36E+09		1	0.1	Chloronitrobenzene, o-	88-73-3	2.3E+02	8.2E+02		1.8E+02	2.3E+02	9.9E+02	1.4E+04	1.9E+02
6.0E-02	P			7.0E-04	P	2.0E-03	P				1.36E+09		1	0.1	Chloronitrobenzene, p-	100-00-5	1.2E+03	4.1E+03		9.0E+02	5.5E+01	2.3E+02	2.8E+06	4.4E+01
				5.0E-03	I			V		2.74E+04	1.36E+09	1.39E+05	1		Chlorophenol, 2-	95-57-8					3.9E+02			3.9E+02
						4.0E-04	C	V		6.17E+02	1.36E+09	4.68E+03	1		Chloropicrin	76-06-2							2.0E+00	2.0E+00
1.7E-02	C			1.5E-02	I						1.36E+09		1	0.1	Chlorothalonil	1897-45-6	4.1E+03	1.5E+04		3.2E+03	1.2E+03	4.9E+03		9.5E+02
				2.0E-02	I			V		9.07E+02	1.36E+09	8.12E+03	1		Chlorotoluene, o-	95-49-8					1.6E+03			1.6E+03
				2.0E-02	X			V		2.53E+02	1.36E+09	7.29E+03	1		Chlorotoluene, p-	106-43-4					1.6E+03			1.6E+03
2.4E+02	C	6.9E-02	C								1.36E+09		1	0.1	Chlorozotocin	54749-90-5	2.9E-01	1.0E+00	5.5E+03	2.3E-01				
				5.0E-03	O						1.36E+09		1	0.1	Chlorpropham	101-21-3					3.9E+02	1.6E+03		3.2E+02
				1.0E-03	A						1.36E+09		1	0.1	Chlorpyrifos	2921-88-2					7.8E+01	3.3E+02		6.3E+01

RMLs TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
				1.0E-02	H						1.36E+09		1	0.1	Chlorpyrifos Methyl	5598-13-0					7.8E+02	3.3E+03		6.3E+02
				5.0E-02	O						1.36E+09		1	0.1	Chlorsulfuron	64902-72-3					3.9E+03	1.6E+04		3.2E+03
				1.0E-02	I						1.36E+09		1	0.1	Chlorthal-dimethyl	1861-32-1					7.8E+02	3.3E+03		6.3E+02
				8.0E-04	H						1.36E+09		1	0.1	Chlorthiophos	60238-56-4					6.3E+01	2.6E+02		5.1E+01
				1.5E+00	I						1.36E+09		0.013		Chromium(III), Insoluble Salts	16065-83-1					1.2E+05			1.2E+05
5.0E-01	C	8.4E-02	G	3.0E-03	I	1.0E-04	I		M		1.36E+09		0.025		Chromium(VI)	18540-29-9	3.1E+01		1.6E+03	3.0E+01	2.3E+02		1.4E+05	2.3E+02
											1.36E+09		0.013		Chromium, Total	7440-47-3								
				1.3E-02	I						1.36E+09		1	0.1	Clofentezine	74115-24-5					1.0E+03	4.3E+03		8.2E+02
		9.0E-03	P	3.0E-04	P	6.0E-06	P				1.36E+09		1		Cobalt	7440-48-4			4.2E+04	4.2E+04	2.3E+01		8.5E+03	2.3E+01
		6.2E-04	I					V	M				1		Coke Oven Emissions	E649830								
				4.0E-02	H						1.36E+09		1		Copper	7440-50-8					3.1E+03			3.1E+03
				5.0E-02	I	6.0E-01	C				1.36E+09		1	0.1	Cresol, m-	108-39-4					3.9E+03	1.6E+04	8.5E+08	3.2E+03
				5.0E-02	I	6.0E-01	C				1.36E+09		1	0.1	Cresol, o-	95-48-7					3.9E+03	1.6E+04	8.5E+08	3.2E+03
				2.0E-02	P	6.0E-01	C				1.36E+09		1	0.1	Cresol, p-	106-44-5					1.6E+03	6.6E+03	8.5E+08	1.3E+03
				1.0E-01	A						1.36E+09		1	0.1	Cresol, p-chloro-m-	59-50-7					7.8E+03	3.3E+04		6.3E+03
				1.0E-01	A	6.0E-01	C				1.36E+09		1	0.1	Cresols	1319-77-3					7.8E+03	3.3E+04	8.5E+08	6.3E+03
1.9E+00	H			1.0E-03	P			V		1.66E+04	1.36E+09	1.89E+04	1		Crotonaldehyde, trans-	123-73-9	3.7E+01			3.7E+01	7.8E+01			7.8E+01
				1.0E-01	I	4.0E-01	I	V		2.68E+02	1.36E+09	6.21E+03	1		Cumene	98-82-8					7.8E+03		2.6E+03	1.9E+03
2.2E-01	C	6.3E-05	C								1.36E+09		1	0.1	Cupferron	135-20-6	3.2E+02	1.1E+03	6.1E+06	2.5E+02				
8.4E-01	H			2.0E-03	H						1.36E+09		1	0.1	Cyanazine	21725-46-2	8.3E+01	2.9E+02		6.5E+01	1.6E+02	6.6E+02		1.3E+02
															Cyanides									
				1.0E-03	I						1.36E+09		1		~Calcium Cyanide	592-01-8					7.8E+01			7.8E+01
				5.0E-03	I						1.36E+09		1		~Copper Cyanide	544-92-3					3.9E+02			3.9E+02
				6.0E-04	I	8.0E-04	G	V		9.54E+05	1.36E+09	5.33E+04	1		~Cyanide (CN-)	57-12-5					4.7E+01		4.4E+01	2.3E+01
				1.0E-03	I			V			1.36E+09		1		~Cyanogen	460-19-5					7.8E+01			7.8E+01
				9.0E-02	I			V			1.36E+09		1		~Cyanogen Bromide	506-68-3					7.0E+03			7.0E+03
				5.0E-02	I			V			1.36E+09		1		~Cyanogen Chloride	506-77-4					3.9E+03			3.9E+03
				6.0E-04	I	8.0E-04	I	V		1.00E+07	1.36E+09	5.22E+04	1		~Hydrogen Cyanide	74-90-8					4.7E+01		4.4E+01	2.3E+01
				2.0E-03	I						1.36E+09		1		~Potassium Cyanide	151-50-8					1.6E+02			1.6E+02
				5.0E-03	I						1.36E+09		0.04		~Potassium Silver Cyanide	506-61-6					3.9E+02			3.9E+02
				1.0E-01	I						1.36E+09		0.04		~Silver Cyanide	506-64-9					7.8E+03			7.8E+03
				1.0E-03	I						1.36E+09		1		~Sodium Cyanide	143-33-9					7.8E+01			7.8E+01
				2.0E-04	P						1.36E+09		1		~Thiocyanates	E1790665					1.6E+01			1.6E+01
				2.0E-04	X			V			1.36E+09		1		~Thiocyanic Acid	463-56-9					1.6E+01			1.6E+01
				5.0E-02	I						1.36E+09		1		~Zinc Cyanide	557-21-1					3.9E+03			3.9E+03
						6.0E+00	I	V		1.17E+02	1.36E+09	1.04E+03	1		Cyclohexane	110-82-7							6.5E+03	6.5E+03
2.0E-02	X			2.0E-02	X						1.36E+09		1	0.1	Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	3.5E+03	1.2E+04		2.7E+03	1.6E+03	6.6E+03		1.3E+03
				5.0E+00	I	7.0E-01	P	V		5.11E+03	1.36E+09	4.17E+04	1		Cyclohexanone	108-94-1					3.9E+05		3.0E+04	2.8E+04
				5.0E-03	P	1.0E+00	X	V		2.83E+02	1.36E+09	1.46E+03	1		Cyclohexene	110-83-8					3.9E+02		1.5E+03	3.1E+02
				2.0E-01	I			V		2.93E+05	1.36E+09	7.46E+04	1		Cyclohexylamine	108-91-8					1.6E+04			1.6E+04
				2.5E-02	I						1.36E+09		1	0.1	Cyfluthrin	68359-37-5					2.0E+03	8.2E+03		1.6E+03
				7.2E-02	O						1.36E+09		1	0.1	Cypermethrin	52315-07-8					5.6E+03	2.4E+04		4.5E+03
				5.0E-01	O						1.36E+09		1	0.1	Cyromazine	66215-27-8					3.9E+04	1.6E+05		3.2E+04
2.4E-01	I	6.9E-05	C	3.0E-05	X						1.36E+09		1	0.1	DDD, p,p`- (DDD)	72-54-8	2.9E+02	1.0E+03	5.5E+06	2.3E+02	2.3E+00	9.9E+00		1.9E+00
3.4E-01	I	9.7E-05	C	3.0E-04	X			V			1.36E+09	2.10E+06	1		DDE, p,p'-	72-55-9	2.0E+02		6.1E+03	2.0E+02	2.3E+01			2.3E+01
3.4E-01	I	9.7E-05	I	5.0E-04	I						1.36E+09		1	0.03	DDT	50-29-3	2.0E+02	2.4E+03	3.9E+06	1.9E+02	3.9E+01	5.5E+02		3.7E+01
				3.0E-02	I						1.36E+09		1	0.1	Dalapon	75-99-0					2.3E+03	9.9E+03		1.9E+03
1.8E-02	C	5.1E-06	C	1.5E-01	I						1.36E+09		1	0.1	Daminozide	1596-84-5	3.9E+03	1.4E+04	7.5E+07	3.0E+03	1.2E+04	4.9E+04		9.5E+03
7.0E-04	I			7.0E-03	I						1.36E+09		1	0.1	Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5	9.9E+04	3.5E+05		7.8E+04	5.5E+02	2.3E+03		4.4E+02
				4.0E-05	I						1.36E+09		1	0.1	Demeton	8065-48-3					3.1E+00	1.3E+01		2.5E+00
1.2E-03	I			6.0E-01	I						1.36E+09		1	0.1	Di(2-ethylhexyl)adipate	103-23-1	5.8E+04	2.1E+05		4.5E+04	4.7E+04	2.0E+05		3.8E+04
6.1E-02	H										1.36E+09		1	0.1	Diallate	2303-16-4	1.1E+03	4.1E+03		8.9E+02				
				7.0E-04	A						1.36E+09		1	0.1	Diazinon	333-41-5					5.5E+01	2.3E+02		4.4E+01
				1.0E-02	X			V			1.36E+09	5.24E+05	1		Dibenzothiophene	132-65-0					7.8E+02			7.8E+02
8.0E-01	P	6.0E-03	P	2.0E-04	P	2.0E-04	I	V	M	9.79E+02	1.36E+09	3.20E+04	1		Dibromo-3-chloropropane, 1,2-	96-12-8	1.9E+01		5.4E-01	5.3E-01	1.6E+01		6.7E+00	4.7E+00
											1.36E+09		1	0.1	Dibromoacetic acid	631-64-1								
				4.0E-04	X			V		1.59E+02	1.36E+09	1.93E+04	1		Dibromobenzene, 1,3-	108-36-1					3.1E+01			3.1E+01
				1.0E-02	I			V			1.36E+09	2.20E+04	1		Dibromobenzene, 1,4-	106-37-6					7.8E+02			7.8E+02
8.4E-02	I			2.0E-02	I			V		8.02E+02	1.36E+09	7.95E+03	1		Dibromochloromethane	124-48-1	8.3E+02			8.3E+02	1.6E+03			1.6E+03
2.0E+00	I	6.0E-04	I	9.0E-03	I	9.0E-03	I	V		1.34E+03	1.36E+09	8.64E+03	1		Dibromoethane, 1,2-	106-93-4	3.5E+01		4.0E+00	3.6E+00	7.0E+02		8.1E+01	7.3E+01
						4.0E-03	X	V		2.82E+03	1.36E+09	5.64E+03	1		Dibromomethane (Methylene Bromide)	74-95-3							2.4E+01	2.4E+01
				3.0E-04	P						1.36E+09		1	0.1	Dibutyltin Compounds	E1790661					2.3E+01	9.9E+01		1.9E+01
				3.0E-02	I						1.36E+09		1	0.1	Dicamba	1918-00-9					2.3E+03	9.9E+03		1.9E+03
											1.36E+09		1	0.1	Dichloramine	3400-09-7								
		4.2E-03	P					V		5.54E+02	1.36E+09	3.21E+03	1		Dichloro-2-butene, 1,4-	764-41-0			2.1E-01	2.1E-01				
		4.2E-03	P					V		5.19E+02	1.36E+09	1.11E+04	1		Dichloro-2-butene, cis-1,4-	1476-11-5			7.4E-01	7.4E-01				
		4.2E-03	P					V		7.60E+02	1.36E+09	1.11E+04	1		Dichloro-2-butene, trans-1,4-	110-57-6			7.4E-01	7.4E-01				
5.0E-02	I			4.0E-03	I						1.36E+09		1	0.1	Dichloroacetic Acid	79-43-6	1.4E+03	4.9E+03		1.1E+03	3.1E+02	1.3E+03		2.5E+02
				9.0E-02	I	2.0E-01	H	V		3.76E+02	1.36E+09	1.17E+04	1		Dichlorobenzene, 1,2-	95-50-1					7.0E+03		2.4E+03	1.8E+03
5.4E-03	C	1.1E-05	C	7.0E-02	A	8.0E-01	I	V			1.36E+09	1.04E+04	1		Dichlorobenzene, 1,4-	106-46-7	1.3E+04		2.7E+02	2.6E+02	5.5E+03		8.7E+03	3.4E+03
4.5E-01	I	3.4E-04	C								1.36E+09		1	0.1	Dichlorobenzidine, 3,3'-	91-94-1	1.5E+02	5.5E+02	1.1E+06	1.2E+02				
				9.0E-03	X						1.36E+09		1	0.1	Dichlorobenzophenone, 4,4'-	90-98-2					7.0E+02	3.0E+03		5.7E+02
				2.0E-01	I	1.0E-01	X	V		8.45E+02	1.36E+09	8.41E+02	1		Dichlorodifluoromethane	75-71-8					1.6E+04		8.8E+01	8.7E+01
5.7E-03	C	1.6E-06	C	2.0E-01	P			V		1.69E+03	1.36E+09	2.08E+03	1		Dichloroethane, 1,1-	75-34-3	1.2E+04		3.7E+02	3.6E+02	1.6E+04			1.6E+04
9.1E-02	I	2.6E-05	I	6.0E-03	X	7.0E-03	P	V		2.98E+03	1.36E+09	4.57E+03	1		Dichloroethane, 1,2-	107-06-2	7.6E+02		4.9E+01	4.6E+01	4.7E+02		3.3E+01	3.1E+01
				5.0E-02	I	2.0E-01	I	V		1.19E+03	1.36E+09	1.16E+03	1		Dichloroethylene, 1,1-	75-35-4					3.9E+03		2.4E+02	2.3E+02
				2.0E-03	I			V		2.37E+03	1.36E+09	2.50E+03	1		Dichloroethylene, cis-1,2-	156-59-2					1.6E+02			1.6E+02

RMLS TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information															Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
				2.0E-02	I	4.0E-02	X	V		1.85E+03	1.36E+09	1.75E+03	1		Dichloroethylene, trans-1,2-	156-60-5					1.6E+03		7.3E+01	7.0E+01
				3.0E-03	I						1.36E+09		1	0.1	Dichlorophenol, 2,4-	120-83-2					2.3E+02	9.9E+02		1.9E+02
				1.0E-02	I						1.36E+09		1	0.05	Dichlorophenoxy Acetic Acid, 2,4-	94-75-7					7.8E+02	6.6E+03		7.0E+02
3.7E-02	P	3.7E-06	P	4.0E-02	P	4.0E-03	I	V		1.36E+03	1.36E+09	3.79E+03	1		Dichloropropane, 1,2-	78-87-5	1.9E+03		2.9E+02	2.5E+02	3.1E+03		1.6E+01	1.6E+01
				2.0E-02	P			V		1.49E+03	1.36E+09	6.76E+03	1		Dichloropropane, 1,3-	142-28-9					1.6E+03			1.6E+03
				3.0E-03	I						1.36E+09		1	0.1	Dichloropropanol, 2,3-	616-23-9					2.3E+02	9.9E+02		1.9E+02
1.0E-01	I	4.0E-06	I	3.0E-02	I	2.0E-02	I	V		1.57E+03	1.36E+09	3.55E+03	1		Dichloropropene, 1,3-	542-75-6	7.0E+02		2.5E+02	1.8E+02	2.3E+03		7.4E+01	7.2E+01
2.9E-01	I	8.3E-05	C	5.0E-04	I	5.0E-04	I				1.36E+09		1	0.1	Dichlorvos	62-73-7	2.4E+02	8.5E+02	4.6E+06	1.9E+02	3.9E+01	1.6E+02	7.1E+05	3.2E+01
				3.0E-05	O						1.36E+09		1	0.1	Dicrotophos	141-66-2					2.3E+00	9.9E+00		1.9E+00
				8.0E-02	P	3.0E-04	X	V		2.56E+02	1.36E+09	4.11E+03	1		Dicyclopentadiene	77-73-6					6.3E+03		1.3E+00	1.3E+00
1.6E+01	I	4.6E-03	I	5.0E-05	I						1.36E+09		1	0.1	Dieldrin	60-57-1	4.3E+00	1.5E+01	8.3E+04	3.4E+00	3.9E+00	1.6E+01		3.2E+00
		3.0E-04	C			5.0E-03	I						1	0.1	Diesel Engine Exhaust	E17136615								
				2.0E-03	P	2.0E-04	P				1.36E+09		1	0.1	Diethanolamine	111-42-2					1.6E+02	6.6E+02	2.8E+05	1.3E+02
				3.0E-02	P	1.0E-04	P				1.36E+09		1	0.1	Diethylene Glycol Monobutyl Ether	112-34-5					2.3E+03	9.9E+03	1.4E+05	1.9E+03
				6.0E-02	P	3.0E-04	P				1.36E+09		1	0.1	Diethylene Glycol Monoethyl Ether	111-90-0					4.7E+03	2.0E+04	4.3E+05	3.8E+03
				1.0E-03	P			V		1.12E+05	1.36E+09	1.39E+05	1		Diethylformamide	617-84-5					7.8E+01			7.8E+01
3.5E+02	C	1.0E-01	C								1.36E+09		1	0.1	Diethylstilbestrol	56-53-1	2.0E-01	7.1E-01	3.8E+03	1.6E-01				
				8.3E-02	O						1.36E+09		1	0.1	Difenzoquat	43222-48-6					6.5E+03	2.7E+04		5.2E+03
				2.0E-02	I						1.36E+09		1	0.1	Diflubenzuron	35367-38-5					1.6E+03	6.6E+03		1.3E+03
						4.0E+01	I	V		1.43E+03	1.36E+09	1.15E+03	1		Difluoroethane, 1,1-	75-37-6							4.8E+04	4.8E+04
						3.0E+01	X	V		6.91E+02	1.36E+09	7.58E+02	1		Difluoropropane, 2,2-	420-45-1							2.4E+04	2.4E+04
4.4E-02	C	1.3E-05	C					V			1.36E+09	1.23E+05	1		Dihydrosafrole	94-58-6	1.6E+03		2.7E+03	9.9E+02				
						7.0E-01	P	V		2.26E+03	1.36E+09	3.06E+03	1		Diisopropyl Ether	108-20-3							2.2E+03	2.2E+03
				8.0E-02	I			V		5.30E+02	1.36E+09	3.81E+04	1		Diisopropyl Methylphosphonate	1445-75-6					6.3E+03			6.3E+03
				2.2E-02	O						1.36E+09		1	0.1	Dimethipin	55290-64-7					1.7E+03	7.2E+03		1.4E+03
				2.2E-03	O						1.36E+09		1	0.1	Dimethoate	60-51-5					1.7E+02	7.3E+02		1.4E+02
1.6E+00	P										1.36E+09		1	0.1	Dimethoxybenzidine, 3,3'-	119-90-4	4.3E+01	1.5E+02		3.4E+01				
1.7E-03	P			6.0E-02	P						1.36E+09		1	0.1	Dimethyl methylphosphonate	756-79-6	4.1E+04	1.5E+05		3.2E+04	4.7E+03	2.0E+04		3.8E+03
4.6E+00	C	1.3E-03	C								1.36E+09		1	0.1	Dimethylamino azobenzene [p-]	60-11-7	1.5E+01	5.4E+01	2.9E+05	1.2E+01				
5.8E-01	H										1.36E+09		1	0.1	Dimethylaniline HCl, 2,4-	21436-96-4	1.2E+02	4.3E+02		9.4E+01				
2.0E-01	P			2.0E-03	X						1.36E+09		1	0.1	Dimethylaniline, 2,4-	95-68-1	3.5E+02	1.2E+03		2.7E+02	1.6E+02	6.6E+02		1.3E+02
2.7E-02	P			2.0E-03	I			V		8.30E+02	1.36E+09	3.13E+04	1		Dimethylaniline, N,N-	121-69-7	2.6E+03			2.6E+03	1.6E+02			1.6E+02
1.1E+01	P										1.36E+09		1	0.1	Dimethylbenzidine, 3,3'-	119-93-7	6.3E+00	2.2E+01		4.9E+00				
				1.0E-01	P	3.0E-02	I	V		1.06E+05	1.36E+09	1.28E+05	1		Dimethylformamide	68-12-2					7.8E+03		4.0E+03	2.6E+03
				1.0E-04	X	2.0E-06	X	V		1.72E+05	1.36E+09	2.77E+04	1		Dimethylhydrazine, 1,1-	57-14-7					7.8E+00		5.8E-02	5.7E-02
5.5E+02	C	1.6E-01	C					V		1.89E+05	1.36E+09	1.68E+05	1		Dimethylhydrazine, 1,2-	540-73-8	1.3E-01		2.9E-01	8.8E-02				
				2.0E-02	I						1.36E+09		1	0.1	Dimethylphenol, 2,4-	105-67-9					1.6E+03	6.6E+03		1.3E+03
				6.0E-04	I						1.36E+09		1	0.1	Dimethylphenol, 2,6-	576-26-1					4.7E+01	2.0E+02		3.8E+01
				1.0E-03	I						1.36E+09		1	0.1	Dimethylphenol, 3,4-	95-65-8					7.8E+01	3.3E+02		6.3E+01
4.5E-02	C	1.3E-05	C					V		4.73E+02	1.36E+09	5.48E+03	1		Dimethylvinylchloride	513-37-1	1.5E+03		1.2E+02	1.1E+02				
				8.0E-05	X						1.36E+09		1	0.1	Dinitro-o-cresol, 4,6-	534-52-1					6.3E+00	2.6E+01		5.1E+00
				2.0E-03	I						1.36E+09		1	0.1	Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5					1.6E+02	6.6E+02		1.3E+02
				4.0E-04	X	2.0E-03	X				1.36E+09		1	0.1	Dinitroaniline, 3,5-	618-87-1					3.1E+01	1.3E+02	2.8E+06	2.5E+01
				1.0E-04	P						1.36E+09		1	0.1	Dinitrobenzene, 1,2-	528-29-0					7.8E+00	3.3E+01		6.3E+00
				1.0E-04	I						1.36E+09		1	0.1	Dinitrobenzene, 1,3-	99-65-0					7.8E+00	3.3E+01		6.3E+00
				1.0E-04	P						1.36E+09		1	0.1	Dinitrobenzene, 1,4-	100-25-4					7.8E+00	3.3E+01		6.3E+00
				2.0E-03	I						1.36E+09		1	0.1	Dinitrophenol, 2,4-	51-28-5					1.6E+02	6.6E+02		1.3E+02
6.8E-01	I										1.36E+09		1	0.1	Dinitrotoluene Mixture, 2,4/2,6-	E1615210	1.0E+02	3.6E+02		8.0E+01				
3.1E-01	C	8.9E-05	C	2.0E-03	I						1.36E+09		1	0.102	Dinitrotoluene, 2,4-	121-14-2	2.2E+02	7.8E+02	4.3E+06	1.7E+02	1.6E+02	6.5E+02		1.3E+02
1.5E+00	P			3.0E-04	X						1.36E+09		1	0.099	Dinitrotoluene, 2,6-	606-20-2	4.6E+01	1.7E+02		3.6E+01	2.3E+01	1.0E+02		1.9E+01
				1.0E-04	X						1.36E+09		1	0.006	Dinitrotoluene, 2-Amino-4,6-	35572-78-2					7.8E+00	5.5E+02		7.7E+00
				1.0E-04	X						1.36E+09		1	0.009	Dinitrotoluene, 4-Amino-2,6-	19406-51-0					7.8E+00	3.7E+02		7.7E+00
4.5E-01	X			9.0E-04	X						1.36E+09		1	0.1	Dinitrotoluene, Technical grade	25321-14-6	1.5E+02	5.5E+02		1.2E+02	7.0E+01	3.0E+02		5.7E+01
				1.0E-03	I						1.36E+09		1	0.1	Dinoseb	88-85-7					7.8E+01	3.3E+02		6.3E+01
1.0E-01	I	5.0E-06	I	3.0E-02	I	3.0E-02	I	V		1.16E+05	1.36E+09	3.96E+04	1		Dioxane, 1,4-	123-91-1	7.0E+02		2.2E+03	5.3E+02	2.3E+03		1.2E+03	8.1E+02
															Dioxins									
6.2E+03	I	1.3E+00	I								1.36E+09		1	0.03	~Hexachlorodibenzo-p-dioxin, Mixture	34465-46-8	1.1E-02	1.3E-01	2.9E+02	1.0E-02				
1.3E+05	C	3.8E+01	C	7.0E-10	I	4.0E-08	C	V			1.36E+09	1.96E+06	1	0.03	~TCDD, 2,3,7,8-	1746-01-6	5.3E-04	6.3E-03	1.4E-02	4.8E-04	5.5E-05	7.7E-04	8.2E-02	5.1E-05
				3.0E-02	I						1.36E+09		1	0.1	Diphenamid	957-51-7					2.3E+03	9.9E+03		1.9E+03
						4.0E-04	X	V			1.36E+09	8.06E+04	1		Diphenyl Ether	101-84-8							3.4E+01	3.4E+01
				8.0E-04	X						1.36E+09		1	0.1	Diphenyl Sulfone	127-63-9					6.3E+01	2.6E+02		5.1E+01
				1.0E-01	O						1.36E+09		1	0.1	Diphenylamine	122-39-4					7.8E+03	3.3E+04		6.3E+03
8.0E-01	I	2.2E-04	I								1.36E+09		1	0.1	Diphenylhydrazine, 1,2-	122-66-7	8.7E+01	3.1E+02	1.7E+06	6.8E+01				
				2.2E-03	I						1.36E+09		1	0.1	Diquat	2764-72-9					1.7E+02	7.3E+02		1.4E+02
7.4E+00	C	2.1E-03	C								1.36E+09		1	0.1	Direct Black 38	1937-37-7	9.4E+00	3.3E+01	1.8E+05	7.3E+00				
7.4E+00	C	2.1E-03	C								1.36E+09		1	0.1	Direct Blue 6	2602-46-2	9.4E+00	3.3E+01	1.8E+05	7.3E+00				
6.7E+00	C	1.9E-03	C								1.36E+09		1	0.1	Direct Brown 95	16071-86-6	1.0E+01	3.7E+01	2.0E+05	8.1E+00				
				4.0E-05	I						1.36E+09		1	0.1	Disulfoton	298-04-4					3.1E+00	1.3E+01		2.5E+00
				1.0E-02	I			V			1.36E+09	4.54E+04	1		Dithiane, 1,4-	505-29-3					7.8E+02			7.8E+02
				2.0E-03	I						1.36E+09		1	0.1	Diuron	330-54-1					1.6E+02	6.6E+02		1.3E+02
				2.0E-02	O						1.36E+09		1	0.1	Dodine	2439-10-3					1.6E+03	6.6E+03		1.3E+03
				5.0E-02	O			V			1.36E+09	1.17E+05	1		EPTC	759-94-4					3.9E+03			3.9E+03
				6.0E-03	I			V			1.36E+09	4.10E+05	1		Endosulfan	115-29-7					4.7E+02			4.7E+02
				6.0E-03	P						1.36E+09		1	0.1	Endosulfan Sulfate	1031-07-8					4.7E+02	2.0E+03		3.8E+02
				2.0E-02	I						1.36E+09		1	0.1	Endothall	145-73-3					1.6E+03	6.6E+03		1.3E+03
				3.0E-04	I						1.36E+09		1	0.1	Endrin	72-20-8					2.3E+01	9.9E+01		1.9E+01
9.9E-03	I	1.2E-06	I	6.0E-03	P	1.0E-03	I	V		1.05E+04	1.36E+09	1.89E+04	1		Epichlorohydrin	106-89-8	7.0E+03		4.4E+03	2.7E+03	4.7E+02		2.0E+01	1.9E+01

RMLs R=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																		Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
						2.0E-02	I	V		1.53E+04	1.36E+09	7.66E+03	1		Epoxybutane, 1,2-	106-88-7							1.6E+02	1.6E+02
				4.0E-02	P						1.36E+09		1	0.1	Ethanol, 2-(2-methoxyethoxy)-	111-77-3					3.1E+03	1.3E+04		2.5E+03
				5.0E-03	I						1.36E+09		1	0.1	Ethephon	16672-87-0					3.9E+02	1.6E+03		3.2E+02
				5.0E-04	I						1.36E+09		1	0.1	Ethion	563-12-2					3.9E+01	1.6E+02		3.2E+01
				1.0E-01	P	6.0E-02	P	V		2.38E+04	1.36E+09	6.15E+04	1		Ethoxyethanol Acetate, 2-	111-15-9					7.8E+03		3.8E+03	2.6E+03
				9.0E-02	P	4.0E-02	P	V		1.06E+05	1.36E+09	9.84E+04	1		Ethoxyethanol, 2-	110-80-5					7.0E+03		4.1E+03	2.6E+03
				7.0E-01	P	7.0E-02	P	V		1.08E+04	1.36E+09	8.62E+03	1		Ethyl Acetate	141-78-6					5.5E+04		6.3E+02	6.2E+02
				5.0E-03	P	8.0E-03	P	V		2.50E+03	1.36E+09	6.34E+03	1		Ethyl Acrylate	140-88-5					3.9E+02		5.3E+01	4.7E+01
						4.0E+00	P	V		2.12E+03	1.36E+09	1.29E+03	1		Ethyl Chloride (Chloroethane)	75-00-3							5.4E+03	5.4E+03
				2.0E-01	I			V		1.01E+04	1.36E+09	3.12E+03	1		Ethyl Ether	60-29-7					1.6E+04			1.6E+04
						3.0E-01	P	V		1.10E+03	1.36E+09	5.77E+03	1		Ethyl Methacrylate	97-63-2							1.8E+03	1.8E+03
		8.0E-08	I	1.0E+00	I	4.0E+01	I	V		2.87E+03	1.36E+09	3.67E+03	1		Ethyl Tertiary Butyl Ether (ETBE)	637-92-3			1.3E+04	1.3E+04	7.8E+04		1.5E+05	5.2E+04
				1.0E-05	I						1.36E+09		1	0.1	Ethyl-p-nitrophenyl Phosphonate	2104-64-5					7.8E-01	3.3E+00		6.3E-01
1.1E-02	C	2.5E-06	C	5.0E-02	P	1.0E+00	I	V		4.80E+02	1.36E+09	5.67E+03	1		Ethylbenzene	100-41-4	6.3E+03		6.4E+02	5.8E+02	3.9E+03		5.9E+03	2.4E+03
				7.0E-02	P						1.36E+09		1	0.1	Ethylene Cyanohydrin	109-78-4					5.5E+03	2.3E+04		4.4E+03
				9.0E-02	P			V		1.89E+05	1.36E+09	1.80E+05	1		Ethylene Diamine	107-15-3					7.0E+03			7.0E+03
				8.0E-01	A	4.0E-01	C				1.36E+09		1	0.1	Ethylene Glycol	107-21-1					6.3E+04	2.6E+05	5.7E+08	5.1E+04
				1.0E-01	I	1.6E+00	I				1.36E+09		1	0.1	Ethylene Glycol Monobutyl Ether	111-76-2					7.8E+03	3.3E+04	2.3E+09	6.3E+03
3.1E-01	C	3.0E-03	I			3.0E-02	C	V	M	1.21E+05	1.36E+09	6.09E+03	1		Ethylene Oxide	75-21-8	4.9E+01		2.1E-01	2.0E-01			1.9E+02	1.9E+02
4.5E-02	C	1.3E-05	C	8.0E-05	I						1.36E+09		1	0.1	Ethylene Thiourea	96-45-7	1.5E+03	5.5E+03	2.9E+07	1.2E+03	6.3E+00	2.6E+01		5.1E+00
6.5E+01	C	1.9E-02	C					V		1.54E+05	1.36E+09	2.39E+04	1		Ethyleneimine	151-56-4	1.1E+00		3.5E-01	2.7E-01				
				3.0E+00	I						1.36E+09		1	0.1	Ethylphthalyl Ethyl Glycolate	84-72-0					2.3E+05	9.9E+05		1.9E+05
				2.5E-04	I						1.36E+09		1	0.1	Fenamiphos	22224-92-6					2.0E+01	8.2E+01		1.6E+01
				2.5E-02	I						1.36E+09		1	0.1	Fenpropathrin	39515-41-8					2.0E+03	8.2E+03		1.6E+03
				2.5E-02	I						1.36E+09		1	0.1	Fenvalerate	51630-58-1					2.0E+03	8.2E+03		1.6E+03
				1.3E-02	I						1.36E+09		1	0.1	Fluometuron	2164-17-2					1.0E+03	4.3E+03		8.2E+02
				4.0E-02	C	1.3E-02	C				1.36E+09		1		Fluoride	16984-48-8					3.1E+03		1.8E+07	3.1E+03
				6.0E-02	I	1.3E-02	C				1.36E+09		1		Fluorine (Soluble Fluoride)	7782-41-4					4.7E+03		1.8E+07	4.7E+03
				8.0E-02	I						1.36E+09		1	0.1	Fluridone	59756-60-4					6.3E+03	2.6E+04		5.1E+03
				4.0E-02	O						1.36E+09		1	0.1	Flurprimidol	56425-91-3					3.1E+03	1.3E+04		2.5E+03
				2.0E-03	O						1.36E+09		1	0.1	Flusilazole	85509-19-9					1.6E+02	6.6E+02		1.3E+02
				5.0E-01	O						1.36E+09		1	0.1	Flutolanil	66332-96-5					3.9E+04	1.6E+05		3.2E+04
				1.0E-02	I						1.36E+09		1	0.1	Fluvalinate	69409-94-5					7.8E+02	3.3E+03		6.3E+02
				9.0E-02	O						1.36E+09		1	0.1	Folpet	133-07-3					7.0E+03	3.0E+04		5.7E+03
				1.0E-02	O						1.36E+09		1	0.1	Fomesafen	72178-02-0					7.8E+02	3.3E+03		6.3E+02
				2.0E-03	I						1.36E+09		1	0.1	Fonofos	944-22-9					1.6E+02	6.6E+02		1.3E+02
2.1E-02	C	1.3E-05	I	2.0E-01	I	9.8E-03	A	V		4.24E+04	1.36E+09	7.77E+04	1		Formaldehyde	50-00-0	3.3E+03		1.7E+03	1.1E+03	1.6E+04		8.0E+02	7.6E+02
				9.0E-01	P	3.0E-04	X	V		1.06E+05	1.36E+09	9.30E+04	1		Formic Acid	64-18-6					7.0E+04		2.9E+01	2.9E+01
				2.5E+00	O						1.36E+09		1	0.1	Fosetyl-AL	39148-24-8					2.0E+05	8.2E+05		1.6E+05
															Furans									
				1.0E-03	X			V			1.36E+09	1.56E+05	1		~Dibenzofuran	132-64-9					7.8E+01			7.8E+01
				1.0E-03	I			V		6.22E+03	1.36E+09	2.62E+03	1		~Furan	110-00-9					7.8E+01			7.8E+01
				9.0E-01	I	2.0E+00	I	V		1.65E+05	1.36E+09	1.20E+04	1		~Tetrahydrofuran	109-99-9					7.0E+04		2.5E+04	1.8E+04
3.8E+00	H										1.36E+09		1	0.1	Furazolidone	67-45-8	1.8E+01	6.5E+01		1.4E+01				
				3.0E-03	I	5.0E-02	H	V		1.01E+04	1.36E+09	4.86E+04	1		Furfural	98-01-1					2.3E+02		2.5E+03	2.1E+02
1.5E+00	C	4.3E-04	C								1.36E+09		1	0.1	Furium	531-82-8	4.6E+01	1.6E+02	8.9E+05	3.6E+01				
3.0E-02	I	8.6E-06	C								1.36E+09		1	0.1	Furmecyclox	60568-05-0	2.3E+03	8.2E+03	4.4E+07	1.8E+03				
				6.0E-03	O						1.36E+09		1	0.1	Glufosinate, Ammonium	77182-82-2					4.7E+02	2.0E+03		3.8E+02
				1.0E-01	A	8.0E-05	C				1.36E+09		1	0.1	Glutaraldehyde	111-30-8					7.8E+03	3.3E+04	1.1E+05	6.0E+03
				4.0E-04	I	1.0E-03	X	V		1.06E+05	1.36E+09	8.43E+04	1		Glycidaldehyde	765-34-4					3.1E+01		8.8E+01	2.3E+01
				1.0E-01	I						1.36E+09		1	0.1	Glyphosate	1071-83-6					7.8E+03	3.3E+04		6.3E+03
				1.0E-02	X			V			1.36E+09	1.45E+05	1		Guanidine	113-00-8					7.8E+02			7.8E+02
				2.0E-02	P						1.36E+09		1	0.1	Guanidine Chloride	50-01-1					1.6E+03	6.6E+03		1.3E+03
				3.0E-02	X						1.36E+09		1	0.1	Guanidine Nitrate	506-93-4					2.3E+03	9.9E+03		1.9E+03
				5.0E-05	I						1.36E+09		1	0.1	Haloxyfop, Methyl	69806-40-2					3.9E+00	1.6E+01		3.2E+00
4.5E+00	I	1.3E-03	I	1.0E-04	A			V			1.36E+09	4.79E+05	1		Heptachlor	76-44-8	1.5E+01		1.0E+02	1.3E+01	7.8E+00			7.8E+00
9.1E+00	I	2.6E-03	I	1.3E-05	I			V			1.36E+09	8.43E+05	1		Heptachlor Epoxide	1024-57-3	7.6E+00		9.1E+01	7.0E+00	1.0E+00			1.0E+00
						3.0E-03	X	V		2.09E+02	1.36E+09	7.80E+03	1		Heptanal, n-	111-71-7							2.4E+01	2.4E+01
				3.0E-04	X	4.0E-01	P	V		5.79E+01	1.36E+09	8.95E+02	1		Heptane, N-	142-82-5					2.3E+01		3.7E+02	2.2E+01
				2.0E-03	I			V			1.36E+09	3.80E+05	1		Hexabromobenzene	87-82-1					1.6E+02			1.6E+02
				2.0E-04	I						1.36E+09		1	0.1	Hexabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-153)	68631-49-2					1.6E+01	6.6E+01		1.3E+01
1.6E+00	I	4.6E-04	I	1.0E-05	P			V			1.36E+09	6.80E+04	1		Hexachlorobenzene	118-74-1	4.3E+01		4.1E+01	2.1E+01	7.8E-01			7.8E-01
7.8E-02	I	2.2E-05	I	1.0E-03	P			V		1.68E+01	1.36E+09	1.08E+04	1		Hexachlorobutadiene	87-68-3	8.9E+02		1.4E+02	1.2E+02	7.8E+01			7.8E+01
6.3E+00	I	1.8E-03	I	8.0E-03	A						1.36E+09		1	0.1	Hexachlorocyclohexane, Alpha-	319-84-6	1.1E+01	3.9E+01	2.1E+05	8.6E+00	6.3E+02	2.6E+03		5.1E+02
1.8E+00	I	5.3E-04	I								1.36E+09		1	0.1	Hexachlorocyclohexane, Beta-	319-85-7	3.9E+01	1.4E+02	7.2E+05	3.0E+01				
1.1E+00	C	3.1E-04	C	1.0E-05	A						1.36E+09		1	0.04	Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	6.3E+01	5.6E+02	1.2E+06	5.7E+01	7.8E-01	8.2E+00		7.1E-01
1.8E+00	I	5.1E-04	I								1.36E+09		1	0.1	Hexachlorocyclohexane, Technical	608-73-1	3.9E+01	1.4E+02	7.5E+05	3.0E+01				
				6.0E-03	I	2.0E-04	I	V		1.57E+01	1.36E+09	8.51E+03	1		Hexachlorocyclopentadiene	77-47-4					4.7E+02		1.8E+00	1.8E+00
4.0E-02	I	1.1E-05	C	7.0E-04	I	3.0E-02	I	V			1.36E+09	8.01E+03	1		Hexachloroethane	67-72-1	1.7E+03		2.0E+02	1.8E+02	5.5E+01		2.5E+02	4.5E+01
				3.0E-04	I						1.36E+09		1	0.1	Hexachlorophene	70-30-4					2.3E+01	9.9E+01		1.9E+01
8.0E-02	I			4.0E-03	I						1.36E+09		1	0.015	Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	8.7E+02	2.1E+04		8.3E+02	3.1E+02	8.8E+03		3.0E+02
						1.0E-05	I	V		3.39E+03	1.36E+09	3.00E+05	1		Hexamethylene Diisocyanate, 1,6-	822-06-0							3.1E+00	3.1E+00
						4.0E-04	C				1.36E+09		1	0.1	Hexamethylene diisocyanate biuret	4035-89-6							5.7E+05	5.7E+05
						4.0E-04	C				1.36E+09		1	0.1	Hexamethylene diisocyanate isocyanurate	3779-63-3							5.7E+05	5.7E+05
				4.0E-04	P						1.36E+09		1	0.1	Hexamethylphosphoramide	680-31-9					3.1E+01	1.3E+02		2.5E+01
		2.0E-07	X			6.0E-01	P	V		1.41E+02	1.36E+09	8.29E+02	1		Hexane, Commercial	E5241997			1.2E+03	1.2E+03			5.2E+02	5.2E+02
						7.0E-01	I	V		1.41E+02	1.36E+09	8.29E+02	1		Hexane, N-	110-54-3							6.1E+02	6.1E+02

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
				2.0E+00	P						1.36E+09		1	0.1	Hexanedioic Acid	124-04-9					1.6E+05	6.6E+05		1.3E+05
9.5E-03	P			7.0E-02	P	4.0E-04	P	V		2.74E+02	1.36E+09	3.62E+04	1		Hexanol, 1-,2-ethyl- (2-Ethyl-1-hexanol)	104-76-7	7.3E+03			7.3E+03	5.5E+03		1.5E+01	1.5E+01
				5.0E-03	I	3.0E-02	I	V		3.28E+03	1.36E+09	1.33E+04	1		Hexanone, 2-	591-78-6					3.9E+02		4.2E+02	2.0E+02
				3.3E-02	I						1.36E+09		1	0.1	Hexazinone	51235-04-2					2.6E+03	1.1E+04		2.1E+03
				2.5E-02	I						1.36E+09		1	0.1	Hexythiazox	78587-05-0					2.0E+03	8.2E+03		1.6E+03
				1.7E-02	O						1.36E+09		1	0.1	Hydramethylnon	67485-29-4					1.3E+03	5.6E+03		1.1E+03
3.0E+00	I	4.9E-03	I			3.0E-05	P	V		1.12E+05	1.36E+09	6.52E+04	1		Hydrazine	302-01-2	2.3E+01		3.7E+00	3.2E+00			2.0E+00	2.0E+00
3.0E+00	I	4.9E-03	I								1.36E+09		1		Hydrazine Sulfate	10034-93-2	2.3E+01		7.8E+04	2.3E+01				
						2.0E-02	I	V			1.36E+09		1		Hydrogen Chloride	7647-01-0							2.8E+07	2.8E+07
				4.0E-02	C	1.4E-02	C	V			1.36E+09		1		Hydrogen Fluoride	7664-39-3					3.1E+03		2.0E+07	3.1E+03
						2.0E-03	I	V			1.36E+09		1		Hydrogen Sulfide	7783-06-4							2.8E+06	2.8E+06
6.0E-02	P			4.0E-02	P						1.36E+09		1	0.1	Hydroquinone	123-31-9	1.2E+03	4.1E+03		9.0E+02	3.1E+03	1.3E+04		2.5E+03
6.1E-02	O			1.1E-01	O						1.36E+09		1	0.1	Imazalil	35554-44-0	1.1E+03	4.0E+03		8.9E+02	8.4E+03	3.6E+04		6.8E+03
				2.5E-01	I						1.36E+09		1	0.1	Imazaquin	81335-37-7					2.0E+04	8.2E+04		1.6E+04
				2.5E+00	O						1.36E+09		1	0.1	Imazethapyr	81335-77-5					2.0E+05	8.2E+05		1.6E+05
				1.0E-02	A						1.36E+09		1		Iodine	7553-56-2					7.8E+02			7.8E+02
				4.0E-02	I						1.36E+09		1	0.1	Iprodione	36734-19-7					3.1E+03	1.3E+04		2.5E+03
				7.0E-01	P						1.36E+09		1		Iron	7439-89-6					5.5E+04			5.5E+04
				3.0E-01	I			V		1.00E+04	1.36E+09	2.81E+04	1		Isobutyl Alcohol	78-83-1					2.3E+04			2.3E+04
9.5E-04	I			2.0E-01	I	2.0E+00	C				1.36E+09		1	0.1	Isophorone	78-59-1	7.3E+04	2.6E+05		5.7E+04	1.6E+04	6.6E+04	2.8E+09	1.3E+04
				1.5E-02	I			V			1.36E+09	4.20E+05	1		Isopropalin	33820-53-0					1.2E+03			1.2E+03
				2.0E+00	P	2.0E-01	P	V		1.09E+05	1.36E+09	2.77E+04	1		Isopropanol	67-63-0					1.6E+05		5.8E+03	5.6E+03
				1.0E-01	I						1.36E+09		1	0.1	Isopropyl Methyl Phosphonic Acid	1832-54-8					7.8E+03	3.3E+04		6.3E+03
				5.0E-02	I						1.36E+09		1	0.1	Isoxaben	82558-50-7					3.9E+03	1.6E+04		3.2E+03
						3.0E-01	A	V			1.36E+09		1		JP-7	E1737665							4.3E+08	4.3E+08
				8.0E-03	O						1.36E+09		1	0.1	Lactofen	77501-63-4					6.3E+02	2.6E+03		5.1E+02
				2.0E-04	X						1.36E+09		1	0.1	Lactonitrile	78-97-7					1.6E+01	6.6E+01		1.3E+01
				5.0E-05	P						1.36E+09		1		Lanthanum	7439-91-0					3.9E+00			3.9E+00
				2.1E-05	P						1.36E+09		1	0.1	Lanthanum Acetate Hydrate	100587-90-4					1.6E+00	6.9E+00		1.3E+00
				1.9E-05	P						1.36E+09		1		Lanthanum Chloride Heptahydrate	10025-84-0					1.5E+00			1.5E+00
				2.8E-05	P						1.36E+09		1		Lanthanum Chloride, Anhydrous	10099-58-8					2.2E+00			2.2E+00
				1.6E-05	P						1.36E+09		1		Lanthanum Nitrate Hexahydrate	10277-43-7					1.3E+00			1.3E+00
															Lead Compounds									
8.5E-03	C	1.2E-05	C								1.36E+09		1		~Lead Phosphate	7446-27-7	8.2E+03		3.2E+07	8.2E+03				
2.1E-01	C	8.0E-05	C								1.36E+09		1	0.1	~Lead acetate	301-04-2	3.3E+02	1.2E+03	4.8E+06	2.6E+02				
											1.36E+09		1		~Lead and Compounds	7439-92-1								4.0E+02
3.8E-02	C	1.1E-05	C								1.36E+09		1	0.1	~Lead subacetate	1335-32-6	1.8E+03	6.5E+03	3.5E+07	1.4E+03				
				1.0E-07	I			V		2.43E+00	1.36E+09	1.91E+03	1		~Tetraethyl Lead	78-00-2					7.8E-03			7.8E-03
				5.0E-06	P			V		3.83E+02	1.36E+09	2.55E+04	1		Lewisite	541-25-3					3.9E-01			3.9E-01
				7.7E-03	O						1.36E+09		1	0.1	Linuron	330-55-2					6.0E+02	2.5E+03		4.9E+02
				2.0E-03	P						1.36E+09		1		Lithium	7439-93-2					1.6E+02			1.6E+02
				5.0E-04	I						1.36E+09		1	0.1	MCPA	94-74-6					3.9E+01	1.6E+02		3.2E+01
				4.4E-03	O						1.36E+09		1	0.1	MCPB	94-81-5					3.4E+02	1.5E+03		2.8E+02
				1.0E-03	I						1.36E+09		1	0.1	MCPP	93-65-2					7.8E+01	3.3E+02		6.3E+01
				2.0E-02	I						1.36E+09		1	0.1	Malathion	121-75-5					1.6E+03	6.6E+03		1.3E+03
				1.0E-01	I	7.0E-04	C				1.36E+09		1	0.1	Maleic Anhydride	108-31-6					7.8E+03	3.3E+04	9.9E+05	6.3E+03
				5.0E-01	I						1.36E+09		1	0.1	Maleic Hydrazide	123-33-1					3.9E+04	1.6E+05		3.2E+04
				1.0E-04	P						1.36E+09		1	0.1	Malononitrile	109-77-3					7.8E+00	3.3E+01		6.3E+00
				3.0E-02	H						1.36E+09		1	0.1	Mancozeb	8018-01-7					2.3E+03	9.9E+03		1.9E+03
				5.0E-03	I						1.36E+09		1	0.1	Maneb	12427-38-2					3.9E+02	1.6E+03		3.2E+02
				1.4E-01	I	5.0E-05	I						1		Manganese (Diet)	7439-96-5								
				2.4E-02	G	5.0E-05	I				1.36E+09		0.04		Manganese (Non-diet)	7439-96-5					1.9E+03		7.1E+04	1.8E+03
				9.0E-05	H						1.36E+09		1	0.1	Mephosfolan	950-10-7					7.0E+00	3.0E+01		5.7E+00
				3.0E-02	I						1.36E+09		1	0.1	Mepiquat Chloride	24307-26-4					2.3E+03	9.9E+03		1.9E+03
1.1E-02	P			4.0E-03	P						1.36E+09		1	0.1	Mercaptobenzothiazole, 2-	149-30-4	6.3E+03	2.2E+04		4.9E+03	3.1E+02	1.3E+03		2.5E+02
															Mercury Compounds									
				3.0E-04	I	3.0E-04	G				1.36E+09		0.07		~Mercuric Chloride (and other Mercury salts)	7487-94-7					2.3E+01		4.3E+05	2.3E+01
						3.0E-04	I	V		3.13E+00	1.36E+09	3.47E+04	1		~Mercury (elemental)	7439-97-6							1.1E+01	1.1E+01
				1.0E-04	I						1.36E+09		1		~Methyl Mercury	22967-92-6					7.8E+00			7.8E+00
				8.0E-05	I						1.36E+09		1	0.1	~Phenylmercuric Acetate	62-38-4					6.3E+00	2.6E+01		5.1E+00
				3.0E-05	I			V			1.36E+09	1.94E+06	1		Merphos	150-50-5					2.3E+00			2.3E+00
				6.0E-02	I						1.36E+09		1	0.1	Metalaxyl	57837-19-1					4.7E+03	2.0E+04		3.8E+03
				1.0E-04	I	3.0E-02	P	V		4.58E+03	1.36E+09	6.79E+03	1		Methacrylonitrile	126-98-7					7.8E+00		2.1E+02	7.5E+00
				5.0E-05	I						1.36E+09		1	0.1	Methamidophos	10265-92-6					3.9E+00	1.6E+01		3.2E+00
				2.0E+00	I	2.0E+01	I	V		1.06E+05	1.36E+09	2.90E+04	1		Methanol	67-56-1					1.6E+05		6.1E+05	1.2E+05
				1.5E-03	O						1.36E+09		1	0.1	Methidathion	950-37-8					1.2E+02	4.9E+02		9.5E+01
				2.5E-02	I						1.36E+09		1	0.1	Methomyl	16752-77-5					2.0E+03	8.2E+03		1.6E+03
4.9E-02	C										1.36E+09		1	0.1	Methoxy-5-nitroaniline, 2-	99-59-2	1.4E+03	5.0E+03		1.1E+03				
				5.0E-03	I						1.36E+09		1	0.1	Methoxychlor	72-43-5					3.9E+02	1.6E+03		3.2E+02
				8.0E-03	P	1.0E-03	P	V		1.15E+05	1.36E+09	1.24E+05	1		Methoxyethanol Acetate, 2-	110-49-6					6.3E+02		1.3E+02	1.1E+02
				5.0E-03	P	7.0E-03	P	V		1.06E+05	1.36E+09	1.01E+05	1		Methoxyethanol, 2-	109-86-4					3.9E+02		7.4E+02	2.6E+02
				1.0E+00	X			V		2.90E+04	1.36E+09	8.12E+03	1		Methyl Acetate	79-20-9					7.8E+04			7.8E+04
						2.0E-02	P	V		6.75E+03	1.36E+09	6.97E+03	1		Methyl Acrylate	96-33-3							1.5E+02	1.5E+02
				6.0E-01	I	5.0E+00	I	V		2.84E+04	1.36E+09	1.22E+04	1		Methyl Ethyl Ketone (2-Butanone)	78-93-3					4.7E+04		6.4E+04	2.7E+04
		1.0E-03	X	1.0E-03	P	2.0E-05	X	V		1.80E+05	1.36E+09	5.04E+04	1		Methyl Hydrazine	60-34-4			1.4E+01	1.4E+01	7.8E+01		1.1E+00	1.0E+00
						3.0E+00	I	V		3.36E+03	1.36E+09	1.06E+04	1		Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1							3.3E+04	3.3E+04
						1.0E-03	C	V		1.01E+04	1.36E+09	4.42E+03	1		Methyl Isocyanate	624-83-9							4.6E+00	4.6E+00

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
				1.4E+00	I	7.0E-01	I	V		2.36E+03	1.36E+09	6.33E+03	1		Methyl Methacrylate	80-62-6					1.1E+05		4.6E+03	4.4E+03
				2.5E-04	I						1.36E+09		1	0.1	Methyl Parathion	298-00-0					2.0E+01	8.2E+01		1.6E+01
				6.0E-02	X						1.36E+09		1	0.1	Methyl Phosphonic Acid	993-13-5					4.7E+03	2.0E+04		3.8E+03
				6.0E-03	H	4.0E-02	H	V		3.93E+02	1.36E+09	2.43E+04	1		Methyl Styrene (Mixed Isomers)	25013-15-4					4.7E+02		1.0E+03	3.2E+02
9.9E-02	C	2.8E-05	C								1.36E+09		1	0.1	Methyl methanesulfonate	66-27-3	7.0E+02	2.5E+03	1.4E+07	5.5E+02				
1.8E-03	C	2.6E-07	C			3.0E+00	I	V		8.87E+03	1.36E+09	4.90E+03	1		Methyl tert-Butyl Ether (MTBE)	1634-04-4	3.9E+04		5.3E+03	4.7E+03			1.5E+04	1.5E+04
				3.0E-04	X						1.36E+09		1	0.1	Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2					2.3E+01	9.9E+01		1.9E+01
						3.0E+00	X	V		2.45E+03	1.36E+09	1.72E+04	1		Methyl-2-Pentanol, 4-	108-11-2							5.4E+04	5.4E+04
9.0E-03	P			2.0E-02	X						1.36E+09		1	0.1	Methyl-5-Nitroaniline, 2-	99-55-8	7.7E+03	2.7E+04		6.0E+03	1.6E+03	6.6E+03		1.3E+03
8.3E+00	C	2.4E-03	C								1.36E+09		1	0.1	Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	8.4E+00	3.0E+01	1.6E+05	6.5E+00				
1.3E-01	C	3.7E-05	C								1.36E+09		1	0.1	Methylaniline Hydrochloride, 2-	636-21-5	5.3E+02	1.9E+03	1.0E+07	4.2E+02				
				1.0E-02	A						1.36E+09		1	0.1	Methylarsonic acid	124-58-3					7.8E+02	3.3E+03		6.3E+02
				2.0E-04	X						1.36E+09		1	0.1	Methylbenzene,1-4-diamine monohydrochloride, 2-	74612-12-7					1.6E+01	6.6E+01		1.3E+01
1.0E-01	X			3.0E-04	X						1.36E+09		1	0.1	Methylbenzene-1,4-diamine sulfate, 2-	615-50-9	7.0E+02	2.5E+03		5.4E+02	2.3E+01	9.9E+01		1.9E+01
2.2E+01	C	6.3E-03	C						M		1.36E+09		1	0.1	Methylcholanthrene, 3-	56-49-5	7.0E-01	2.7E+00	2.2E+04	5.5E-01				
2.0E-03	I	1.0E-08	I	6.0E-03	I	6.0E-01	I	V	M	3.32E+03	1.36E+09	2.19E+03	1		Methylene Chloride	75-09-2	7.7E+03		2.2E+04	5.7E+03	4.7E+02		1.4E+03	3.5E+02
1.0E-01	P	4.3E-04	C	2.0E-03	P				M		1.36E+09		1	0.1	Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	1.5E+02	6.0E+02	3.2E+05	1.2E+02	1.6E+02	6.6E+02		1.3E+02
4.6E-02	I	1.3E-05	C								1.36E+09		1	0.1	Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	1.5E+03	5.4E+03	2.9E+07	1.2E+03				
1.6E+00	C	4.6E-04	C			2.0E-02	C				1.36E+09		1	0.1	Methylenebisbenzenamine, 4,4'-	101-77-9	4.3E+01	1.5E+02	8.3E+05	3.4E+01			2.8E+07	2.8E+07
						6.0E-04	I				1.36E+09		1	0.1	Methylenediphenyl Diisocyanate	101-68-8							8.5E+05	8.5E+05
				7.0E-02	H			V		5.00E+02	1.36E+09	1.28E+04	1		Methylstyrene, Alpha-	98-83-9					5.5E+03			5.5E+03
				1.5E-01	I						1.36E+09		1	0.1	Metolachlor	51218-45-2					1.2E+04	4.9E+04		9.5E+03
				2.5E-02	I						1.36E+09		1	0.1	Metribuzin	21087-64-9					2.0E+03	8.2E+03		1.6E+03
				2.5E-01	I						1.36E+09		1	0.1	Metsulfuron-methyl	74223-64-6					2.0E+04	8.2E+04		1.6E+04
		4.5E-06	X	1.0E-02	X	1.0E-01	P	V		6.86E+00	1.36E+09	1.04E+03	1		Midrange Aliphatic Hydrocarbon Streams	E1790669			6.5E+01	6.5E+01	7.8E+02		1.1E+02	9.6E+01
				3.0E+00	P			V		3.42E-01	1.36E+09	1.38E+03	1		Mineral oils	8012-95-1					2.3E+05			2.3E+05
1.8E+01	C	5.1E-03	C	2.0E-04	I			V			1.36E+09	8.58E+05	1		Mirex	2385-85-5	3.9E+00		4.7E+01	3.6E+00	1.6E+01			1.6E+01
				2.0E-03	I						1.36E+09		1	0.1	Molinate	2212-67-1					1.6E+02	6.6E+02		1.3E+02
				5.0E-03	I	2.0E-03	A				1.36E+09		1		Molybdenum	7439-98-7					3.9E+02		2.8E+06	3.9E+02
				1.0E-01	I						1.36E+09		1		Monochloramine	10599-90-3					7.8E+03			7.8E+03
				2.0E-03	P						1.36E+09		1	0.1	Monomethylaniline	100-61-8					1.6E+02	6.6E+02		1.3E+02
				2.5E-02	I						1.36E+09		1	0.1	Myclobutanil	88671-89-0					2.0E+03	8.2E+03		1.6E+03
				3.0E-04	X						1.36E+09		1	0.1	N,N'-Diphenyl-1,4-benzenediamine	74-31-7					2.3E+01	9.9E+01		1.9E+01
				2.0E-03	I			V			1.36E+09	5.70E+04	1		Naled	300-76-5					1.6E+02			1.6E+02
				3.0E-02	X	1.0E-01	P	V			1.36E+09		1		Naphtha, High Flash Aromatic (HFAN)	64742-95-6					2.3E+03		1.4E+08	2.3E+03
1.8E+00	C	0.0E+00	C								1.36E+09		1	0.1	Naphthylamine, 2-	91-59-8	3.9E+01	1.4E+02		3.0E+01				
				1.2E-01	O						1.36E+09		1	0.1	Napropamide	15299-99-7					9.4E+03	4.0E+04		7.6E+03
		2.6E-04	C	1.1E-02	C	1.4E-05	C				1.36E+09		1	0.1	Nickel Acetate	373-02-4			1.5E+06	1.5E+06	8.6E+02	3.6E+03	2.0E+04	6.7E+02
		2.6E-04	C	1.1E-02	C	1.4E-05	C				1.36E+09		1	0.1	Nickel Carbonate	3333-67-3			1.5E+06	1.5E+06	8.6E+02	3.6E+03	2.0E+04	6.7E+02
		2.6E-04	C	1.1E-02	C	1.4E-05	C	V			1.36E+09		1		Nickel Carbonyl	13463-39-3			1.5E+06	1.5E+06	8.6E+02		2.0E+04	8.2E+02
		2.6E-04	C	1.1E-02	C	1.4E-05	C				1.36E+09		0.04		Nickel Hydroxide	12054-48-7			1.5E+06	1.5E+06	8.6E+02		2.0E+04	8.2E+02
		2.6E-04	C	1.1E-02	C	2.0E-05	C				1.36E+09		0.04		Nickel Oxide	1313-99-1			1.5E+06	1.5E+06	8.6E+02		2.8E+04	8.4E+02
		2.4E-04	I	1.1E-02	C	1.4E-05	C				1.36E+09		0.04		Nickel Refinery Dust	E715532			1.6E+06	1.6E+06	8.6E+02		2.0E+04	8.2E+02
		2.6E-04	C	2.0E-02	I	9.0E-05	A				1.36E+09		0.04		Nickel Soluble Salts	7440-02-0			1.5E+06	1.5E+06	1.6E+03		1.3E+05	1.5E+03
1.7E+00	C	4.8E-04	I	1.1E-02	C	1.4E-05	C				1.36E+09		0.04		Nickel Subsulfide	12035-72-2	4.1E+01		8.0E+05	4.1E+01	8.6E+02		2.0E+04	8.2E+02
9.1E-01	C	2.6E-04	C	1.1E-02	C	1.4E-05	C				1.36E+09		1	0.1	Nickelocene	1271-28-9	7.6E+01	2.7E+02	1.5E+06	6.0E+01	8.6E+02	3.6E+03	2.0E+04	6.7E+02
				1.6E+00	I						1.36E+09		1		Nitrate (measured as nitrogen)	14797-55-8					1.3E+05			1.3E+05
											1.36E+09		1		Nitrate + Nitrite (measured as nitrogen)	E701177								
				1.0E-01	I						1.36E+09		1		Nitrite (measured as nitrogen)	14797-65-0					7.8E+03			7.8E+03
				1.0E-02	X	5.0E-05	X				1.36E+09		1	0.1	Nitroaniline, 2-	88-74-4					7.8E+02	3.3E+03	7.1E+04	6.3E+02
2.0E-02	P			4.0E-03	P	6.0E-03	P				1.36E+09		1	0.1	Nitroaniline, 4-	100-01-6	3.5E+03	1.2E+04		2.7E+03	3.1E+02	1.3E+03	8.5E+06	2.5E+02
		4.0E-05	I	2.0E-03	I	9.0E-03	I	V		3.05E+03	1.36E+09	7.32E+04	1		Nitrobenzene	98-95-3			5.1E+02	5.1E+02	1.6E+02		6.9E+02	1.3E+02
				3.0E+03	P						1.36E+09		1	0.1	Nitrocellulose	9004-70-0					2.3E+08	9.9E+08		1.9E+08
				7.0E-02	H						1.36E+09		1	0.1	Nitrofurantoin	67-20-9					5.5E+03	2.3E+04		4.4E+03
1.3E+00	C	3.7E-04	C								1.36E+09		1	0.1	Nitrofurazone	59-87-0	5.3E+01	1.9E+02	1.0E+06	4.2E+01				
1.7E-02	P			1.0E-04	P						1.36E+09		1	0.1	Nitroglycerin	55-63-0	4.1E+03	1.5E+04		3.2E+03	7.8E+00	3.3E+01		6.3E+00
				1.0E-01	I						1.36E+09		1	0.1	Nitroguanidine	556-88-7					7.8E+03	3.3E+04		6.3E+03
		8.8E-06	P			5.0E-03	P	V		1.80E+04	1.36E+09	1.69E+04	1		Nitromethane	75-52-5			5.4E+02	5.4E+02			8.8E+01	8.8E+01
		5.8E-04	X			2.0E-02	I	V		4.86E+03	1.36E+09	1.31E+04	1		Nitropropane, 2-	79-46-9			6.4E+00	6.4E+00			2.7E+02	2.7E+02
2.7E+01	C	7.7E-03	C						M		1.36E+09		1	0.1	Nitroso-N-ethylurea, N-	759-73-9	5.7E-01	2.2E+00	1.8E+04	4.5E-01				
1.2E+02	C	3.4E-02	C						M		1.36E+09		1	0.1	Nitroso-N-methylurea, N-	684-93-5	1.3E-01	5.0E-01	4.1E+03	1.0E-01				
5.4E+00	I	1.6E-03	I					V			1.36E+09	2.43E+05	1		Nitroso-di-N-butylamine, N-	924-16-3	1.3E+01		4.3E+01	9.9E+00				
7.0E+00	I	2.0E-03	C								1.36E+09		1	0.1	Nitroso-di-N-propylamine, N-	621-64-7	9.9E+00	3.5E+01	1.9E+05	7.8E+00				
2.8E+00	I	8.0E-04	C								1.36E+09		1	0.1	Nitrosodiethanolamine, N-	1116-54-7	2.5E+01	8.8E+01	4.8E+05	1.9E+01				
1.5E+02	I	4.3E-02	I						M		1.36E+09		1	0.1	Nitrosodiethylamine, N-	55-18-5	1.0E-01	4.0E-01	3.2E+03	8.1E-02				
5.1E+01	I	1.4E-02	I	8.0E-06	P	4.0E-05	X	V	M	2.37E+05	1.36E+09	8.23E+04	1		Nitrosodimethylamine, N-	62-75-9	3.0E-01		6.0E-01	2.0E-01	6.3E-01		3.4E+00	5.3E-01
4.9E-03	I	2.6E-06	C								1.36E+09		1	0.1	Nitrosodiphenylamine, N-	86-30-6	1.4E+04	5.0E+04	1.5E+08	1.1E+04				
2.2E+01	I	6.3E-03	C					V		1.08E+05	1.36E+09	1.21E+05	1		Nitrosomethylethylamine, N-	10595-95-6	3.2E+00		5.4E+00	2.0E+00				
6.7E+00	C	1.9E-03	C								1.36E+09		1	0.1	Nitrosomorpholine [N-]	59-89-2	1.0E+01	3.7E+01	2.0E+05	8.1E+00				
9.4E+00	C	2.7E-03	C								1.36E+09		1	0.1	Nitrosopiperidine [N-]	100-75-4	7.4E+00	2.6E+01	1.4E+05	5.8E+00				
2.1E+00	I	6.1E-04	I								1.36E+09		1	0.1	Nitrosopyrrolidine, N-	930-55-2	3.3E+01	1.2E+02	6.3E+05	2.6E+01				
				1.0E-04	X						1.36E+09		1	0.1	Nitrotoluene, m-	99-08-1					7.8E+00	3.3E+01		6.3E+00
2.2E-01	P			9.0E-04	P			V		1.51E+03	1.36E+09	1.37E+05	1		Nitrotoluene, o-	88-72-2	3.2E+02			3.2E+02	7.0E+01			7.0E+01
1.6E-02	P			4.0E-03	P						1.36E+09		1	0.1	Nitrotoluene, p-	99-99-0	4.3E+03	1.5E+04		3.4E+03	3.1E+02	1.3E+03		2.5E+02
				3.0E-04	X	2.0E-02	P	V		6.86E+00	1.36E+09	1.04E+03	1		Nonane, n-	111-84-2					2.3E+01		2.2E+01	1.1E+01
				1.5E-03	O						1.36E+09		1	0.1	Norflurazon	27314-13-2					1.2E+02	4.9E+02		9.5E+01
				3.0E-03	I						1.36E+09		1	0.1	Octabromodiphenyl Ether	32536-52-0					2.3E+02	9.9E+02		1.9E+02

RMLS TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
				5.0E-02	I						1.36E+09		1	0.006	Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0					3.9E+03	2.7E+05		3.9E+03
				2.0E-03	H						1.36E+09		1	0.1	Octamethylpyrophosphoramide	152-16-9					1.6E+02	6.6E+02		1.3E+02
7.8E-03	O			1.9E-01	O						1.36E+09		1	0.1	Oryzalin	19044-88-3	8.9E+03	3.2E+04		7.0E+03	1.5E+04	6.3E+04		1.2E+04
				5.0E-03	I						1.36E+09		1	0.1	Oxadiazon	19666-30-9					3.9E+02	1.6E+03		3.2E+02
				2.5E-02	I						1.36E+09		1	0.1	Oxamyl	23135-22-0					2.0E+03	8.2E+03		1.6E+03
7.3E-02	O			4.0E-02	O						1.36E+09		1	0.1	Oxyfluorfen	42874-03-3	9.5E+02	3.4E+03		7.4E+02	3.1E+03	1.3E+04		2.5E+03
				1.3E-02	I						1.36E+09		1	0.1	Paclobutrazol	76738-62-0					1.0E+03	4.3E+03		8.2E+02
				4.5E-03	I						1.36E+09		1	0.1	Paraquat Dichloride	1910-42-5					3.5E+02	1.5E+03		2.8E+02
				6.0E-03	H						1.36E+09		1	0.1	Parathion	56-38-2					4.7E+02	2.0E+03		3.8E+02
				5.0E-02	H			V			1.36E+09	4.49E+04	1		Pebulate	1114-71-2					3.9E+03			3.9E+03
				3.0E-01	O						1.36E+09		1	0.1	Pendimethalin	40487-42-1					2.3E+04	9.9E+04		1.9E+04
				2.0E-03	I			V		3.12E-01	1.36E+09	5.13E+05	1		Pentabromodiphenyl Ether	32534-81-9					1.6E+02			1.6E+02
				1.0E-04	I						1.36E+09		1	0.1	Pentabromodiphenyl ether, 2,2',4,4',5- (BDE-99)	60348-60-9					7.8E+00	3.3E+01		6.3E+00
				8.0E-04	I			V			1.36E+09	8.12E+04	1		Pentachlorobenzene	608-93-5					6.3E+01			6.3E+01
9.0E-02	P							V		4.57E+02	1.36E+09	9.65E+03	1		Pentachloroethane	76-01-7	7.7E+02			7.7E+02				
2.6E-01	H			3.0E-03	I			V			1.36E+09	4.32E+05	1		Pentachloronitrobenzene	82-68-8	2.7E+02			2.7E+02	2.3E+02			2.3E+02
4.0E-01	I	5.1E-06	C	5.0E-03	I						1.36E+09		1	0.25	Pentachlorophenol	87-86-5	1.7E+02	2.5E+02	7.5E+07	1.0E+02	3.9E+02	6.6E+02		2.5E+02
4.3E-03	X			9.0E-03	P						1.36E+09		1	0.1	Pentaerythritol tetranitrate (PETN)	78-11-5	1.6E+04	5.7E+04		1.3E+04	7.0E+02	3.0E+03		5.7E+02
				1.0E-04	X						1.36E+09		1	0.1	Pentamethylphosphoramide (PMPA)	10159-46-3					7.8E+00	3.3E+01		6.3E+00
						1.0E+00	P	V		3.88E+02	1.36E+09	7.79E+02	1		Pentane, n-	109-66-0							8.1E+02	8.1E+02
															Per- and Polyfluoroalkyl Substances (PFAS)									
				3.0E-06	D						1.36E+09		1	0.1	~Ammonium perfluoro-2-methyl-3-oxahexanoate	62037-80-3					2.3E-01	9.9E-01		1.9E-01
				3.0E-06	D			V			1.36E+09		1		~Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6					2.3E-01			2.3E-01
				3.0E-04	P						1.36E+09		1	0.1	~Perfluorobutanesulfonate	45187-15-3					2.3E+01	9.9E+01		1.9E+01
				3.0E-04	P						1.36E+09		1	0.1	~Perfluorobutanesulfonic acid (PFBS)	375-73-5					2.3E+01	9.9E+01		1.9E+01
				2.0E-05	A						1.36E+09		1	0.1	~Perfluorohexanesulfonate	108427-53-8					1.6E+00	6.6E+00		1.3E+00
				2.0E-05	A						1.36E+09		1	0.1	~Perfluorohexanesulfonic acid (PFHxS)	355-46-4					1.6E+00	6.6E+00		1.3E+00
				3.0E-06	A						1.36E+09		1	0.1	~Perfluorononanoate	72007-68-2					2.3E-01	9.9E-01		1.9E-01
				3.0E-06	A						1.36E+09		1	0.1	~Perfluorononanoic acid (PFNA)	375-95-1					2.3E-01	9.9E-01		1.9E-01
				2.0E-06	A						1.36E+09		1	0.1	~Perfluorooctanesulfonate	45298-90-6					1.6E-01	6.6E-01		1.3E-01
				2.0E-06	A						1.36E+09		1	0.1	~Perfluorooctanesulfonic acid (PFOS)	1763-23-1					1.6E-01	6.6E-01		1.3E-01
7.0E-02	D			3.0E-06	A						1.36E+09		1	0.1	~Perfluorooctanoate	45285-51-6	9.9E+02	3.5E+03		7.8E+02	2.3E-01	9.9E-01		1.9E-01
7.0E-02	D			3.0E-06	A						1.36E+09		1	0.1	~Perfluorooctanoic acid (PFOA)	335-67-1	9.9E+02	3.5E+03		7.8E+02	2.3E-01	9.9E-01		1.9E-01
				3.0E-04	P						1.36E+09		1	0.1	~Potassium perfluorobutanesulfonate	29420-49-3					2.3E+01	9.9E+01		1.9E+01
				2.0E-06	A						1.36E+09		1	0.1	~Potassium perfluorooctanesulfonate	2795-39-3					1.6E-01	6.6E-01		1.3E-01
															Perchlorates									
				7.0E-04	I						1.36E+09		1		~Ammonium Perchlorate	7790-98-9					5.5E+01			5.5E+01
				7.0E-04	I						1.36E+09		1		~Lithium Perchlorate	7791-03-9					5.5E+01			5.5E+01
				7.0E-04	I						1.36E+09		1		~Perchlorate and Perchlorate Salts	14797-73-0					5.5E+01			5.5E+01
				7.0E-04	I						1.36E+09		1		~Potassium Perchlorate	7778-74-7					5.5E+01			5.5E+01
				7.0E-04	I						1.36E+09		1		~Sodium Perchlorate	7601-89-0					5.5E+01			5.5E+01
				5.0E-02	I						1.36E+09		1	0.1	Permethrin	52645-53-1					3.9E+03	1.6E+04		3.2E+03
2.2E-03	C	6.3E-07	C								1.36E+09		1	0.1	Phenacetin	62-44-2	3.2E+04	1.1E+05	6.1E+08	2.5E+04				
				2.4E-01	O						1.36E+09		1	0.1	Phenmedipham	13684-63-4					1.9E+04	7.9E+04		1.5E+04
				3.0E-01	I	2.0E-01	C				1.36E+09		1	0.1	Phenol	108-95-2					2.3E+04	9.9E+04	2.8E+08	1.9E+04
				4.0E-03	I						1.36E+09		1	0.1	Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1					3.1E+02	1.3E+03		2.5E+02
				5.0E-04	X						1.36E+09		1	0.1	Phenothiazine	92-84-2					3.9E+01	1.6E+02		3.2E+01
				2.0E-04	X			V		1.29E+02	1.36E+09	7.06E+03	1		Phenyl Isothiocyanate	103-72-0					1.6E+01			1.6E+01
				6.0E-03	I						1.36E+09		1	0.1	Phenylenediamine, m-	108-45-2					4.7E+02	2.0E+03		3.8E+02
1.2E-01	P			4.0E-03	P						1.36E+09		1	0.1	Phenylenediamine, o-	95-54-5	5.8E+02	2.1E+03		4.5E+02	3.1E+02	1.3E+03		2.5E+02
				1.0E-03	X						1.36E+09		1	0.1	Phenylenediamine, p-	106-50-3					7.8E+01	3.3E+02		6.3E+01
1.9E-03	H										1.36E+09		1	0.1	Phenylphenol, 2-	90-43-7	3.6E+04	1.3E+05		2.8E+04				
				2.0E-04	H						1.36E+09		1	0.1	Phorate	298-02-2					1.6E+01	6.6E+01		1.3E+01
						3.0E-04	I	V		1.61E+03	1.36E+09	9.81E+02	1		Phosgene	75-44-5							3.1E-01	3.1E-01
				2.0E-02	I						1.36E+09		1	0.1	Phosmet	732-11-6					1.6E+03	6.6E+03		1.3E+03
				3.0E-04	I	3.0E-04	I	V			1.36E+09		1		Phosphine	7803-51-2					2.3E+01		4.3E+05	2.3E+01
						1.0E-02	I				1.36E+09		1		Phosphoric Acid	7664-38-2							1.4E+07	1.4E+07
				2.0E-05	I			V			1.36E+09	6.92E+03	1		Phosphorus, White	7723-14-0					1.6E+00			1.6E+00
															Phthalates									
1.4E-02	I	2.4E-06	C	2.0E-02	I						1.36E+09		1	0.1	~Bis(2-ethylhexyl)phthalate	117-81-7	5.0E+03	1.8E+04	1.6E+08	3.9E+03	1.6E+03	6.6E+03		1.3E+03
1.9E-03	P			2.0E-01	I						1.36E+09		1	0.1	~Butyl Benzyl Phthalate	85-68-7	3.7E+04	1.3E+05		2.9E+04	1.6E+04	6.6E+04		1.3E+04
				1.0E+00	I						1.36E+09		1	0.1	~Butylphthalyl Butylglycolate	85-70-1					7.8E+04	3.3E+05		6.3E+04
				1.0E-01	I						1.36E+09		1	0.1	~Dibutyl Phthalate	84-74-2					7.8E+03	3.3E+04		6.3E+03
				8.0E-01	I						1.36E+09		1	0.1	~Diethyl Phthalate	84-66-2					6.3E+04	2.6E+05		5.1E+04
				1.0E-01	I			V			1.36E+09	2.13E+04	1		~Dimethylterephthalate	120-61-6					7.8E+03			7.8E+03
				1.0E-02	P						1.36E+09		1	0.1	~Octyl Phthalate, di-N-	117-84-0					7.8E+02	3.3E+03		6.3E+02
				5.0E-01	X						1.36E+09		1	0.1	~Phthalic Acid, p-	100-21-0					3.9E+04	1.6E+05		3.2E+04
				2.0E+00	I	2.0E-02	C				1.36E+09		1	0.1	~Phthalic Anhydride	85-44-9					1.6E+05	6.6E+05	2.8E+07	1.3E+05
				7.0E-02	I						1.36E+09		1	0.1	Picloram	1918-02-1					5.5E+03	2.3E+04		4.4E+03
				1.0E-04	X						1.36E+09		1	0.1	Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3					7.8E+00	3.3E+01		6.3E+00
				2.0E-03	X						1.36E+09		1	0.1	Picric Acid (2,4,6-Trinitrophenol)	88-89-1					1.6E+02	6.6E+02		1.3E+02
				7.3E-04	O						1.36E+09		1	0.1	Pirimiphos, Methyl	29232-93-7					5.7E+01	2.4E+02		4.6E+01
3.0E+01	C	8.6E-03	C	7.0E-06	H						1.36E+09		1	0.1	Polybrominated Biphenyls	36355-01-8	2.3E+00	8.2E+00	4.4E+04	1.8E+00	5.5E-01	2.3E+00		4.4E-01
															Polychlorinated Biphenyls (PCBs)									
7.0E-02	G	2.0E-05	G	7.0E-05	I			V			1.36E+09	5.86E+05	1	0.14	~Aroclor 1016	12674-11-2	9.9E+02	2.5E+03	8.2E+03	6.6E+02	5.5E+00	1.6E+01		4.1E+00
2.0E+00	G	5.7E-04	G					V			1.36E+09	2.04E+05	1	0.14	~Aroclor 1221	11104-28-2	3.5E+01	8.8E+01	1.0E+02	2.0E+01				
2.0E+00	G	5.7E-04	G					V			1.36E+09	1.12E+05	1	0.14	~Aroclor 1232	11141-16-5	3.5E+01	8.8E+01	5.5E+01	1.7E+01				

RMLs TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
2.0E+00	G	5.7E-04	G					V			1.36E+09	5.91E+05	1	0.14	~Aroclor 1242	53469-21-9	3.5E+01	8.8E+01	2.9E+02	2.3E+01				
2.0E+00	G	5.7E-04	G					V			1.36E+09	5.14E+05	1	0.14	~Aroclor 1248	12672-29-6	3.5E+01	8.8E+01	2.5E+02	2.3E+01				
2.0E+00	G	5.7E-04	G	2.0E-05	I			V			1.36E+09	8.43E+05	1	0.14	~Aroclor 1254	11097-69-1	3.5E+01	8.8E+01	4.1E+02	2.4E+01	1.6E+00	4.7E+00		1.2E+00
2.0E+00	G	5.7E-04	G					V			1.36E+09	1.31E+06	1	0.14	~Aroclor 1260	11096-82-5	3.5E+01	8.8E+01	6.5E+02	2.4E+01				
				6.0E-04	X			V			1.36E+09	1.15E+06	1	0.14	~Aroclor 5460	11126-42-4					4.7E+01	1.4E+02		3.5E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	2.43E+06	1	0.14	~Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	1.8E+01	4.5E+01	6.0E+02	1.3E+01	1.8E+00	5.5E+00	3.4E+03	1.4E+00
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.58E+06	1	0.14	~Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167)	52663-72-6	1.8E+01	4.5E+01	3.9E+02	1.2E+01	1.8E+00	5.5E+00	2.2E+03	1.4E+00
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.04E+06	1	0.14	~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	1.8E+01	4.5E+01	2.6E+02	1.2E+01	1.8E+00	5.5E+00	1.4E+03	1.4E+00
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.11E+06	1	0.14	~Hexachlorobiphenyl, 2,3,3',4,4',5- (PCB 156)	38380-08-4	1.8E+01	4.5E+01	2.7E+02	1.2E+01	1.8E+00	5.5E+00	1.5E+03	1.4E+00
3.9E+03	W	1.1E+00	W	2.3E-08	W	1.3E-06	W	V			1.36E+09	1.58E+06	1	0.14	~Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	1.8E-02	4.5E-02	3.9E-01	1.2E-02	1.8E-03	5.5E-03	2.2E+00	1.4E-03
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	7.33E+05	1	0.14	~Pentachlorobiphenyl, 2',3,4,4',5- (PCB 123)	65510-44-3	1.8E+01	4.5E+01	1.8E+02	1.2E+01	1.8E+00	5.5E+00	1.0E+03	1.4E+00
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	5.90E+05	1	0.14	~Pentachlorobiphenyl, 2,3',4,4',5- (PCB 118)	31508-00-6	1.8E+01	4.5E+01	1.5E+02	1.2E+01	1.8E+00	5.5E+00	8.2E+02	1.4E+00
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	6.01E+05	1	0.14	~Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	1.8E+01	4.5E+01	1.5E+02	1.2E+01	1.8E+00	5.5E+00	8.4E+02	1.4E+00
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.05E+06	1	0.14	~Pentachlorobiphenyl, 2,3,4,4',5- (PCB 114)	74472-37-0	1.8E+01	4.5E+01	2.6E+02	1.2E+01	1.8E+00	5.5E+00	1.5E+03	1.4E+00
1.3E+04	W	3.8E+00	W	7.0E-09	W	4.0E-07	W	V			1.36E+09	7.26E+05	1	0.14	~Pentachlorobiphenyl, 3,3',4,4',5- (PCB 126)	57465-28-8	5.3E-03	1.4E-02	5.4E-02	3.6E-03	5.5E-04	1.6E-03	3.0E-01	4.1E-04
2.0E+00	I	5.7E-04	I					V			1.36E+09	5.32E+05	1	0.14	~Polychlorinated Biphenyls (high risk)	1336-36-3	3.5E+01	8.8E+01	2.6E+02	2.3E+01				
4.0E-01	I	1.0E-04	I					V					1	0.14	~Polychlorinated Biphenyls (low risk)	1336-36-3								
7.0E-02	I	2.0E-05	I					V					1	0.14	~Polychlorinated Biphenyls (lowest risk)	1336-36-3								
1.3E+01	W	3.8E-03	W	7.0E-06	W	4.0E-04	W				1.36E+09		1	0.14	~Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	5.3E+00	1.4E+01	1.0E+05	3.8E+00	5.5E-01	1.6E+00	5.7E+05	4.1E-01
3.9E+01	W	1.1E-02	W	2.3E-06	W	1.3E-04	W	V			1.36E+09	5.09E+05	1	0.14	~Tetrachlorobiphenyl, 3,4,4',5- (PCB 81)	70362-50-4	1.8E+00	4.5E+00	1.3E+01	1.2E+00	1.8E-01	5.5E-01	7.1E+01	1.4E-01
						6.0E-04	I				1.36E+09		1	0.1	Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9							8.5E+05	8.5E+05
															Polynuclear Aromatic Hydrocarbons (PAHs)									
				6.0E-02	I			V			1.36E+09	1.41E+05	1	0.13	~Acenaphthene	83-32-9					4.7E+03	1.5E+04		3.6E+03
				3.0E-01	I			V			1.36E+09	5.23E+05	1	0.13	~Anthracene	120-12-7					2.3E+04	7.6E+04		1.8E+04
1.0E-01	E	6.0E-05	E					V	M		1.36E+09	4.41E+06	1	0.13	~Benz[a]anthracene	56-55-3	1.5E+02	4.6E+02	7.4E+03	1.1E+02				
				9.0E-05	X	2.0E-06	X				1.36E+09		1	0.1	~Benzo(e)pyrene	192-97-2					7.0E+00	3.0E+01	2.8E+03	5.7E+00
1.2E+00	C	1.1E-04	C								1.36E+09		1	0.13	~Benzo(j)fluoranthene	205-82-3	5.8E+01	1.6E+02	3.5E+06	4.2E+01				
1.0E+00	I	6.0E-04	I	3.0E-04	I	2.0E-06	I		M		1.36E+09		1	0.13	~Benzo[a]pyrene	50-32-8	1.5E+01	4.6E+01	2.3E+05	1.1E+01	2.3E+01	7.6E+01	2.8E+03	1.8E+01
1.0E-01	E	6.0E-05	E						M		1.36E+09		1	0.13	~Benzo[b]fluoranthene	205-99-2	1.5E+02	4.6E+02	2.3E+06	1.1E+02				
1.0E-02	E	6.0E-06	E						M		1.36E+09		1	0.13	~Benzo[k]fluoranthene	207-08-9	1.5E+03	4.6E+03	2.3E+07	1.1E+03				
				8.0E-02	I			V			1.36E+09	7.99E+04	1	0.13	~Chloronaphthalene, Beta-	91-58-7					6.3E+03	2.0E+04		4.8E+03
1.0E-03	E	6.0E-07	E						M		1.36E+09		1	0.13	~Chrysene	218-01-9	1.5E+04	4.6E+04	2.3E+08	1.1E+04				
1.0E+00	E	6.0E-04	E						M		1.36E+09		1	0.13	~Dibenz[a,h]anthracene	53-70-3	1.5E+01	4.6E+01	2.3E+05	1.1E+01				
1.2E+01	C	1.1E-03	C								1.36E+09		1	0.13	~Dibenzo(a,e)pyrene	192-65-4	5.8E+00	1.6E+01	3.5E+05	4.2E+00				
2.5E+02	C	7.1E-02	C						M		1.36E+09		1	0.13	~Dimethylbenz(a)anthracene, 7,12-	57-97-6	6.1E-02	1.8E-01	1.9E+03	4.6E-02				
				4.0E-02	I						1.36E+09		1	0.13	~Fluoranthene	206-44-0					3.1E+03	1.0E+04		2.4E+03
				4.0E-02	I			V			1.36E+09	2.81E+05	1	0.13	~Fluorene	86-73-7					3.1E+03	1.0E+04		2.4E+03
1.0E-01	E	6.0E-05	E						M		1.36E+09		1	0.13	~Indeno[1,2,3-cd]pyrene	193-39-5	1.5E+02	4.6E+02	2.3E+06	1.1E+02				
2.9E-02	P			7.0E-02	A			V		3.94E+02	1.36E+09	5.86E+04	1	0.13	~Methylnaphthalene, 1-	90-12-0	2.4E+03	6.6E+03		1.8E+03	5.5E+03	1.8E+04		4.2E+03
				4.0E-03	I			V			1.36E+09	5.80E+04	1	0.13	~Methylnaphthalene, 2-	91-57-6					3.1E+02	1.0E+03		2.4E+02
1.2E-01	C	3.4E-05	C	2.0E-02	I	3.0E-03	I	V			1.36E+09	4.63E+04	1	0.13	~Naphthalene	91-20-3	5.8E+02	1.6E+03	3.8E+02	2.0E+02	1.6E+03	5.1E+03	1.4E+02	1.3E+02
1.2E+00	C	1.1E-04	C								1.36E+09		1	0.13	~Nitropyrene, 4-	57835-92-4	5.8E+01	1.6E+02	3.5E+06	4.2E+01				
				3.0E-02	I			V			1.36E+09	2.38E+06	1	0.13	~Pyrene	129-00-0					2.3E+03	7.6E+03		1.8E+03
1.5E-01	I			9.0E-03	I						1.36E+09		1	0.1	Prochloraz	67747-09-5	4.6E+02	1.6E+03		3.6E+02	7.0E+02	3.0E+03		5.7E+02
				6.0E-03	H			V			1.36E+09	4.20E+05	1		Profluralin	26399-36-0					4.7E+02			4.7E+02
				1.5E-02	I						1.36E+09		1	0.1	Prometon	1610-18-0					1.2E+03	4.9E+03		9.5E+02
				4.0E-02	O						1.36E+09		1	0.1	Prometryn	7287-19-6					3.1E+03	1.3E+04		2.5E+03
				7.5E-02	I						1.36E+09		1	0.1	Pronamide	23950-58-5					5.9E+03	2.5E+04		4.7E+03
				1.3E-02	I						1.36E+09		1	0.1	Propachlor	1918-16-7					1.0E+03	4.3E+03		8.2E+02
				5.0E-03	I						1.36E+09		1	0.1	Propanil	709-98-8					3.9E+02	1.6E+03		3.2E+02
1.9E-01	O			4.0E-02	O						1.36E+09		1	0.1	Propargite	2312-35-8	3.6E+02	1.3E+03		2.8E+02	3.1E+03	1.3E+04		2.5E+03
				2.0E-03	I			V		1.11E+05	1.36E+09	6.27E+04	1		Propargyl Alcohol	107-19-7					1.6E+02			1.6E+02
				2.0E-02	I						1.36E+09		1	0.1	Propazine	139-40-2					1.6E+03	6.6E+03		1.3E+03
				2.0E-02	I						1.36E+09		1	0.1	Propham	122-42-9					1.6E+03	6.6E+03		1.3E+03
				1.0E-01	O						1.36E+09		1	0.1	Propiconazole	60207-90-1					7.8E+03	3.3E+04		6.3E+03
						8.0E-03	I	V		3.26E+04	1.36E+09	8.94E+03	1		Propionaldehyde	123-38-6							7.5E+01	7.5E+01
				1.0E-01	X	1.0E+00	X	V		2.64E+02	1.36E+09	6.99E+03	1		Propyl benzene	103-65-1					7.8E+03		7.3E+03	3.8E+03
						3.0E+00	C	V		3.49E+02	1.36E+09	7.04E+02	1		Propylene	115-07-1							2.2E+03	2.2E+03
				2.0E+01	P						1.36E+09		1	0.1	Propylene Glycol	57-55-6					1.6E+06	6.6E+06		1.3E+06
						2.7E-04	A				1.36E+09		1	0.1	Propylene Glycol Dinitrate	6423-43-4							3.9E+05	3.9E+05
				7.0E-01	H	2.0E+00	I	V		1.06E+05	1.36E+09	7.83E+04	1		Propylene Glycol Monomethyl Ether	107-98-2					5.5E+04		1.6E+05	4.1E+04
2.4E-01	I	3.7E-06	I			3.0E-02	I	V		7.77E+04	1.36E+09	1.03E+04	1		Propylene Oxide	75-56-9	2.9E+02		7.8E+02	2.1E+02			3.2E+02	3.2E+02
				1.0E-03	I			V		5.30E+05	1.36E+09	5.54E+04	1		Pyridine	110-86-1					7.8E+01			7.8E+01
				5.0E-04	I						1.36E+09		1	0.1	Quinalphos	13593-03-8					3.9E+01	1.6E+02		3.2E+01
3.0E+00	I										1.36E+09		1	0.1	Quinoline	91-22-5	2.3E+01	8.2E+01		1.8E+01				
				9.0E-03	I						1.36E+09		1	0.1	Quizalofop-ethyl	76578-14-8					7.0E+02	3.0E+03		5.7E+02
						3.0E+04	A						1		Refractory Ceramic Fibers (units in fibers)	E715557								
				3.0E-02	I						1.36E+09		1	0.1	Resmethrin	10453-86-8					2.3E+03	9.9E+03		1.9E+03
				5.0E-02	H			V			1.36E+09	4.65E+05	1		Ronnel	299-84-3					3.9E+03			3.9E+03
				4.0E-03	I						1.36E+09		1	0.1	Rotenone	83-79-4					3.1E+02	1.3E+03		2.5E+02
2.2E-01	C	6.3E-05	C						M		1.36E+09		1	0.1	Safrole	94-59-7	7.0E+01	2.7E+02	2.2E+06	5.5E+01				
				5.0E-03	I						1.36E+09		1		Selenious Acid	7783-00-8					3.9E+02			3.9E+02
				5.0E-03	I	2.0E-02	C				1.36E+09		1		Selenium	7782-49-2					3.9E+02		2.8E+07	3.9E+02
				5.0E-03	C	2.0E-02	C				1.36E+09		1		Selenium Sulfide	7446-34-6					3.9E+02		2.8E+07	3.9E+02
				1.4E-01	O						1.36E+09		1	0.1	Sethoxydim	74051-80-2					1.1E+04	4.6E+04		8.8E+03
						3.0E-03	C				1.36E+09		1		Silica (crystalline, respirable)	7631-86-9							4.3E+06	4.3E+06
				5.0E-03	I						1.36E+09		0.04		Silver	7440-22-4					3.9E+02			3.9E+02

RMLs TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
1.2E-01	H			5.0E-03	I						1.36E+09		1	0.1	Simazine	122-34-9	5.8E+02	2.1E+03		4.5E+02	3.9E+02	1.6E+03		3.2E+02
				1.3E-02	I						1.36E+09		1	0.1	Sodium Acifluorfen	62476-59-9					1.0E+03	4.3E+03		8.2E+02
				4.0E-03	I						1.36E+09		1		Sodium Azide	26628-22-8					3.1E+02			3.1E+02
2.7E-01	H			3.0E-02	I						1.36E+09		1	0.1	Sodium Diethyldithiocarbamate	148-18-5	2.6E+02	9.2E+02		2.0E+02	2.3E+03	9.9E+03		1.9E+03
				5.0E-02	A	1.3E-02	C				1.36E+09		1		Sodium Fluoride	7681-49-4					3.9E+03		1.8E+07	3.9E+03
				2.0E-05	I						1.36E+09		1	0.1	Sodium Fluoroacetate	62-74-8					1.6E+00	6.6E+00		1.3E+00
				1.0E-03	H						1.36E+09		1		Sodium Metavanadate	13718-26-8					7.8E+01			7.8E+01
				8.0E-04	P						1.36E+09		1		Sodium Tungstate	13472-45-2					6.3E+01			6.3E+01
				8.0E-04	P						1.36E+09		1		Sodium Tungstate Dihydrate	10213-10-2					6.3E+01			6.3E+01
2.4E-02	H			3.0E-02	I						1.36E+09		1	0.1	Stirofos (Tetrachlorovinphos)	961-11-5	2.9E+03	1.0E+04		2.3E+03	2.3E+03	9.9E+03		1.9E+03
				6.0E-01	I						1.36E+09		1		Strontium, Stable	7440-24-6					4.7E+04			4.7E+04
				3.0E-04	I						1.36E+09		1	0.1	Strychnine	57-24-9					2.3E+01	9.9E+01		1.9E+01
				2.0E-01	I	1.0E+00	I	V		8.67E+02	1.36E+09	9.35E+03	1		Styrene	100-42-5					1.6E+04		9.7E+03	6.0E+03
				3.0E-03	P						1.36E+09		1	0.1	Styrene-Acrylonitrile (SAN) Trimer (THNA isomer)	57964-39-3					2.3E+02	9.9E+02		1.9E+02
				3.0E-03	P						1.36E+09		1	0.1	Styrene-Acrylonitrile (SAN) Trimer (THNP isomer)	57964-40-6					2.3E+02	9.9E+02		1.9E+02
				1.0E-03	P	2.0E-03	X				1.36E+09		1	0.1	Sulfolane	126-33-0					7.8E+01	3.3E+02	2.8E+06	6.3E+01
				8.0E-04	P						1.36E+09		1	0.1	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9					6.3E+01	2.6E+02		5.1E+01
						1.0E-03	C	V			1.36E+09		1		Sulfur Trioxide	7446-11-9							1.4E+06	1.4E+06
						1.0E-03	C				1.36E+09		1		Sulfuric Acid	7664-93-9							1.4E+06	1.4E+06
2.5E-02	I	7.1E-06	I	5.0E-02	H						1.36E+09		1	0.1	Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl	140-57-8	2.8E+03	9.9E+03	5.4E+07	2.2E+03	3.9E+03	1.6E+04		3.2E+03
				3.0E-02	H						1.36E+09		1	0.1	TCMTB	21564-17-0					2.3E+03	9.9E+03		1.9E+03
				7.0E-02	I						1.36E+09		1	0.1	Tebuthiuron	34014-18-1					5.5E+03	2.3E+04		4.4E+03
				2.0E-02	H						1.36E+09		1	0.1	Temephos	3383-96-8					1.6E+03	6.6E+03		1.3E+03
				1.3E-02	I						1.36E+09		1	0.1	Terbacil	5902-51-2					1.0E+03	4.3E+03		8.2E+02
				2.5E-05	H			V		3.09E+01	1.36E+09	2.64E+05	1		Terbufos	13071-79-9					2.0E+00			2.0E+00
				1.0E-03	I						1.36E+09		1	0.1	Terbutryn	886-50-0					7.8E+01	3.3E+02		6.3E+01
5.0E-03	C	1.3E-06	C					V			1.36E+09	3.99E+03	1		Tert-Butyl Acetate	540-88-5	1.4E+04		8.6E+02	8.1E+02				
				1.0E-04	I						1.36E+09		1	0.1	Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1					7.8E+00	3.3E+01		6.3E+00
				3.0E-05	P			V			1.36E+09	5.07E+04	1		Tetrachlorobenzene, 1,2,4,5-	95-94-3					2.3E+00			2.3E+00
2.6E-02	I	7.4E-06	I	3.0E-02	I			V		6.80E+02	1.36E+09	5.68E+03	1		Tetrachloroethane, 1,1,1,2-	630-20-6	2.7E+03		2.2E+02	2.0E+02	2.3E+03			2.3E+03
2.0E-01	I	5.8E-05	C	2.0E-02	I			V		1.90E+03	1.36E+09	1.51E+04	1		Tetrachloroethane, 1,1,2,2-	79-34-5	3.5E+02		7.3E+01	6.0E+01	1.6E+03			1.6E+03
2.1E-03	I	2.6E-07	I	6.0E-03	I	4.0E-02	I	V		1.66E+02	1.36E+09	2.35E+03	1		Tetrachloroethylene	127-18-4	3.3E+04		2.5E+03	2.4E+03	4.7E+02		9.8E+01	8.1E+01
				3.0E-02	I						1.36E+09		1	0.1	Tetrachlorophenol, 2,3,4,6-	58-90-2					2.3E+03	9.9E+03		1.9E+03
1.6E+01	X			6.0E-05	X			V			1.36E+09	1.05E+05	1		Tetrachlorotoluene, p- alpha, alpha, alpha-	5216-25-1	4.3E+00			4.3E+00	4.7E+00			4.7E+00
				5.0E-04	I						1.36E+09		1	0.1	Tetraethyl Dithiopyrophosphate	3689-24-5					3.9E+01	1.6E+02		3.2E+01
						8.0E+01	I	V		2.05E+03	1.36E+09	1.22E+03	1		Tetrafluoroethane, 1,1,1,2-	811-97-2							1.0E+05	1.0E+05
				1.0E-04	X						1.36E+09		1	0.1	Tetramethylphosphoramide, -N,N,N',N" (TMPA)	16853-36-4					7.8E+00	3.3E+01		6.3E+00
				2.0E-03	P						1.36E+09		1	0.00065	Tetryl (Trinitrophenylmethylnitramine)	479-45-8					1.6E+02	1.0E+05		1.6E+02
				2.0E-05	G						1.36E+09		1		Thallic Oxide	1314-32-5					1.6E+00			1.6E+00
				1.0E-05	X						1.36E+09		1		Thallium (I) Nitrate	10102-45-1					7.8E-01			7.8E-01
				1.0E-05	X						1.36E+09		1		Thallium (Soluble Salts)	7440-28-0					7.8E-01			7.8E-01
				1.0E-05	X			V			1.36E+09	1.40E+05	1		Thallium Acetate	563-68-8					7.8E-01			7.8E-01
				2.0E-05	X						1.36E+09		1	0.1	Thallium Carbonate	6533-73-9					1.6E+00	6.6E+00		1.3E+00
				1.0E-05	X						1.36E+09		1		Thallium Chloride	7791-12-0					7.8E-01			7.8E-01
				1.0E-05	G						1.36E+09		1		Thallium Selenite	12039-52-0					7.8E-01			7.8E-01
				2.0E-05	X						1.36E+09		1		Thallium Sulfate	7446-18-6					1.6E+00			1.6E+00
				4.3E-02	O						1.36E+09		1	0.1	Thifensulfuron-methyl	79277-27-3					3.4E+03	1.4E+04		2.7E+03
				1.0E-02	I						1.36E+09		1	0.1	Thiobencarb	28249-77-6					7.8E+02	3.3E+03		6.3E+02
				7.0E-02	X						1.36E+09		1	0.0075	Thiodiglycol	111-48-8					5.5E+03	3.1E+05		5.4E+03
				3.0E-04	H						1.36E+09		1	0.1	Thiofanox	39196-18-4					2.3E+01	9.9E+01		1.9E+01
1.2E-02	O			1.6E-01	O						1.36E+09		1	0.1	Thiophanate, Methyl	23564-05-8	6.0E+03	2.1E+04		4.7E+03	1.3E+04	5.3E+04		1.0E+04
				1.5E-02	O						1.36E+09		1	0.1	Thiram	137-26-8					1.2E+03	4.9E+03		9.5E+02
				6.0E-01	H						1.36E+09		1		Tin	7440-31-5					4.7E+04			4.7E+04
						1.0E-04	A	V			1.36E+09		1		Titanium Tetrachloride	7550-45-0							1.4E+05	1.4E+05
				8.0E-02	I	5.0E+00	I	V		8.18E+02	1.36E+09	4.29E+03	1		Toluene	108-88-3					6.3E+03		2.2E+04	4.9E+03
3.9E-02	C	1.1E-05	C			8.0E-06	C	V			1.36E+09	7.62E+05	1		Toluene-2,4-diisocyanate	584-84-9	1.8E+03		1.9E+04	1.6E+03			6.4E+00	6.4E+00
1.8E-01	X			2.0E-04	X						1.36E+09		1	0.1	Toluene-2,5-diamine	95-70-5	3.9E+02	1.4E+03		3.0E+02	1.6E+01	6.6E+01		1.3E+01
3.9E-02	C	1.1E-05	C			8.0E-06	C	V		1.71E+03	1.36E+09	6.32E+05	1		Toluene-2,6-diisocyanate	91-08-7	1.8E+03		1.6E+04	1.6E+03			5.3E+00	5.3E+00
				1.0E-04	X						1.36E+09		1	0.1	Toluenediamine, 2,3-	2687-25-4					7.8E+00	3.3E+01		6.3E+00
				1.0E-04	X						1.36E+09		1	0.1	Toluenediamine, 3,4-	496-72-0					7.8E+00	3.3E+01		6.3E+00
				5.0E-03	P						1.36E+09		1	0.1	Toluic Acid, p-	99-94-5					3.9E+02	1.6E+03		3.2E+02
1.6E-02	P	5.1E-05	C								1.36E+09		1	0.1	Toluidine, o- (Methylaniline, 2-)	95-53-4	4.3E+03	1.5E+04	7.5E+06	3.4E+03				
3.0E-02	P			4.0E-03	X						1.36E+09		1	0.1	Toluidine, p-	106-49-0	2.3E+03	8.2E+03		1.8E+03	3.1E+02	1.3E+03		2.5E+02
				3.0E+00	P			V		3.42E-01	1.36E+09	1.38E+03	1		Total Petroleum Hydrocarbons (Aliphatic High)	E1790670					2.3E+05			2.3E+05
						6.0E-01	P	V		1.41E+02	1.36E+09	8.29E+02	1		Total Petroleum Hydrocarbons (Aliphatic Low)	E1790666							5.2E+02	5.2E+02
				1.0E-02	X	1.0E-01	P	V		6.86E+00	1.36E+09	1.04E+03	1		Total Petroleum Hydrocarbons (Aliphatic Medium)	E1790668					7.8E+02		1.1E+02	9.6E+01
				4.0E-02	P						1.36E+09		1	0.13	Total Petroleum Hydrocarbons (Aromatic High)	E1790676					3.1E+03	1.0E+04		2.4E+03
				4.0E-03	P	3.0E-02	P	V		1.82E+03	1.36E+09	3.54E+03	1		Total Petroleum Hydrocarbons (Aromatic Low)	E1790672					3.1E+02		1.1E+02	8.2E+01
				4.0E-03	P	3.0E-03	P	V			1.36E+09	5.24E+04	1	0.13	Total Petroleum Hydrocarbons (Aromatic Medium)	E1790674					3.1E+02	1.0E+03	1.6E+02	9.7E+01
1.1E+00	I	3.2E-04	I	9.0E-05	P						1.36E+09		1	0.1	Toxaphene	8001-35-2	6.3E+01	2.2E+02	1.2E+06	4.9E+01	7.0E+00	3.0E+01		5.7E+00
				3.0E-05	X						1.36E+09		1	0.1	Toxaphene, Weathered	E1841606					2.3E+00	9.9E+00		1.9E+00
				7.5E-03	I						1.36E+09		1	0.1	Tralomethrin	66841-25-6					5.9E+02	2.5E+03		4.7E+02
				3.0E-04	A			V			1.36E+09	3.36E+03	1		Tri-n-butyltin	688-73-3					2.3E+01			2.3E+01
				8.0E+01	X						1.36E+09		1	0.1	Triacetin	102-76-1					6.3E+06	2.6E+07		5.1E+06
				3.4E-02	O						1.36E+09		1	0.1	Triadimefon	43121-43-3					2.7E+03	1.1E+04		2.1E+03
7.2E-02	O			2.5E-02	O			V			1.36E+09	3.62E+05	1		Triallate	2303-17-5	9.7E+02			9.7E+02	2.0E+03			2.0E+03
				1.0E-02	I						1.36E+09		1	0.1	Triasulfuron	82097-50-5					7.8E+02	3.3E+03		6.3E+02

RMLs TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
				8.0E-03	I						1.36E+09		1	0.1	Tribenuron-methyl	101200-48-0					6.3E+02	2.6E+03		5.1E+02
				5.0E-03	I			V			1.36E+09	4.83E+04	1		Tribromobenzene, 1,2,4-	615-54-3					3.9E+02			3.9E+02
				9.0E-03	X						1.36E+09		1	0.1	Tribromophenol, 2,4,6-	118-79-6					7.0E+02	3.0E+03		5.7E+02
				2.0E-04	O						1.36E+09		1	0.1	Tribufos	78-48-8					1.6E+01	6.6E+01		1.3E+01
9.0E-03	P			1.0E-02	P						1.36E+09		1	0.1	Tributyl Phosphate	126-73-8	7.7E+03	2.7E+04		6.0E+03	7.8E+02	3.3E+03		6.3E+02
				3.0E-04	P						1.36E+09		1	0.1	Tributyltin Compounds	E1790679					2.3E+01	9.9E+01		1.9E+01
				3.0E-04	I						1.36E+09		1	0.1	Tributyltin Oxide	56-35-9					2.3E+01	9.9E+01		1.9E+01
											1.36E+09		1	0.1	Trichloramine	10025-85-1								
				3.0E+01	I	5.0E+00	P	V		9.10E+02	1.36E+09	1.29E+03	1		Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1					2.3E+06		6.7E+03	6.7E+03
7.0E-02	I			2.0E-02	I						1.36E+09		1	0.1	Trichloroacetic Acid	76-03-9	9.9E+02	3.5E+03		7.8E+02	1.6E+03	6.6E+03		1.3E+03
2.9E-02	H										1.36E+09		1	0.1	Trichloroaniline HCl, 2,4,6-	33663-50-2	2.4E+03	8.5E+03		1.9E+03				
7.0E-03	X			3.0E-05	X						1.36E+09		1	0.1	Trichloroaniline, 2,4,6-	634-93-5	9.9E+03	3.5E+04		7.8E+03	2.3E+00	9.9E+00		1.9E+00
				8.0E-04	X			V			1.36E+09	3.22E+04	1		Trichlorobenzene, 1,2,3-	87-61-6					6.3E+01			6.3E+01
2.9E-02	P			1.0E-02	I	2.0E-03	P	V		4.04E+02	1.36E+09	2.99E+04	1		Trichlorobenzene, 1,2,4-	120-82-1	2.4E+03			2.4E+03	7.8E+02		6.2E+01	5.8E+01
				2.0E+00	I	5.0E+00	I	V		6.40E+02	1.36E+09	1.65E+03	1		Trichloroethane, 1,1,1-	71-55-6					1.6E+05		8.6E+03	8.1E+03
5.7E-02	I	1.6E-05	I	4.0E-03	I	2.0E-04	X	V		2.16E+03	1.36E+09	7.22E+03	1		Trichloroethane, 1,1,2-	79-00-5	1.2E+03		1.3E+02	1.1E+02	3.1E+02		1.5E+00	1.5E+00
4.6E-02	I	4.1E-06	I	5.0E-04	I	2.0E-03	I	V	M	6.92E+02	1.36E+09	2.21E+03	1		Trichloroethylene	79-01-6	8.8E+02		1.1E+02	9.4E+01	3.9E+01		4.6E+00	4.1E+00
				3.0E-01	I			V		1.23E+03	1.36E+09	1.04E+03	1		Trichlorofluoromethane	75-69-4					2.3E+04			2.3E+04
				1.0E-01	I						1.36E+09		1	0.1	Trichlorophenol, 2,4,5-	95-95-4					7.8E+03	3.3E+04		6.3E+03
1.1E-02	I	3.1E-06	I	1.0E-03	P						1.36E+09		1	0.1	Trichlorophenol, 2,4,6-	88-06-2	6.3E+03	2.2E+04	1.2E+08	4.9E+03	7.8E+01	3.3E+02		6.3E+01
				1.0E-02	I						1.36E+09		1	0.1	Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5					7.8E+02	3.3E+03		6.3E+02
				8.0E-03	I						1.36E+09		1	0.1	Trichlorophenoxypropionic acid, -2,4,5	93-72-1					6.3E+02	2.6E+03		5.1E+02
				5.0E-03	I			V		1.28E+03	1.36E+09	1.50E+04	1		Trichloropropane, 1,1,2-	598-77-6					3.9E+02			3.9E+02
3.0E+01	I			4.0E-03	I	3.0E-04	I	V	M	1.40E+03	1.36E+09	1.57E+04	1		Trichloropropane, 1,2,3-	96-18-4	5.1E-01			5.1E-01	3.1E+02		4.9E+00	4.8E+00
				3.0E-03	X	3.0E-04	P	V		3.11E+02	1.36E+09	2.34E+03	1		Trichloropropene, 1,2,3-	96-19-5					2.3E+02		7.3E-01	7.3E-01
				2.0E-02	A						1.36E+09		1	0.1	Tricresyl Phosphate (TCP)	1330-78-5					1.6E+03	6.6E+03		1.3E+03
				3.0E-03	I						1.36E+09		1	0.1	Tridiphane	58138-08-2					2.3E+02	9.9E+02		1.9E+02
						7.0E-03	I	V		2.79E+04	1.36E+09	1.58E+04	1		Triethylamine	121-44-8							1.2E+02	1.2E+02
				2.0E+00	P						1.36E+09		1	0.1	Triethylene Glycol	112-27-6					1.6E+05	6.6E+05		1.3E+05
						2.0E+01	P	V		4.81E+03	1.36E+09	7.12E+02	1		Trifluoroethane, 1,1,1-	420-46-2							1.5E+04	1.5E+04
7.7E-03	I			7.5E-03	I			V			1.36E+09	5.13E+05	1		Trifluralin	1582-09-8	9.0E+03			9.0E+03	5.9E+02			5.9E+02
2.0E-02	P			1.0E-02	P						1.36E+09		1	0.1	Trimethyl Phosphate	512-56-1	3.5E+03	1.2E+04		2.7E+03	7.8E+02	3.3E+03		6.3E+02
				1.0E-02	I	6.0E-02	I	V		2.93E+02	1.36E+09	9.44E+03	1		Trimethylbenzene, 1,2,3-	526-73-8					7.8E+02		5.9E+02	3.4E+02
				1.0E-02	I	6.0E-02	I	V		2.19E+02	1.36E+09	7.91E+03	1		Trimethylbenzene, 1,2,4-	95-63-6					7.8E+02		5.0E+02	3.0E+02
				1.0E-02	I	6.0E-02	I	V		1.82E+02	1.36E+09	6.61E+03	1		Trimethylbenzene, 1,3,5-	108-67-8					7.8E+02		4.1E+02	2.7E+02
				1.0E-02	X			V		2.96E+01	1.36E+09	1.00E+03	1		Trimethylpentene, 2,4,4-	25167-70-8					7.8E+02			7.8E+02
				3.0E-02	I						1.36E+09		1	0.019	Trinitrobenzene, 1,3,5-	99-35-4					2.3E+03	5.2E+04		2.2E+03
3.0E-02	I			5.0E-04	I						1.36E+09		1	0.032	Trinitrotoluene, 2,4,6-	118-96-7	2.3E+03	2.6E+04		2.1E+03	3.9E+01	5.2E+02		3.6E+01
				2.0E-02	P						1.36E+09		1	0.1	Triphenylphosphine Oxide	791-28-6					1.6E+03	6.6E+03		1.3E+03
				2.0E-02	A						1.36E+09		1	0.1	Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8					1.6E+03	6.6E+03		1.3E+03
				1.0E-02	X						1.36E+09		1	0.1	Tris(1-chloro-2-propyl)phosphate	13674-84-5					7.8E+02	3.3E+03		6.3E+02
2.3E+00	C	6.6E-04	C					V		4.67E+02	1.36E+09	9.03E+05	1		Tris(2,3-dibromopropyl)phosphate	126-72-7	3.0E+01		3.8E+02	2.8E+01				
2.0E-02	P			7.0E-03	P						1.36E+09		1	0.1	Tris(2-chloroethyl)phosphate	115-96-8	3.5E+03	1.2E+04		2.7E+03	5.5E+02	2.3E+03		4.4E+02
3.2E-03	P			1.0E-01	P						1.36E+09		1	0.1	Tris(2-ethylhexyl)phosphate	78-42-2	2.2E+04	7.7E+04		1.7E+04	7.8E+03	3.3E+04		6.3E+03
				8.0E-04	P						1.36E+09		1		Tungsten	7440-33-7					6.3E+01			6.3E+01
				2.0E-04	A	4.0E-05	A				1.36E+09		1		Uranium	7440-61-1					1.6E+01		5.7E+04	1.6E+01
1.0E+00	C	2.9E-04	C						M		1.36E+09		1	0.1	Urethane	51-79-6	1.5E+01	6.0E+01	4.8E+05	1.2E+01				
		8.3E-03	P	9.0E-03	I	7.0E-06	P				1.36E+09		0.026		Vanadium Pentoxide	1314-62-1			4.6E+04	4.6E+04	7.0E+02		9.9E+03	6.6E+02
				5.0E-03	G	1.0E-04	A				1.36E+09		0.026		Vanadium and Compounds	7440-62-2					3.9E+02		1.4E+05	3.9E+02
				1.0E-03	I			V			1.36E+09	1.23E+05	1		Vernolate	1929-77-7					7.8E+01			7.8E+01
				1.2E-03	O						1.36E+09		1	0.1	Vinclozolin	50471-44-8					9.4E+01	4.0E+02		7.6E+01
				1.0E+00	H	2.0E-01	I	V		2.75E+03	1.36E+09	4.40E+03	1		Vinyl Acetate	108-05-4					7.8E+04		9.2E+02	9.1E+02
		1.5E-05	P			3.0E-03	I	V		2.47E+03	1.36E+09	1.37E+03	1		Vinyl Bromide	593-60-2			2.6E+01	2.6E+01			4.3E+00	4.3E+00
7.2E-01	I	4.4E-06	I	3.0E-03	I	8.0E-02	A	V	M	3.92E+03	1.36E+09	9.56E+02	1		Vinyl Chloride	75-01-4	9.4E+00		1.6E+01	5.9E+00	2.3E+02		8.0E+01	6.0E+01
				3.0E-04	I						1.36E+09		1	0.1	Warfarin	81-81-2					2.3E+01	9.9E+01		1.9E+01
				2.0E-01	G	1.0E-01	G	V		3.88E+02	1.36E+09	5.47E+03	1		Xylene, m-	108-38-3					1.6E+04		5.7E+02	5.5E+02
				2.0E-01	G	1.0E-01	G	V		4.34E+02	1.36E+09	6.45E+03	1		Xylene, o-	95-47-6					1.6E+04		6.7E+02	6.4E+02
				2.0E-01	G	1.0E-01	G	V		3.90E+02	1.36E+09	5.58E+03	1		Xylene, p-	106-42-3					1.6E+04		5.8E+02	5.6E+02
				2.0E-01	I	1.0E-01	I	V		2.60E+02	1.36E+09	5.74E+03	1		Xylenes	1330-20-7					1.6E+04		6.0E+02	5.8E+02
				3.0E-04	I						1.36E+09		1		Zinc Phosphide	1314-84-7					2.3E+01			2.3E+01
				3.0E-01	I						1.36E+09		1		Zinc and Compounds	7440-66-6					2.3E+04			2.3E+04
				5.0E-02	I						1.36E+09		1	0.1	Zineb	12122-67-7					3.9E+03	1.6E+04		3.2E+03
				8.0E-05	X						1.36E+09		1		Zirconium	7440-67-7					6.3E+00			6.3E+00

RMLs TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1				
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	log K_{ow} (unitless)	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
				3.0E-04	O					-0.85	1	1.0	Yes	Acephate	30560-19-1					6.0E+00	8.6E+03		6.0E+00	
		2.2E-06	I			9.0E-03	I	V		-0.34	1	1.0	Yes	Acetaldehyde	75-07-0			2.6E+02	2.6E+02			1.9E+01	1.9E+01	
				2.0E-02	I					3.03	1	0.9	Yes	Acetochlor	34256-82-1					4.0E+02	2.9E+03		3.5E+02	
				9.0E-01	I			V		-0.24	1	1.0	Yes	Acetone	67-64-1					1.8E+04	4.4E+06		1.8E+04	
						2.0E-03	X			-0.03	1	1.0	Yes	Acetone Cyanohydrin	75-86-5									
						6.0E-02	I	V		-0.34	1	1.0	Yes	Acetonitrile	75-05-8							1.3E+02	1.3E+02	
				1.0E-01	I			V		1.58	1	1.0	Yes	Acetophenone	98-86-2					2.0E+03	4.6E+04		1.9E+03	
3.8E+00	C	1.3E-03	C							3.12	1	1.0	Yes	Acetylaminofluorene, 2-	53-96-3	2.1E+00	8.0E+00		1.6E+00					
				5.0E-04	I	2.0E-05	I	V		-0.01	1	1.0	Yes	Acrolein	107-02-8					1.0E+01	1.7E+03	4.2E-02	4.2E-02	
5.0E-01	I	1.0E-04	I	2.0E-03	I	6.0E-03	I		M	-0.67	1	1.0	Yes	Acrylamide	79-06-1	5.0E+00	2.3E+03		5.0E+00	4.0E+01	2.1E+04		4.0E+01	
				5.0E-01	I	2.0E-04	P	V		0.35	1	1.0	Yes	Acrylic Acid	79-10-7					1.0E+04	1.1E+06	4.2E-01	4.2E-01	
5.4E-01	I	6.8E-05	I	1.0E-02	A	2.0E-03	I	V		0.25	1	1.0	Yes	Acrylonitrile	107-13-1	1.4E+01	1.4E+03	8.3E+00	5.2E+00	2.0E+02	2.2E+04	4.2E+00	4.1E+00	
						6.0E-03	P			-0.32	1	1.0	Yes	Adiponitrile	111-69-3									
5.6E-02	C			1.0E-02	I					3.52	1	0.9	Yes	Alachlor	15972-60-8	1.4E+02	4.4E+02		1.1E+02	2.0E+02	6.9E+02		1.6E+02	2.0E+00
				1.0E-03	I					1.13	1	1.0	Yes	Aldicarb	116-06-3					2.0E+01	1.4E+03		2.0E+01	3.0E+00
				1.0E-03	I					-0.57	1	1.0	Yes	Aldicarb Sulfone	1646-88-4					2.0E+01	2.4E+04		2.0E+01	2.0E+00
										-0.78	1	1.0	Yes	Aldicarb sulfoxide	1646-87-3									4.0E+00
1.7E+01	I	4.9E-03	I	3.0E-05	I			V		6.5	1	1.0	No	Aldrin	309-00-2	4.6E-01		1.1E-01	9.2E-02	6.0E-01			6.0E-01	
				4.0E-03	P	1.0E-04	X	V		0.17	1	1.0	Yes	Allyl Alcohol	107-18-6					8.0E+01	1.0E+04	2.1E-01	2.1E-01	
2.1E-02	C	6.0E-06	C			1.0E-03	I	V		1.93	1	1.0	Yes	Allyl Chloride	107-05-1	3.7E+02	3.5E+03	9.4E+01	7.3E+01			2.1E+00	2.1E+00	
				1.0E+00	P	5.0E-03	P				1	1.0	Yes	Aluminum	7429-90-5					2.0E+04	4.6E+06		2.0E+04	
				4.0E-04	I						1	1.0	Yes	Aluminum Phosphide	20859-73-8					8.0E+00	1.8E+03		8.0E+00	
				9.0E-03	I					2.98	1	1.0	Yes	Ametryn	834-12-8					1.8E+02	9.8E+02		1.5E+02	
2.1E+01	C	6.0E-03	C							2.86	1	1.0	Yes	Aminobiphenyl, 4-	92-67-1	3.7E-01	1.5E+00		3.0E-01					
				8.0E-02	P					0.21	1	1.0	Yes	Aminophenol, m-	591-27-5					1.6E+03	2.8E+05		1.6E+03	
				4.0E-03	X					0.62	1	1.0	Yes	Aminophenol, o-	95-55-6					8.0E+01	7.5E+03		7.9E+01	
				2.0E-02	P					0.04	1	1.0	Yes	Aminophenol, p-	123-30-8					4.0E+02	9.1E+04		4.0E+02	
				2.5E-03	I					5.5	1	0.9	Yes	Amitraz	33089-61-1					5.0E+01	9.8E+00		8.2E+00	
						5.0E-01	I	V		0.23	1	1.0		Ammonia	7664-41-7									
				2.0E-03	X					1.44	1	1.0	Yes	Ammonium Picrate	131-74-8					4.0E+01	2.7E+03		4.0E+01	
				2.0E-01	I						1	1.0	Yes	Ammonium Sulfamate	7773-06-0					4.0E+03	9.1E+05		4.0E+03	
						3.0E-03	X	V		0.89	1	1.0	Yes	Amyl Alcohol, tert-	75-85-4							6.3E+00	6.3E+00	
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I			0.9	1	1.0	Yes	Aniline	62-53-3	1.4E+03	6.9E+04		1.3E+03	1.4E+02	7.7E+03		1.4E+02	
4.0E-02	P			2.0E-03	X					3.39	1	0.9	Yes	Anthraquinone, 9,10-	84-65-1	1.9E+02	5.1E+02		1.4E+02	4.0E+01	1.1E+02		3.0E+01	
				4.0E-04	I	3.0E-04	A				0.15	1.0	Yes	Antimony (metallic)	7440-36-0					8.0E+00	2.7E+02		7.8E+00	6.0E+00
				5.0E-04	H						0.15	1.0	Yes	Antimony Pentoxide	1314-60-9					1.0E+01	3.4E+02		9.7E+00	
				4.0E-04	H						0.15	1.0	Yes	Antimony Tetroxide	1332-81-6					8.0E+00	2.7E+02		7.8E+00	
						2.0E-04	I				0.15	1.0	Yes	Antimony Trioxide	1309-64-4									
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C				1	1.0	Yes	Arsenic, Inorganic	7440-38-2	5.2E+00	9.7E+02		5.2E+00	6.0E+00	1.4E+03		6.0E+00	1.0E+01
				3.5E-06	C	5.0E-05	I				1	1.0	Yes	Arsine	7784-42-1					7.0E-02	1.6E+01		7.0E-02	
											1	0.0		Asbestos (units in fibers)	1332-21-4									7.0E+06(G)
				3.6E-02	O					-0.27	1	1.0	Yes	Asulam	3337-71-1					7.2E+02	5.8E+05		7.2E+02	
2.3E-01	C			3.0E-03	A					2.61	1	1.0	Yes	Atrazine	1912-24-9	3.4E+01	2.8E+02		3.0E+01	6.0E+01	5.3E+02		5.4E+01	3.0E+00
8.8E-01	C	2.5E-04	C							2.98	1	0.9	Yes	Auramine	492-80-8	8.9E+00	6.3E+01		7.8E+00					
				4.0E-04	I					4.48	1	1.0	No	Avermectin B1	65195-55-3					8.0E+00			8.0E+00	
				3.0E-03	A	1.0E-02	A			2.75	1	1.0	Yes	Azinphos-methyl	86-50-0					6.0E+01	8.3E+02		5.6E+01	
1.1E-01	I	3.1E-05	I					V		3.82	1	1.0	Yes	Azobenzene	103-33-3	7.1E+01	7.3E+01	1.8E+01	1.2E+01					
				1.0E+00	P	7.0E-06	P			-1.7	1	1.0	Yes	Azodicarbonamide	123-77-3					2.0E+04	6.8E+07		2.0E+04	
				2.0E-01	I	5.0E-04	H				0.07	1.0	Yes	Barium	7440-39-3					4.0E+03	6.4E+04		3.8E+03	2.0E+03
				5.0E-03	O			V		5.29	1	0.8	Yes	Benfluralin	1861-40-1					1.0E+02	4.0E+01		2.8E+01	
				5.0E-02	I					2.12	1	1.0	Yes	Benomyl	17804-35-2					1.0E+03	3.0E+04		9.7E+02	
				2.0E-01	I					2.18	1	1.0	Yes	Bensulfuron-methyl	83055-99-6					4.0E+03	2.4E+05		3.9E+03	
				3.0E-02	I					2.34	1	1.0	Yes	Bentazon	25057-89-0					6.0E+02	9.4E+03		5.7E+02	
4.0E-03	P			1.0E-01	I			V		1.48	1	1.0	Yes	Benzaldehyde	100-52-7	1.9E+03	4.4E+04		1.9E+03	2.0E+03	4.9E+04		1.9E+03	
5.5E-02	I	7.8E-06	I	4.0E-03	I	3.0E-02	I	V		2.13	1	1.0	Yes	Benzene	71-43-2	1.4E+02	9.8E+02	7.2E+01	4.6E+01	8.0E+01	6.1E+02	6.3E+01	3.3E+01	5.0E+00
1.0E-01	X			3.0E-04	X					-3.7267	1	1.0	No	Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1	7.8E+01			7.8E+01	6.0E+00			6.0E+00	
				1.0E-03	P			V		2.52	1	1.0	Yes	Benzenethiol	108-98-5					2.0E+01	1.0E+02		1.7E+01	
2.3E+02	I	6.7E-02	I	3.0E-03	I				M	1.34	1	1.0	Yes	Benzidine	92-87-5	1.1E-02	5.0E-01		1.1E-02	6.0E+01	3.0E+03		5.9E+01	
				4.0E+00	I					1.87	1	1.0	Yes	Benzoic Acid	65-85-0					8.0E+04	1.2E+06		7.5E+04	
1.3E+01	I							V		3.9	1	1.0	Yes	Benzotrichloride	98-07-7	6.0E-01	6.0E-01		3.0E-01					
				1.0E-01	P					1.1	1	1.0	Yes	Benzyl Alcohol	100-51-6					2.0E+03	8.9E+04		2.0E+03	
1.7E-01	I	4.9E-05	C	2.0E-03	P	1.0E-03	P	V		2.3	1	1.0	Yes	Benzyl Chloride	100-44-7	4.6E+01	3.4E+02	1.1E+01	8.9E+00	4.0E+01	3.2E+02	2.1E+00	2.0E+00	
		2.4E-03	I	2.0E-03	I	2.0E-05	I				0.007	1.0	Yes	Beryllium and compounds	7440-41-7					4.0E+01	6.4E+01		2.5E+01	4.0E+00
				9.0E-03	P					4.48	1	0.9	Yes	Bifenox	42576-02-3					1.8E+02	2.3E+02		1.0E+02	
				1.5E-02	I					6	1	0.0	Yes	Biphenthrin	82657-04-3					3.0E+02			3.0E+02	
8.0E-03	I			5.0E-01	I	4.0E-04	X	V		4.01	1	1.0	Yes	Biphenyl, 1,1'-	92-52-4	9.7E+02	6.5E+02		3.9E+02	1.0E+04	7.3E+03	8.3E-01	8.3E-01	
				4.0E-02	I			V		2.48	1	1.0	Yes	Bis(2-chloro-1-methylethyl) ether	108-60-1					8.0E+02	6.5E+03		7.1E+02	
				3.0E-03	P					1.3	1	1.0	Yes	Bis(2-chloroethoxy)methane	111-91-1					6.0E+01	3.0E+03		5.9E+01	
1.1E+00	I	3.3E-04	I					V		1.29	1	1.0	Yes	Bis(2-chloroethyl)ether	111-44-4	7.1E+00	2.7E+02	1.7E+00	1.4E+00					
2.2E+02	I	6.2E-02	I					V		0.57	1	1.0	Yes	Bis(chloromethyl)ether	542-88-1	3.5E-02	3.4E+00	9.1E-03	7.2E-03					
				5.0E-02	I					3.32	1	1.0	Yes	Bisphenol A	80-05-7					1.0E+03	3.2E+03		7.7E+02	
				2.0E-01	I	2.0E-02	H				1	1.0	Yes	Boron And Borates Only	7440-42-8					4.0E+03	9.1E+05		4.0E+03	
				2.0E+00	P	2.0E-02	P	V		1.16	1	1.0	Yes	Boron Trichloride	10294-34-5					4.0E+04	9.1E+06	4.2E+01	4.2E+01	
				4.0E-02	C	1.3E-02	C	V		0.22	1	1.0	Yes	Boron Trifluoride	7637-07-2					8.0E+02	1.8E+05	2.7E+01	2.6E+01	
7.0E-01	I			4.0E-03	I						1	1.0	Yes	Bromate	15541-45-4	1.1E+01	2.1E+03		1.1E+01	8.0E+01	1.8E+04		8.0E+01	1.0E+01
				1.0E-04	X	6.0E-05	X	V		1.92	1	1.0	Yes	Bromo-2-chloroethane, 1-	107-04-0					2.0E+00	3.2E+01	1.3E-01	1.2E-01	

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1				
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	log K_{ow} (unitless)	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
				3.0E-04	X			V		2.92	1	1.0	Yes	Bromo-3-fluorobenzene, 1-	1073-06-9					6.0E+00	2.6E+01		4.9E+00	
				3.0E-04	X			V		3.08	1	1.0	Yes	Bromo-4-fluorobenzene, 1-	460-00-4					6.0E+00	2.0E+01		4.6E+00	
										0.41	1	1.0	Yes	Bromoacetic acid	79-08-3									6.0E+01(G)
				8.0E-03	I	6.0E-02	I	V		2.99	1	1.0	Yes	Bromobenzene	108-86-1					1.6E+02	5.4E+02	1.3E+02	6.2E+01	
						4.0E-02	X	V		1.41	1	1.0	Yes	Bromochloromethane	74-97-5							8.3E+01	8.3E+01	
6.2E-02	I	3.7E-05	C	8.0E-03	P			V		2	1	1.0	Yes	Bromodichloromethane	75-27-4	1.3E+02	1.9E+03	1.5E+01	1.3E+01	1.6E+02	2.6E+03		1.5E+02	8.0E+01(G)
7.9E-03	I	1.1E-06	I	2.0E-02	I			V		2.4	1	1.0	Yes	Bromoform	75-25-2	9.9E+02	1.4E+04	5.1E+02	3.3E+02	4.0E+02	6.2E+03		3.8E+02	8.0E+01(G)
				1.4E-03	I	5.0E-03	I	V		1.19	1	1.0	Yes	Bromomethane	74-83-9					2.8E+01	1.0E+03	1.0E+01	7.5E+00	
				5.0E-03	H			V		5.21	1	0.8	Yes	Bromophos	2104-96-3					1.0E+02	5.5E+01		3.5E+01	
						1.0E-01	A	V		2.1	1	1.0	Yes	Bromopropane, 1-	106-94-5							2.1E+02	2.1E+02	
1.0E-01	O			1.5E-02	O					2.8	1	0.9	Yes	Bromoxynil	1689-84-5	7.6E+01	3.1E+02		6.1E+01	3.0E+02	1.3E+03		2.5E+02	
1.0E-01	O			1.5E-02	O			V		5.4	1	0.8	Yes	Bromoxynil Octanoate	1689-99-2	7.6E+01	3.6E+01		2.4E+01	3.0E+02	1.6E+02		1.0E+02	
6.0E-01	C	3.0E-05	I			2.0E-03	I	V		1.99	1	1.0	Yes	Butadiene, 1,3-	106-99-0	1.3E+01	9.2E+01	1.9E+01	7.1E+00			4.2E+00	4.2E+00	
				1.0E-01	I			V		0.88	1	1.0	Yes	Butanol, N-	71-36-3					2.0E+03	1.0E+05		2.0E+03	
5.0E-04	I			4.0E-01	I	5.0E+00	I	V		0.35	1	1.0	Yes	Butyl Alcohol, t-	75-65-0	1.6E+04	1.6E+06		1.5E+04	8.0E+03	9.0E+05	1.0E+04	4.5E+03	
				2.0E+00	P	3.0E+01	P	V		0.61	1	1.0	Yes	Butyl alcohol, sec-	78-92-2					4.0E+04	3.0E+06	6.3E+04	2.4E+04	
				5.0E-02	I			V		4.15	1	1.0	Yes	Butylate	2008-41-5					1.0E+03	8.5E+02		4.6E+02	
2.0E-04	C	5.7E-08	C							3.5	1	0.8	Yes	Butylated hydroxyanisole	25013-16-5	3.9E+04	2.5E+04		1.5E+04					
3.6E-03	P			3.0E-01	P					5.1	1	1.0	Yes	Butylated hydroxytoluene	128-37-0	2.2E+03	4.0E+02		3.4E+02	6.0E+03	1.2E+03		1.0E+03	
				5.0E-02	P			V		4.38	1	1.0	No	Butylbenzene, n-	104-51-8					1.0E+03			1.0E+03	
				1.0E-01	X			V		4.57	1	1.0	No	Butylbenzene, sec-	135-98-8					2.0E+03			2.0E+03	
				1.0E-01	X			V		4.11	1	1.0	Yes	Butylbenzene, tert-	98-06-6					2.0E+03	1.1E+03		6.9E+02	
				2.0E-02	A					0.36	1	1.0	Yes	Cacodylic Acid	75-60-5					4.0E+02	6.7E+04		4.0E+02	
		1.8E-03	I	1.0E-04	A	1.0E-05	A				0.025	1.0		Cadmium (Diet)	7440-43-9									
		1.8E-03	I	1.0E-04	A	1.0E-05	A				0.05	1.0	Yes	Cadmium (Water)	7440-43-9					2.0E+00	2.3E+01		1.8E+00	5.0E+00
				5.0E-01	I	2.2E-03	C			-0.19	1	1.0	Yes	Caprolactam	105-60-2					1.0E+04	9.0E+05		9.9E+03	
1.5E-01	C	4.3E-05	C	2.0E-03	I					3.8	1	0.9	Yes	Captafol	2425-06-1	5.2E+01	1.8E+02		4.0E+01	4.0E+01	1.5E+02		3.2E+01	
2.3E-03	C	6.6E-07	C	1.3E-01	I					2.8	1	1.0	Yes	Captan	133-06-2	3.4E+03	3.6E+04		3.1E+03	2.6E+03	3.0E+04		2.4E+03	
				1.0E-01	I					2.36	1	1.0	Yes	Carbaryl	63-25-2					2.0E+03	2.4E+04		1.8E+03	
				5.0E-03	I					2.32	1	1.0	Yes	Carbofuran	1563-66-2					1.0E+02	1.4E+03		9.4E+01	4.0E+01
				1.0E-01	I	7.0E-01	I	V		1.94	1	1.0	Yes	Carbon Disulfide	75-15-0					2.0E+03	2.0E+04	1.5E+03	8.1E+02	
7.0E-02	I	6.0E-06	I	4.0E-03	I	1.0E-01	I	V		2.83	1	1.0	Yes	Carbon Tetrachloride	56-23-5	1.1E+02	4.3E+02	9.4E+01	4.6E+01	8.0E+01	3.4E+02	2.1E+02	4.9E+01	5.0E+00
						1.0E-01	P	V		-1.33	1	1.0	Yes	Carbonyl Sulfide	463-58-1							2.1E+02	2.1E+02	
				1.0E-02	I					5.57	1	0.8	Yes	Carbosulfan	55285-14-8					2.0E+02	6.9E+01		5.1E+01	
				1.0E-01	I					2.14	1	1.0	Yes	Carboxin	5234-68-4					2.0E+03	4.1E+04		1.9E+03	
						9.0E-04	I				1	1.0	Yes	Ceric oxide	1306-38-3									
				1.0E-01	I			V		0.99	1	1.0	Yes	Chloral Hydrate	302-17-0					2.0E+03	1.5E+05		2.0E+03	
				1.5E-02	I					1.9	1	1.0	Yes	Chloramben	133-90-4					3.0E+02	7.4E+03		2.9E+02	
											1	0.0	No	Chloramines, Organic	E701235									4.0E+03(G)
4.0E-01	H									2.22	1	1.0	Yes	Chloranil	118-75-2	1.9E+01	3.5E+02		1.8E+01					
				5.0E-04	G			V		6.1	1	0.7	Yes	Chlordane (alpha)	5103-71-9					1.0E+01	5.6E+00		3.6E+00	
				5.0E-04	G			V		6.22	1	0.7	No	Chlordane (gamma)	5103-74-2					1.0E+01			1.0E+01	
3.5E-01	I	1.0E-04	I	5.0E-04	I	7.0E-04	I	V		6.16	1	0.7	Yes	Chlordane (technical mixture)	12789-03-6	2.2E+01	3.6E+00	5.6E+00	2.0E+00	1.0E+01	1.8E+00	1.5E+00	7.4E-01	2.0E+00
1.0E+01	I	4.6E-03	C	3.0E-04	I					5.41	1	0.8	Yes	Chlordecone (Kepone)	143-50-0	7.8E-01	6.5E-01		3.5E-01	6.0E+00	5.4E+00		2.9E+00	
				7.0E-04	A					3.81	1	0.9	Yes	Chlorfenvinphos	470-90-6					1.4E+01	5.6E+01		1.1E+01	
				9.0E-02	O					2.5	1	1.0	Yes	Chlorimuron, Ethyl-	90982-32-4					1.8E+03	6.8E+04		1.8E+03	
				1.0E-01	I	1.5E-04	A	V		0.85	1	1.0	Yes	Chlorine	7782-50-5					2.0E+03	4.6E+05	3.0E-01	3.0E-01	4.0E+03(G)
				3.0E-02	I	2.0E-04	I	V			1	1.0	Yes	Chlorine Dioxide	10049-04-4					6.0E+02	1.4E+05	4.2E-01	4.2E-01	8.0E+02(G)
				3.0E-02	I						1	1.0	Yes	Chlorite (Sodium Salt)	7758-19-2					6.0E+02	1.4E+05		6.0E+02	1.0E+03
						5.0E+01	I	V		2.05	1	1.0	Yes	Chloro-1,1-difluoroethane, 1-	75-68-3							1.0E+05	1.0E+05	
		3.0E-04	I	2.0E-02	H	2.0E-02	I	V		2.53	1	1.0	Yes	Chloro-1,3-butadiene, 2-	126-99-8			1.9E+00	1.9E+00	4.0E+02	1.8E+03	4.2E+01	3.7E+01	
4.6E-01	H									2.27	1	1.0	Yes	Chloro-2-methylaniline HCl, 4-	3165-93-3	1.7E+01	5.1E+04		1.7E+01					
1.0E-01	P	7.7E-05	C	3.0E-03	X					2.27	1	1.0	Yes	Chloro-2-methylaniline, 4-	95-69-2	7.8E+01	6.6E+02		7.0E+01	6.0E+01	5.6E+02		5.4E+01	
2.7E-01	X							V		0.09	1	1.0	Yes	Chloroacetaldehyde, 2-	107-20-0	2.9E+01	4.6E+03		2.9E+01					
										0.22	1	1.0	Yes	Chloroacetic Acid	79-11-8									6.0E+01(G)
						3.0E-05	I			1.93	1	1.0	Yes	Chloroacetophenone, 2-	532-27-4									
2.0E-01	P			5.0E-04	P					1.83	1	1.0	Yes	Chloroaniline, p-	106-47-8	3.9E+01	5.9E+02		3.7E+01	1.0E+01	1.7E+02		9.5E+00	
				2.0E-02	I	5.0E-02	P	V		2.84	1	1.0	Yes	Chlorobenzene	108-90-7					4.0E+02	1.3E+03	1.0E+02	7.8E+01	1.0E+02
				1.0E-01	X					-0.52	1	1.0	Yes	Chlorobenzene sulfonic acid, p-	98-66-8					2.0E+03	1.8E+06		2.0E+03	
1.1E-01	C	3.1E-05	C	2.0E-02	I					4.74	1	0.8	Yes	Chlorobenzilate	510-15-6	7.1E+01	5.6E+01		3.1E+01	4.0E+02	3.5E+02		1.9E+02	
				3.0E-02	X					2.65	1	1.0	Yes	Chlorobenzoic Acid, p-	74-11-3					6.0E+02	3.4E+03		5.1E+02	
		8.6E-06	C	3.0E-03	P	3.0E-01	P	V		3.6	1	1.0	Yes	Chlorobenzotrifluoride, 4-	98-56-6			6.5E+01	6.5E+01	6.0E+01	9.3E+01	6.3E+02	3.5E+01	
				4.0E-02	P			V		2.64	1	1.0	Yes	Chlorobutane, 1-	109-69-3					8.0E+02	3.1E+03		6.4E+02	
						5.0E+01	I	V		1.08	1	1.0	Yes	Chlorodifluoromethane	75-45-6							1.0E+05	1.0E+05	
				2.0E-02	P			V		0.03	1	1.0	Yes	Chloroethanol, 2-	107-07-3					4.0E+02	7.7E+04		4.0E+02	
3.1E-02	C	2.3E-05	I	1.0E-02	I	9.8E-02	A	V		1.97	1	1.0	Yes	Chloroform	67-66-3	2.5E+02	2.9E+03	2.4E+01	2.2E+01	2.0E+02	2.5E+03	2.0E+02	9.7E+01	8.0E+01(G)
						9.0E-02	I	V		0.91	1	1.0	Yes	Chloromethane	74-87-3							1.9E+02	1.9E+02	
2.4E+00	C	6.9E-04	C					V		0.32	1	1.0	Yes	Chloromethyl Methyl Ether	107-30-2	3.2E+00	3.7E+02	8.1E-01	6.5E-01					
3.0E-01	P			3.0E-03	P	1.0E-05	X			2.24	1	1.0	Yes	Chloronitrobenzene, o-	88-73-3	2.6E+01	2.6E+02		2.4E+01	6.0E+01	6.4E+02		5.5E+01	
6.0E-02	P			7.0E-04	P	2.0E-03	P			2.39	1	1.0	Yes	Chloronitrobenzene, p-	100-00-5	1.3E+02	1.0E+03		1.2E+02	1.4E+01	1.2E+02		1.3E+01	
				5.0E-03	I			V		2.15	1	1.0	Yes	Chlorophenol, 2-	95-57-8					1.0E+02	1.0E+03		9.1E+01	
						4.0E-04	C	V		2.09	1	1.0	Yes	Chloropicrin	76-06-2							8.3E-01	8.3E-01	
1.7E-02	C			1.5E-02	I					3.05	1	0.9	Yes	Chlorothalonil	1897-45-6	4.6E+02	2.9E+03		4.0E+02	3.0E+02	2.1E+03		2.6E+02	
				2.0E-02	I			V		3.42	1	1.0	Yes	Chlorotoluene, o-	95-49-8					4.0E+02	5.8E+02		2.4E+02	
				2.0E-02	X			V		3.33	1	1.0	Yes	Chlorotoluene, p-	106-43-4					4.0E+02	6.6E+02		2.5E+02	
2.4E+02	C	6.9E-02	C							-1.02	1	1.0	Yes	Chlorozotocin	54749-90-5	3.2E-02	1.0E+02		3.2E-02					

RMLS TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1				
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	log K_{ow} (unitless)	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
				5.0E-03	O					3.51	1	0.9	Yes	Chlorpropham	101-21-3					1.0E+02	2.5E+02		7.1E+01	
				1.0E-03	A					4.96	1	0.8	Yes	Chlorpyrifos	2921-88-2					2.0E+01	1.5E+01		8.4E+00	
				1.0E-02	H					4.31	1	0.9	Yes	Chlorpyrifos Methyl	5598-13-0					2.0E+02	2.9E+02		1.2E+02	
				5.0E-02	O					2	1	1.0	Yes	Chlorsulfuron	64902-72-3					1.0E+03	5.7E+04		9.9E+02	
				1.0E-02	I					4.28	1	0.9	Yes	Chlorthal-dimethyl	1861-32-1					2.0E+02	3.3E+02		1.2E+02	
				8.0E-04	H					5.8	1	0.8	Yes	Chlorthiophos	60238-56-4					1.6E+01	3.4E+00		2.8E+00	
				1.5E+00	I						0.013	1.0	Yes	Chromium(III), Insoluble Salts	16065-83-1					3.0E+04	8.9E+04		2.2E+04	
5.0E-01	C	8.4E-02	G	3.0E-03	I	1.0E-04	I		M		0.025	1.0	Yes	Chromium(VI)	18540-29-9	5.0E+00	1.2E+01		3.5E+00	6.0E+01	1.7E+02		4.4E+01	
											0.013	1.0	Yes	Chromium, Total	7440-47-3									1.0E+02
				1.3E-02	I					3.1	1	0.9	Yes	Clofentezine	74115-24-5					2.6E+02	2.1E+03		2.3E+02	
		9.0E-03	P	3.0E-04	P	6.0E-06	P				1	1.0	Yes	Cobalt	7440-48-4					6.0E+00	3.4E+03		6.0E+00	
		6.2E-04	I					V	M		1	0.0		Coke Oven Emissions	E649830									
				4.0E-02	H						1	1.0	Yes	Copper	7440-50-8					8.0E+02	1.8E+05		8.0E+02	1.3E+03
				5.0E-02	I	6.0E-01	C			1.96	1	1.0	Yes	Cresol, m-	108-39-4					1.0E+03	1.2E+04		9.3E+02	
				5.0E-02	I	6.0E-01	C			1.95	1	1.0	Yes	Cresol, o-	95-48-7					1.0E+03	1.2E+04		9.3E+02	
				2.0E-02	P	6.0E-01	C			1.94	1	1.0	Yes	Cresol, p-	106-44-5					4.0E+02	4.9E+03		3.7E+02	
				1.0E-01	A					3.1	1	1.0	Yes	Cresol, p-chloro-m-	59-50-7					2.0E+03	5.2E+03		1.4E+03	
				1.0E-01	A	6.0E-01	C			1.95	1	0.9	Yes	Cresols	1319-77-3					2.0E+03	6.7E+03		1.5E+03	
1.9E+00	H			1.0E-03	P			V		0.6	1	1.0	Yes	Crotonaldehyde, trans-	123-73-9	4.1E+00	2.7E+02		4.0E+00	2.0E+01	1.5E+03		2.0E+01	
				1.0E-01	I	4.0E-01	I	V		3.66	1	1.0	Yes	Cumene	98-82-8					2.0E+03	1.9E+03	8.3E+02	4.5E+02	
2.2E-01	C	6.3E-05	C							-1.73	1	1.0	Yes	Cupferron	135-20-6	3.5E+01	1.3E+06		3.5E+01					
8.4E-01	H			2.0E-03	H					2.22	1	1.0	Yes	Cyanazine	21725-46-2	9.3E+00	1.6E+02		8.8E+00	4.0E+01	7.6E+02		3.8E+01	
														Cyanides										
				1.0E-03	I						1	1.0	Yes	~Calcium Cyanide	592-01-8					2.0E+01	4.6E+03		2.0E+01	
				5.0E-03	I						1	1.0	Yes	~Copper Cyanide	544-92-3					1.0E+02	2.3E+04		1.0E+02	
				6.0E-04	I	8.0E-04	G	V			1	1.0	Yes	~Cyanide (CN-)	57-12-5					1.2E+01	2.7E+03	1.7E+00	1.5E+00	2.0E+02
				1.0E-03	I			V		0.07	1	1.0	Yes	~Cyanogen	460-19-5					2.0E+01	5.1E+03		2.0E+01	
				9.0E-02	I			V			1	1.0	Yes	~Cyanogen Bromide	506-68-3					1.8E+03	1.6E+06		1.8E+03	
				5.0E-02	I			V			1	1.0	Yes	~Cyanogen Chloride	506-77-4					1.0E+03	5.8E+05		1.0E+03	
				6.0E-04	I	8.0E-04	I	V		-0.25	1	1.0	Yes	~Hydrogen Cyanide	74-90-8					1.2E+01	2.7E+03	1.7E+00	1.5E+00	
				2.0E-03	I						1	1.0	Yes	~Potassium Cyanide	151-50-8					4.0E+01	4.6E+03		4.0E+01	
				5.0E-03	I						0.04	1.0	Yes	~Potassium Silver Cyanide	506-61-6					1.0E+02	4.6E+02		8.2E+01	
				1.0E-01	I						0.04	1.0	Yes	~Silver Cyanide	506-64-9					2.0E+03	1.8E+04		1.8E+03	
				1.0E-03	I						1	1.0	Yes	~Sodium Cyanide	143-33-9					2.0E+01	4.6E+03		2.0E+01	2.0E+02
				2.0E-04	P						1	0.0	Yes	~Thiocyanates	E1790665					4.0E+00	9.1E+02		4.0E+00	
				2.0E-04	X			V		0.58	1	1.0	Yes	~Thiocyanic Acid	463-56-9					4.0E+00	9.1E+02		4.0E+00	
				5.0E-02	I						1	1.0	Yes	~Zinc Cyanide	557-21-1					1.0E+03	3.8E+05		1.0E+03	
						6.0E+00	I	V		3.44	1	1.0	Yes	Cyclohexane	110-82-7							1.3E+04	1.3E+04	
2.0E-02	X			2.0E-02	X					4.72	1	0.9	Yes	Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	3.9E+02	9.6E+02		2.8E+02	4.0E+02	1.1E+03		2.9E+02	
				5.0E+00	I	7.0E-01	P	V		0.81	1	1.0	Yes	Cyclohexanone	108-94-1					1.0E+05	6.5E+06	1.5E+03	1.4E+03	
				5.0E-03	P	1.0E+00	X	V		2.86	1	1.0	Yes	Cyclohexene	110-83-8					1.0E+02	2.5E+02	2.1E+03	7.0E+01	
				2.0E-01	I			V		1.49	1	1.0	Yes	Cyclohexylamine	108-91-8					4.0E+03	9.3E+04		3.8E+03	
				2.5E-02	I					5.95	1	0.7	Yes	Cyfluthrin	68359-37-5					5.0E+02	1.6E+02		1.2E+02	
				7.2E-02	O					6.6	1	0.7	No	Cypermethrin	52315-07-8					1.4E+03			1.4E+03	
				5.0E-01	O					-0.061	1	1.0	Yes	Cyromazine	66215-27-8					1.0E+04	8.0E+05		9.9E+03	
2.4E-01	I	6.9E-05	C	3.0E-05	X					6.02	1	0.8	Yes	DDD, p,p`- (DDD)	72-54-8	3.2E+01	3.5E+00		3.2E+00	6.0E-01	7.1E-02		6.3E-02	
3.4E-01	I	9.7E-05	C	3.0E-04	X			V		6.51	1	0.8	No	DDE, p,p'-	72-55-9	2.3E+01		5.8E+00	4.6E+00	6.0E+00			6.0E+00	
3.4E-01	I	9.7E-05	I	5.0E-04	I					6.91	1	0.7	No	DDT	50-29-3	2.3E+01			2.3E+01	1.0E+01			1.0E+01	
				3.0E-02	I					0.78	1	1.0	Yes	Dalapon	75-99-0					6.0E+02	5.5E+04		6.0E+02	2.0E+02
1.8E-02	C	5.1E-06	C	1.5E-01	I					-1.5	1	1.0	Yes	Daminozide	1596-84-5	4.3E+02	1.3E+06		4.3E+02	3.0E+03	1.0E+07		3.0E+03	
7.0E-04	I			7.0E-03	I					12.11	1	0.0	No	Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5	1.1E+04			1.1E+04	1.4E+02			1.4E+02	
				4.0E-05	I					3.21	1	0.8	Yes	Demeton	8065-48-3					8.0E-01	8.8E-01		4.2E-01	
1.2E-03	I			6.0E-01	I					6.11	1	0.0	Yes	Di(2-ethylhexyl)adipate	103-23-1	6.5E+03			6.5E+03	1.2E+04			1.2E+04	4.0E+02
6.1E-02	H									4.49	1	0.9	Yes	Diallate	2303-16-4	1.3E+02	9.2E+01		5.4E+01					
				7.0E-04	A					3.81	1	0.9	Yes	Diazinon	333-41-5					1.4E+01	3.9E+01		1.0E+01	
				1.0E-02	X			V		4.38	1	1.0	Yes	Dibenzothiophene	132-65-0					2.0E+02	9.6E+01		6.5E+01	
8.0E-01	P	6.0E-03	P	2.0E-04	P	2.0E-04	I	V	M	2.96	1	1.0	Yes	Dibromo-3-chloropropane, 1,2-	96-12-8	3.1E+00	1.7E+01	3.4E-02	3.3E-02	4.0E+00	2.4E+01	4.2E-01	3.7E-01	2.0E-01
										0.7	1	1.0	Yes	Dibromoacetic acid	631-64-1									6.0E+01(G)
				4.0E-04	X			V		3.75	1	0.9	Yes	Dibromobenzene, 1,3-	108-36-1					8.0E+00	1.6E+01		5.3E+00	
				1.0E-02	I			V		3.79	1	0.9	Yes	Dibromobenzene, 1,4-	106-37-6					2.0E+02	3.7E+02		1.3E+02	
8.4E-02	I			2.0E-02	I			V		2.16	1	1.0	Yes	Dibromochloromethane	124-48-1	9.3E+01	1.4E+03		8.7E+01	4.0E+02	6.7E+03		3.8E+02	8.0E+01(G)
2.0E+00	I	6.0E-04	I	9.0E-03	I	9.0E-03	I	V		1.96	1	1.0	Yes	Dibromoethane, 1,2-	106-93-4	3.9E+00	7.1E+01	9.4E-01	7.5E-01	1.8E+02	3.6E+03	1.9E+01	1.7E+01	5.0E-02
						4.0E-03	X	V		1.7	1	1.0	Yes	Dibromomethane (Methylene Bromide)	74-95-3							8.3E+00	8.3E+00	
				3.0E-04	P						1	0.0	No	Dibutyltin Compounds	E1790661					6.0E+00			6.0E+00	
				3.0E-02	I					2.21	1	1.0	Yes	Dicamba	1918-00-9					6.0E+02	1.0E+04		5.7E+02	
											1	0.0	No	Dichloramine	3400-09-7									4.0E+03(G)
		4.2E-03	P					V		2.6	1	1.0	Yes	Dichloro-2-butene, 1,4-	764-41-0			1.3E-01	1.3E-01					
		4.2E-03	P					V		2.6	1	1.0	Yes	Dichloro-2-butene, cis-1,4-	1476-11-5			1.3E-01	1.3E-01					
		4.2E-03	P					V		2.6	1	1.0	Yes	Dichloro-2-butene, trans-1,4-	110-57-6			1.3E-01	1.3E-01					
5.0E-02	I			4.0E-03	I					0.92	1	1.0	Yes	Dichloroacetic Acid	79-43-6	1.6E+02	9.6E+03		1.5E+02	8.0E+01	5.4E+03		7.9E+01	6.0E+01(G)
				9.0E-02	I	2.0E-01	H	V		3.43	1	1.0	Yes	Dichlorobenzene, 1,2-	95-50-1					1.8E+03	2.9E+03	4.2E+02	3.0E+02	6.0E+02
5.4E-03	C	1.1E-05	C	7.0E-02	A	8.0E-01	I	V		3.44	1	1.0	Yes	Dichlorobenzene, 1,4-	106-46-7	1.4E+03	2.1E+03	5.1E+01	4.8E+01	1.4E+03	2.2E+03	1.7E+03	5.7E+02	7.5E+01
4.5E-01	I	3.4E-04	C							3.51	1	1.0	Yes	Dichlorobenzidine, 3,3'-	91-94-1	1.7E+01	4.5E+01		1.3E+01					
				9.0E-03	X					4.44	1	0.9	Yes	Dichlorobenzophenone, 4,4'-	90-98-2					1.8E+02	1.4E+02		7.8E+01	
				2.0E-01	I	1.0E-01	X	V		2.16	1	1.0	Yes	Dichlorodifluoromethane	75-71-8					4.0E+03	3.8E+04	2.1E+02	2.0E+02	
5.7E-03	C	1.6E-06	C	2.0E-01	P			V		1.79	1	1.0	Yes	Dichloroethane, 1,1-	75-34-3	1.4E+03	1.8E+04	3.5E+02	2.8E+02	4.0E+03	5.8E+04		3.8E+03	

RMLS TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1				
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	log K_{ow} (unitless)	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
9.1E-02	I	2.6E-05	I	6.0E-03	X	7.0E-03	P	V		1.48	1	1.0	Yes	Dichloroethane, 1,2-	107-06-2	8.6E+01	1.8E+03	2.2E+01	1.7E+01	1.2E+02	2.8E+03	1.5E+01	1.3E+01	5.0E+00
				5.0E-02	I	2.0E-01	I	V		2.13	1	1.0	Yes	Dichloroethylene, 1,1-	75-35-4					1.0E+03	8.5E+03	4.2E+02	2.8E+02	7.0E+00
				2.0E-03	I			V		1.86	1	1.0	Yes	Dichloroethylene, cis-1,2-	156-59-2					4.0E+01	3.6E+02		3.6E+01	7.0E+01
				2.0E-02	I	4.0E-02	X	V		2.09	1	1.0	Yes	Dichloroethylene, trans-1,2-	156-60-5					4.0E+02	3.6E+03	8.3E+01	6.8E+01	1.0E+02
				3.0E-03	I					3.06	1	1.0	Yes	Dichlorophenol, 2,4-	120-83-2					6.0E+01	1.9E+02		4.6E+01	
				1.0E-02	I					2.81	1	1.0	Yes	Dichlorophenoxy Acetic Acid, 2,4-	94-75-7					2.0E+02	1.4E+03		1.7E+02	7.0E+01
3.7E-02	P	3.7E-06	P	4.0E-02	P	4.0E-03	I	V		1.98	1	1.0	Yes	Dichloropropane, 1,2-	78-87-5	2.1E+02	2.3E+03	1.5E+02	8.5E+01	8.0E+02	9.6E+03	8.3E+00	8.2E+00	5.0E+00
				2.0E-02	P			V		2	1	1.0	Yes	Dichloropropane, 1,3-	142-28-9					4.0E+02	4.6E+03		3.7E+02	
				3.0E-03	I					0.78	1	1.0	Yes	Dichloropropanol, 2,3-	616-23-9					6.0E+01	5.0E+03		5.9E+01	
1.0E-01	I	4.0E-06	I	3.0E-02	I	2.0E-02	I	V		2.04	1	1.0	Yes	Dichloropropene, 1,3-	542-75-6	7.8E+01	7.8E+02	1.4E+02	4.7E+01	6.0E+02	6.6E+03	4.2E+01	3.9E+01	
2.9E-01	I	8.3E-05	C	5.0E-04	I	5.0E-04	I			1.43	1	1.0	Yes	Dichlorvos	62-73-7	2.7E+01	1.4E+03		2.6E+01	1.0E+01	5.6E+02		9.9E+00	
				3.0E-05	O					0	1	1.0	Yes	Dicrotophos	141-66-2					6.0E-01	3.3E+02		6.0E-01	
				8.0E-02	P	3.0E-04	X	V		3.16	1	1.0	Yes	Dicyclopentadiene	77-73-6					1.6E+03	3.5E+03	6.3E-01	6.3E-01	
1.6E+01	I	4.6E-03	I	5.0E-05	I					5.4	1	0.8	Yes	Dieldrin	60-57-1	4.9E-01	2.7E-01		1.8E-01	1.0E+00	6.1E-01		3.8E-01	
		3.0E-04	C			5.0E-03	I				1	0.0		Diesel Engine Exhaust	E17136615									
				2.0E-03	P	2.0E-04	P			-1.43	1	1.0	Yes	Diethanolamine	111-42-2					4.0E+01	8.4E+04		4.0E+01	
				3.0E-02	P	1.0E-04	P			0.56	1	1.0	Yes	Diethylene Glycol Monobutyl Ether	112-34-5					6.0E+02	8.7E+04		6.0E+02	
				6.0E-02	P	3.0E-04	P			-0.54	1	1.0	Yes	Diethylene Glycol Monoethyl Ether	111-90-0					1.2E+03	7.8E+05		1.2E+03	
				1.0E-03	P			V		0.05	1	1.0	Yes	Diethylformamide	617-84-5					2.0E+01	4.3E+03		2.0E+01	
3.5E+02	C	1.0E-01	C							5.07	1	0.9	Yes	Diethylstilbestrol	56-53-1	2.2E-02	6.6E-03		5.1E-03					
				8.3E-02	O					0.65	1	1.0	Yes	Difenzoquat	43222-48-6					1.7E+03	7.5E+05		1.7E+03	
				2.0E-02	I					3.88	1	0.9	Yes	Diflubenzuron	35367-38-5					4.0E+02	1.0E+03		2.9E+02	
						4.0E+01	I	V		0.75	1	1.0	Yes	Difluoroethane, 1,1-	75-37-6							8.3E+04	8.3E+04	
						3.0E+01	X	V		2.29	1	1.0	Yes	Difluoropropane, 2,2-	420-45-1							6.3E+04	6.3E+04	
4.4E-02	C	1.3E-05	C					V		3.58	1	1.0	Yes	Dihydrosafrole	94-58-6	1.8E+02	2.3E+02	4.3E+01	3.0E+01					
						7.0E-01	P	V		1.52	1	1.0	Yes	Diisopropyl Ether	108-20-3							1.5E+03	1.5E+03	
				8.0E-02	I			V		1.03	1	1.0	Yes	Diisopropyl Methylphosphonate	1445-75-6					1.6E+03	1.3E+05		1.6E+03	
				2.2E-02	O					-0.17	1	1.0	Yes	Dimethipin	55290-64-7					4.4E+02	2.6E+05		4.4E+02	
				2.2E-03	O					0.78	1	1.0	Yes	Dimethoate	60-51-5					4.4E+01	7.0E+03		4.4E+01	
1.6E+00	P									1.81	1	1.0	Yes	Dimethoxybenzidine, 3,3'-	119-90-4	4.9E+00	1.6E+02		4.7E+00					
1.7E-03	P			6.0E-02	P					-0.61	1	1.0	Yes	Dimethyl methylphosphonate	756-79-6	4.6E+03	2.8E+06		4.6E+03	1.2E+03	8.1E+05		1.2E+03	
4.6E+00	C	1.3E-03	C							4.58	1	1.0	Yes	Dimethylamino azobenzene [p-]	60-11-7	1.7E+00	7.2E-01		5.0E-01					
5.8E-01	H									2.17	1	1.0	Yes	Dimethylaniline HCl, 2,4-	21436-96-4	1.3E+01	5.2E+04		1.3E+01					
2.0E-01	P			2.0E-03	X					1.68	1	1.0	Yes	Dimethylaniline, 2,4-	95-68-1	3.9E+01	7.1E+02		3.7E+01	4.0E+01	8.0E+02		3.8E+01	
2.7E-02	P			2.0E-03	I			V		2.31	1	1.0	Yes	Dimethylaniline, N,N-	121-69-7	2.9E+02	2.0E+03		2.5E+02	4.0E+01	3.1E+02		3.5E+01	
1.1E+01	P									2.34	1	1.0	Yes	Dimethylbenzidine, 3,3'-	119-93-7	7.1E-01	8.5E+00		6.5E-01					
				1.0E-01	P	3.0E-02	I	V		-1.01	1	1.0	Yes	Dimethylformamide	68-12-2					2.0E+03	1.8E+06	6.3E+01	6.1E+01	
				1.0E-04	X	2.0E-06	X	V		-1.19	1	1.0	Yes	Dimethylhydrazine, 1,1-	57-14-7					2.0E+00	2.2E+03	4.2E-03	4.2E-03	
5.5E+02	C	1.6E-01	C					V		-0.54	1	1.0	Yes	Dimethylhydrazine, 1,2-	540-73-8	1.4E-02	5.0E+00	3.5E-03	2.8E-03					
				2.0E-02	I					2.3	1	1.0	Yes	Dimethylphenol, 2,4-	105-67-9					4.0E+02	3.1E+03		3.6E+02	
				6.0E-04	I					2.36	1	1.0	Yes	Dimethylphenol, 2,6-	576-26-1					1.2E+01	8.5E+01		1.1E+01	
				1.0E-03	I					2.23	1	1.0	Yes	Dimethylphenol, 3,4-	95-65-8					2.0E+01	1.7E+02		1.8E+01	
4.5E-02	C	1.3E-05	C					V		2.58	1	1.0	Yes	Dimethylvinylchloride	513-37-1	1.7E+02	6.5E+02	4.3E+01	3.3E+01					
				8.0E-05	X					2.13	1	1.0	Yes	Dinitro-o-cresol, 4,6-	534-52-1					1.6E+00	2.6E+01		1.5E+00	
				2.0E-03	I					4.12	1	0.9	Yes	Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5					4.0E+01	5.4E+01		2.3E+01	
				4.0E-04	X	2.0E-03	X			1.89	1	1.0	Yes	Dinitroaniline, 3,5-	618-87-1					8.0E+00	1.7E+02		7.7E+00	
				1.0E-04	P					1.69	1	1.0	Yes	Dinitrobenzene, 1,2-	528-29-0					2.0E+00	5.3E+01		1.9E+00	
				1.0E-04	I					1.49	1	1.0	Yes	Dinitrobenzene, 1,3-	99-65-0					2.0E+00	7.3E+01		2.0E+00	
				1.0E-04	P					1.46	1	1.0	Yes	Dinitrobenzene, 1,4-	100-25-4					2.0E+00	7.6E+01		2.0E+00	
				2.0E-03	I					1.67	1	1.0	Yes	Dinitrophenol, 2,4-	51-28-5					4.0E+01	1.2E+03		3.9E+01	
6.8E-01	I									2.18	1	1.0	Yes	Dinitrotoluene Mixture, 2,4/2,6-	E1615210	1.1E+01	1.5E+02		1.1E+01					
3.1E-01	C	8.9E-05	C	2.0E-03	I					1.98	1	1.0	Yes	Dinitrotoluene, 2,4-	121-14-2	2.5E+01	4.3E+02		2.4E+01	4.0E+01	7.5E+02		3.8E+01	
1.5E+00	P			3.0E-04	X					2.1	1	1.0	Yes	Dinitrotoluene, 2,6-	606-20-2	5.2E+00	7.4E+01		4.9E+00	6.0E+00	9.3E+01		5.7E+00	
				1.0E-04	X					1.84	1	1.0	Yes	Dinitrotoluene, 2-Amino-4,6-	35572-78-2					2.0E+00	5.1E+01		1.9E+00	
				1.0E-04	X					1.84	1	1.0	Yes	Dinitrotoluene, 4-Amino-2,6-	19406-51-0					2.0E+00	5.1E+01		1.9E+00	
4.5E-01	X			9.0E-04	X					2.18	1	0.8	Yes	Dinitrotoluene, Technical grade	25321-14-6	1.7E+01	2.6E+01		1.0E+01	1.8E+01	3.0E+01		1.1E+01	
				1.0E-03	I					3.56	1	0.9	Yes	Dinoseb	88-85-7					2.0E+01	5.4E+01		1.5E+01	7.0E+00
1.0E-01	I	5.0E-06	I	3.0E-02	I	3.0E-02	I	V		-0.27	1	1.0	Yes	Dioxane, 1,4-	123-91-1	7.8E+01	2.3E+04	1.1E+02	4.6E+01	6.0E+02	1.9E+05	6.3E+01	5.7E+01	
														Dioxins										
6.2E+03	I	1.3E+00	I							8.21	1	0.0	No	~Hexachlorodibenzo-p-dioxin, Mixture	34465-46-8	1.3E-03			1.3E-03					
1.3E+05	C	3.8E+01	C	7.0E-10	I	4.0E-08	C	V		6.8	1	0.5	No	~TCDD, 2,3,7,8-	1746-01-6	6.0E-05		1.5E-05	1.2E-05	1.4E-05		8.3E-05	1.2E-05	3.0E-05
				3.0E-02	I					2.17	1	1.0	Yes	Diphenamid	957-51-7					6.0E+02	4.2E+03		5.3E+02	
						4.0E-04	X	V		4.21	1	1.0	Yes	Diphenyl Ether	101-84-8							8.3E-01	8.3E-01	
				8.0E-04	X					2.4	1	1.0	Yes	Diphenyl Sulfone	127-63-9					1.6E+01	2.0E+02		1.5E+01	
				1.0E-01	O					3.5	1	1.0	Yes	Diphenylamine	122-39-4					2.0E+03	3.4E+03		1.3E+03	
8.0E-01	I	2.2E-04	I							2.94	1	1.0	Yes	Diphenylhydrazine, 1,2-	122-66-7	9.7E+00	3.9E+01		7.8E+00					
				2.2E-03	I					2.3622	1	1.0	Yes	Diquat	2764-72-9					4.4E+01	4.7E+02		4.0E+01	2.0E+01
7.4E+00	C	2.1E-03	C							4.9	1	1.0	No	Direct Black 38	1937-37-7	1.1E+00			1.1E+00					
7.4E+00	C	2.1E-03	C							2.6	1	1.0	No	Direct Blue 6	2602-46-2	1.1E+00			1.1E+00					
6.7E+00	C	1.9E-03	C							-6.53	1	1.0	No	Direct Brown 95	16071-86-6	1.2E+00			1.2E+00					
				4.0E-05	I					4.02	1	0.9	Yes	Disulfoton	298-04-4					8.0E-01	1.3E+00		5.0E-01	
				1.0E-02	I			V		0.77	1	1.0	Yes	Dithiane, 1,4-	505-29-3					2.0E+02	1.6E+04		2.0E+02	
				2.0E-03	I					2.68	1	1.0	Yes	Diuron	330-54-1					4.0E+01	3.6E+02		3.6E+01	
				2.0E-02	O					1.15	1	1.0	Yes	Dodine	2439-10-3					4.0E+02	5.3E+04		4.0E+02	
				5.0E-02	O			V		3.21	1	1.0	Yes	EPTC	759-94-4					1.0E+03	3.0E+03		7.5E+02	
				6.0E-03	I			V		3.83	1	0.9	Yes	Endosulfan	115-29-7					1.2E+02	6.3E+02		1.0E+02	

RMLS TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1				
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	log K_{ow} (unitless)	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
				6.0E-03	P					3.66	1	0.9	Yes	Endosulfan Sulfate	1031-07-8					1.2E+02	9.1E+02		1.1E+02	
				2.0E-02	I					1.91	1	1.0	Yes	Endothall	145-73-3					4.0E+02	8.5E+03		3.8E+02	1.0E+02
				3.0E-04	I					5.2	1	0.8	Yes	Endrin	72-20-8					6.0E+00	3.7E+00		2.3E+00	2.0E+00
9.9E-03	I	1.2E-06	I	6.0E-03	P	1.0E-03	I	V		0.45	1	1.0	Yes	Epichlorohydrin	106-89-8	7.9E+02	7.9E+04	4.7E+02	2.9E+02	1.2E+02	1.3E+04	2.1E+00	2.0E+00	
						2.0E-02	I	V		0.86	1	1.0	Yes	Epoxybutane, 1,2-	106-88-7							4.2E+01	4.2E+01	
				4.0E-02	P					-1.18	1	1.0	Yes	Ethanol, 2-(2-methoxyethoxy)-	111-77-3					8.0E+02	1.3E+06		8.0E+02	
				5.0E-03	I					-0.22	1	1.0	Yes	Ethephon	16672-87-0					1.0E+02	4.2E+04		1.0E+02	
				5.0E-04	I					5.07	1	0.8	Yes	Ethion	563-12-2					1.0E+01	7.7E+00		4.3E+00	
				1.0E-01	P	6.0E-02	P	V		0.59	1	1.0	Yes	Ethoxyethanol Acetate, 2-	111-15-9					2.0E+03	2.3E+05	1.3E+02	1.2E+02	
				9.0E-02	P	4.0E-02	P	V		-0.32	1	1.0	Yes	Ethoxyethanol, 2-	110-80-5					1.8E+03	6.3E+05	8.3E+01	8.0E+01	
				7.0E-01	P	7.0E-02	P	V		0.73	1	1.0	Yes	Ethyl Acetate	141-78-6					1.4E+04	9.7E+05	1.5E+02	1.4E+02	
				5.0E-03	P	8.0E-03	P	V		1.32	1	1.0	Yes	Ethyl Acrylate	140-88-5					1.0E+02	3.0E+03	1.7E+01	1.4E+01	
						4.0E+00	P	V		1.43	1	1.0	Yes	Ethyl Chloride (Chloroethane)	75-00-3							8.3E+03	8.3E+03	
				2.0E-01	I			V		0.89	1	1.0	Yes	Ethyl Ether	60-29-7					4.0E+03	2.0E+05		3.9E+03	
						3.0E-01	P	V		1.94	1	1.0	Yes	Ethyl Methacrylate	97-63-2							6.3E+02	6.3E+02	
		8.0E-08	I	1.0E+00	I	4.0E+01	I	V		1.9203	1	1.0	Yes	Ethyl Tertiary Butyl Ether (ETBE)	637-92-3			7.0E+03	7.0E+03	2.0E+04	2.4E+05	8.3E+04	1.5E+04	
				1.0E-05	I					4.78	1	0.8	Yes	Ethyl-p-nitrophenyl Phosphonate	2104-64-5					2.0E-01	1.6E-01		8.9E-02	
1.1E-02	C	2.5E-06	C	5.0E-02	P	1.0E+00	I	V		3.15	1	1.0	Yes	Ethylbenzene	100-41-4	7.1E+02	1.2E+03	2.2E+02	1.5E+02	1.0E+03	1.9E+03	2.1E+03	5.0E+02	7.0E+02
				7.0E-02	P					-0.94	1	1.0	Yes	Ethylene Cyanohydrin	109-78-4					1.4E+03	1.1E+06		1.4E+03	
				9.0E-02	P			V		-2.04	1	1.0	No	Ethylene Diamine	107-15-3					1.8E+03			1.8E+03	
				8.0E-01	A	4.0E-01	C			-1.36	1	1.0	Yes	Ethylene Glycol	107-21-1					1.6E+04	2.3E+07		1.6E+04	
				1.0E-01	I	1.6E+00	I			0.83	1	1.0	Yes	Ethylene Glycol Monobutyl Ether	111-76-2					2.0E+03	1.4E+05		2.0E+03	
3.1E-01	C	3.0E-03	I			3.0E-02	C	V	M	-0.3	1	1.0	Yes	Ethylene Oxide	75-21-8	8.1E+00	1.7E+03	6.8E-02	6.7E-02			6.3E+01	6.3E+01	
4.5E-02	C	1.3E-05	C	8.0E-05	I					-0.66	1	1.0	Yes	Ethylene Thiourea	96-45-7	1.7E+02	1.0E+05		1.7E+02	1.6E+00	1.0E+03		1.6E+00	
6.5E+01	C	1.9E-02	C					V		-0.28	1	1.0	Yes	Ethyleneimine	151-56-4	1.2E-01	2.5E+01	3.0E-02	2.4E-02					
				3.0E+00	I					2.19	1	1.0	Yes	Ethylphthalyl Ethyl Glycolate	84-72-0					6.0E+04	1.5E+06		5.8E+04	
				2.5E-04	I					3.23	1	0.9	Yes	Fenamiphos	22224-92-6					5.0E+00	3.4E+01		4.4E+00	
				2.5E-02	I					5.7	1	0.8	Yes	Fenpropathrin	39515-41-8					5.0E+02	7.3E+01		6.4E+01	
				2.5E-02	I					6.2	1	0.7	No	Fenvalerate	51630-58-1					5.0E+02			5.0E+02	
				1.3E-02	I					2.42	1	1.0	Yes	Fluometuron	2164-17-2					2.6E+02	3.4E+03		2.4E+02	
				4.0E-02	C	1.3E-02	C				1	1.0	Yes	Fluoride	16984-48-8					8.0E+02	1.8E+05		8.0E+02	4.0E+03
				6.0E-02	I	1.3E-02	C				1	1.0	Yes	Fluorine (Soluble Fluoride)	7782-41-4					1.2E+03	2.7E+05		1.2E+03	4.0E+03
				8.0E-02	I					3.16	1	0.9	Yes	Fluridone	59756-60-4					1.6E+03	1.4E+04		1.4E+03	
				4.0E-02	O					3.34	1	0.9	Yes	Flurprimidol	56425-91-3					8.0E+02	4.8E+03		6.9E+02	
				2.0E-03	O					3.7	1	0.9	Yes	Flusilazole	85509-19-9					4.0E+01	1.4E+02		3.1E+01	
				5.0E-01	O					3.7	1	0.9	Yes	Flutolanil	66332-96-5					1.0E+04	3.7E+04		7.9E+03	
				1.0E-02	I					6.81	1	0.6	No	Fluvalinate	69409-94-5					2.0E+02			2.0E+02	
				9.0E-02	O					2.85	1	1.0	Yes	Folpet	133-07-3					1.8E+03	1.9E+04		1.6E+03	
				1.0E-02	O					2.9	1	1.0	Yes	Fomesafen	72178-02-0					2.0E+02	4.8E+03		1.9E+02	
				2.0E-03	I					3.94	1	0.9	Yes	Fonofos	944-22-9					4.0E+01	6.3E+01		2.4E+01	
2.1E-02	C	1.3E-05	I	2.0E-01	I	9.8E-03	A	V		0.35	1	1.0	Yes	Formaldehyde	50-00-0	3.7E+02	2.6E+04	4.3E+01	3.9E+01	4.0E+03	3.2E+05	2.0E+01	2.0E+01	
				9.0E-01	P	3.0E-04	X	V		-0.54	1	1.0	Yes	Formic Acid	64-18-6					1.8E+04	6.4E+06	6.3E-01	6.3E-01	
				2.5E+00	O					-2.4	1	1.0	No	Fosetyl-AL	39148-24-8					5.0E+04			5.0E+04	
														Furans										
				1.0E-03	X			V		4.12	1	1.0	Yes	~Dibenzofuran	132-64-9					2.0E+01	1.3E+01		7.9E+00	
				1.0E-03	I			V		1.34	1	1.0	Yes	~Furan	110-00-9					2.0E+01	4.8E+02		1.9E+01	
				9.0E-01	I	2.0E+00	I	V		0.46	1	1.0	Yes	~Tetrahydrofuran	109-99-9					1.8E+04	1.7E+06	4.2E+03	3.4E+03	
3.8E+00	H									-0.04	1	1.0	Yes	Furazolidone	67-45-8	2.1E+00	1.0E+03		2.0E+00					
				3.0E-03	I	5.0E-02	H	V		0.41	1	1.0	Yes	Furfural	98-01-1					6.0E+01	7.1E+03	1.0E+02	3.8E+01	
1.5E+00	C	4.3E-04	C							1.8	1	1.0	Yes	Furium	531-82-8	5.2E+00	1.9E+02		5.1E+00					
3.0E-02	I	8.6E-06	C							4.38	1	0.9	Yes	Furmecyclox	60568-05-0	2.6E+02	2.0E+02		1.1E+02					
				6.0E-03	O					-4.81	1	1.0	No	Glufosinate, Ammonium	77182-82-2					1.2E+02			1.2E+02	
				1.0E-01	A	8.0E-05	C			-0.33	1	1.0	Yes	Glutaraldehyde	111-30-8					2.0E+03	6.0E+05		2.0E+03	
				4.0E-04	I	1.0E-03	X	V		-0.12	1	1.0	Yes	Glycidaldehyde	765-34-4					8.0E+00	1.8E+03	2.1E+00	1.7E+00	
				1.0E-01	I					-3.4	1	1.0	No	Glyphosate	1071-83-6					2.0E+03			2.0E+03	7.0E+02
				1.0E-02	X			V		-1.63	1	1.0	Yes	Guanidine	113-00-8					2.0E+02	4.2E+05		2.0E+02	
				2.0E-02	P					-3.56	1	1.0	No	Guanidine Chloride	50-01-1					4.0E+02			4.0E+02	
				3.0E-02	X					-8.35	1	1.0	No	Guanidine Nitrate	506-93-4					6.0E+02			6.0E+02	
				5.0E-05	I					4.07	1	0.9	Yes	Haloxyfop, Methyl	69806-40-2					1.0E+00	3.1E+00		7.6E-01	
4.5E+00	I	1.3E-03	I	1.0E-04	A			V		6.1	1	0.8	Yes	Heptachlor	76-44-8	1.7E+00	2.3E-01	4.3E-01	1.4E-01	2.0E+00	2.9E-01		2.6E-01	4.0E-01
9.1E+00	I	2.6E-03	I	1.3E-05	I			V		4.98	1	0.8	Yes	Heptachlor Epoxide	1024-57-3	8.6E-01	7.1E-01	2.2E-01	1.4E-01	2.6E-01	2.4E-01		1.2E-01	2.0E-01
						3.0E-03	X	V		2.29	1	1.0	Yes	Heptanal, n-	111-71-7							6.3E+00	6.3E+00	
				3.0E-04	X	4.0E-01	P	V		4.66	1	1.0	No	Heptane, N-	142-82-5					6.0E+00		8.3E+02	6.0E+00	
				2.0E-03	I			V		6.07	1	0.7	No	Hexabromobenzene	87-82-1					4.0E+01			4.0E+01	
				2.0E-04	I						1	0.0	No	Hexabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-153)	68631-49-2					4.0E+00			4.0E+00	
1.6E+00	I	4.6E-04	I	1.0E-05	P			V		5.73	1	0.9	No	Hexachlorobenzene	118-74-1	4.9E+00		1.2E+00	9.8E-01	2.0E-01			2.0E-01	1.0E+00
7.8E-02	I	2.2E-05	I	1.0E-03	P			V		4.78	1	0.9	Yes	Hexachlorobutadiene	87-68-3	1.0E+02	4.4E+01	2.6E+01	1.4E+01	2.0E+01	9.5E+00		6.5E+00	
6.3E+00	I	1.8E-03	I	8.0E-03	A					3.8	1	0.9	Yes	Hexachlorocyclohexane, Alpha-	319-84-6	1.2E+00	1.8E+00		7.2E-01	1.6E+02	2.5E+02		9.7E+01	
1.8E+00	I	5.3E-04	I							3.78	1	0.9	Yes	Hexachlorocyclohexane, Beta-	319-85-7	4.3E+00	6.1E+00		2.5E+00					
1.1E+00	C	3.1E-04	C	1.0E-05	A					3.72	1	0.9	Yes	Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	7.1E+00	1.0E+01		4.2E+00	2.0E-01	3.1E-01		1.2E-01	2.0E-01
1.8E+00	I	5.1E-04	I							4.14	1	0.9	Yes	Hexachlorocyclohexane, Technical	608-73-1	4.3E+00	6.1E+00		2.5E+00					
				6.0E-03	I	2.0E-04	I	V		5.04	1	0.9	Yes	Hexachlorocyclopentadiene	77-47-4					1.2E+02	4.2E+01	4.2E-01	4.1E-01	5.0E+01
4.0E-02	I	1.1E-05	C	7.0E-04	I	3.0E-02	I	V		4.14	1	1.0	Yes	Hexachloroethane	67-72-1	1.9E+02	1.7E+02	5.1E+01	3.3E+01	1.4E+01	1.4E+01	6.3E+01	6.2E+00	
				3.0E-04	I					7.54	1	0.0	No	Hexachlorophene	70-30-4					6.0E+00			6.0E+00	
8.0E-02	I			4.0E-03	I					0.87	1	1.0	Yes	Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	9.7E+01	1.2E+04		9.7E+01	8.0E+01	1.1E+04		8.0E+01	
						1.0E-05	I	V		3.2	1	1.0	Yes	Hexamethylene Diisocyanate, 1,6-	822-06-0							2.1E-02	2.1E-02	

RMLS TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1				
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	log K_{ow} (unitless)	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
						4.0E-04	C			7.5795	1	0.0	No	Hexamethylene diisocyanate biuret	4035-89-6									
						4.0E-04	C			9.814	1	0.0	No	Hexamethylene diisocyanate isocyanurate	3779-63-3									
				4.0E-04	P					0.28	1	1.0	Yes	Hexamethylphosphoramide	680-31-9					8.0E+00	2.0E+03		8.0E+00	
		2.0E-07	X			6.0E-01	P	V		3.9	1	1.0	Yes	Hexane, Commercial	E5241997			2.8E+03	2.8E+03			1.3E+03	1.3E+03	
						7.0E-01	I	V		3.9	1	1.0	Yes	Hexane, N-	110-54-3							1.5E+03	1.5E+03	
				2.0E+00	P					0.08	1	1.0	Yes	Hexanedioic Acid	124-04-9					4.0E+04	1.1E+07		4.0E+04	
9.5E-03	P			7.0E-02	P	4.0E-04	P	V		2.73	1	1.0	Yes	Hexanol, 1-,2-ethyl- (2-Ethyl-1-hexanol)	104-76-7	8.2E+02	3.2E+03		6.5E+02	1.4E+03	5.9E+03	8.3E-01	8.3E-01	
				5.0E-03	I	3.0E-02	I	V		1.38	1	1.0	Yes	Hexanone, 2-	591-78-6					1.0E+02	2.8E+03	6.3E+01	3.8E+01	
				3.3E-02	I					1.85	1	1.0	Yes	Hexazinone	51235-04-2					6.6E+02	2.4E+04		6.4E+02	
				2.5E-02	I					5.57	1	0.8	Yes	Hexythiazox	78587-05-0					5.0E+02	1.4E+02		1.1E+02	
				1.7E-02	O					2.31	1	1.0	Yes	Hydramethylnon	67485-29-4					3.4E+02	2.9E+04		3.4E+02	
3.0E+00	I	4.9E-03	I			3.0E-05	P	V		-2.07	1	1.0	Yes	Hydrazine	302-01-2	2.6E+00	1.1E+04	1.1E-01	1.1E-01			6.3E-02	6.3E-02	
3.0E+00	I	4.9E-03	I								1	1.0	Yes	Hydrazine Sulfate	10034-93-2	2.6E+00	4.9E+02		2.6E+00					
						2.0E-02	I	V			1	1.0	Yes	Hydrogen Chloride	7647-01-0							4.2E+01	4.2E+01	
				4.0E-02	C	1.4E-02	C	V		0.23	1	1.0	Yes	Hydrogen Fluoride	7664-39-3					8.0E+02	1.8E+05	2.9E+01	2.8E+01	
						2.0E-03	I	V		0.23	1	1.0	Yes	Hydrogen Sulfide	7783-06-4							4.2E+00	4.2E+00	
6.0E-02	P			4.0E-02	P					0.59	1	1.0	Yes	Hydroquinone	123-31-9	1.3E+02	1.2E+04		1.3E+02	8.0E+02	7.9E+04		7.9E+02	
6.1E-02	O			1.1E-01	O					3.82	1	0.9	Yes	Imazalil	35554-44-0	1.3E+02	3.1E+02		9.0E+01	2.2E+03	5.7E+03		1.6E+03	
				2.5E-01	I					1.86	1	1.0	Yes	Imazaquin	81335-37-7					5.0E+03	2.6E+05		4.9E+03	
				2.5E+00	O					1.49	1	1.0	Yes	Imazethapyr	81335-77-5					5.0E+04	7.2E+05		4.7E+04	
				1.0E-02	A					2.49	1	1.0	Yes	Iodine	7553-56-2					2.0E+02	4.6E+04		2.0E+02	
				4.0E-02	I					3	1	0.9	Yes	Iprodione	36734-19-7					8.0E+02	9.1E+03		7.4E+02	
				7.0E-01	P						1	1.0	Yes	Iron	7439-89-6					1.4E+04	3.2E+06		1.4E+04	
				3.0E-01	I			V		0.76	1	1.0	Yes	Isobutyl Alcohol	78-83-1					6.0E+03	3.6E+05		5.9E+03	
9.5E-04	I			2.0E-01	I	2.0E+00	C			1.7	1	1.0	Yes	Isophorone	78-59-1	8.2E+03	1.6E+05		7.8E+03	4.0E+03	8.6E+04		3.8E+03	
				1.5E-02	I			V		5.8	1	0.8	Yes	Isopropalin	33820-53-0					3.0E+02	4.6E+01		4.0E+01	
				2.0E+00	P	2.0E-01	P	V		0.05	1	1.0	Yes	Isopropanol	67-63-0					4.0E+04	6.5E+06	4.2E+02	4.1E+02	
				1.0E-01	I					0.27	1	1.0	Yes	Isopropyl Methyl Phosphonic Acid	1832-54-8					2.0E+03	3.9E+05		2.0E+03	
				5.0E-02	I					3.94	1	0.9	Yes	Isoxaben	82558-50-7					1.0E+03	2.7E+03		7.3E+02	
						3.0E-01	A	V		8	1	0.0	No	JP-7	E1737665							6.3E+02	6.3E+02	
				8.0E-03	O					4.81	1	0.9	Yes	Lactofen	77501-63-4					1.6E+02	2.7E+02		1.0E+02	
				2.0E-04	X					-0.94	1	1.0	Yes	Lactonitrile	78-97-7					4.0E+00	3.2E+03		4.0E+00	
				5.0E-05	P						1	1.0	Yes	Lanthanum	7439-91-0					1.0E+00	2.3E+02		1.0E+00	
				2.1E-05	P						1	0.0	No	Lanthanum Acetate Hydrate	100587-90-4					4.2E-01			4.2E-01	
				1.9E-05	P						1	1.0	Yes	Lanthanum Chloride Heptahydrate	10025-84-0					3.8E-01	8.5E+01		3.7E-01	
				2.8E-05	P						1	1.0	Yes	Lanthanum Chloride, Anhydrous	10099-58-8					5.7E-01	1.3E+02		5.7E-01	
				1.6E-05	P						1	0.9	Yes	Lanthanum Nitrate Hexahydrate	10277-43-7					3.2E-01	7.3E+01		3.2E-01	
														Lead Compounds										
8.5E-03	C	1.2E-05	C								1	0.8	Yes	~Lead Phosphate	7446-27-7	9.2E+02	1.7E+05		9.1E+02					
2.1E-01	C	8.0E-05	C							-0.08	1	1.0	Yes	~Lead acetate	301-04-2	3.7E+01	3.7E+04		3.7E+01					
											1	1.0	Yes	~Lead and Compounds	7439-92-1								1.5E+01	1.5E+01
3.8E-02	C	1.1E-05	C							-4	1	1.0	No	~Lead subacetate	1335-32-6	2.1E+02			2.1E+02					
				1.0E-07	I			V		4.15	1	0.9	Yes	~Tetraethyl Lead	78-00-2					2.0E-03	3.8E-03		1.3E-03	
				5.0E-06	P			V		2.56	1	1.0	Yes	Lewisite	541-25-3					1.0E-01	9.1E-01		9.0E-02	
				7.7E-03	O					3.2	1	0.9	Yes	Linuron	330-55-2					1.5E+02	7.6E+02		1.3E+02	
				2.0E-03	P						1	1.0	Yes	Lithium	7439-93-2					4.0E+01	9.1E+03		4.0E+01	
				5.0E-04	I					3.25	1	1.0	Yes	MCPA	94-74-6					1.0E+01	3.0E+01		7.5E+00	
				4.4E-03	O					2.79	1	0.9	Yes	MCPB	94-81-5					8.8E+01	2.4E+02		6.5E+01	
				1.0E-03	I					3.13	1	1.0	Yes	MCPP	93-65-2					2.0E+01	7.1E+01		1.6E+01	
				2.0E-02	I					2.36	1	1.0	Yes	Malathion	121-75-5					4.0E+02	1.1E+04		3.9E+02	
				1.0E-01	I	7.0E-04	C			1.62	1	1.0	Yes	Maleic Anhydride	108-31-6					2.0E+03	3.8E+04		1.9E+03	
				5.0E-01	I					-0.84	1	1.0	Yes	Maleic Hydrazide	123-33-1					1.0E+04	8.9E+06		1.0E+04	
				1.0E-04	P					-0.6	1	1.0	Yes	Malononitrile	109-77-3					2.0E+00	9.2E+02		2.0E+00	
				3.0E-02	H					1.33	1	0.9	Yes	Mancozeb	8018-01-7					6.0E+02	4.9E+03		5.4E+02	
				5.0E-03	I					0.62	1	1.0	Yes	Maneb	12427-38-2					1.0E+02	3.6E+03		9.8E+01	
				1.4E-01	I	5.0E-05	I				1	1.0		Manganese (Diet)	7439-96-5									
				2.4E-02	G	5.0E-05	I				0.04	1.0	Yes	Manganese (Non-diet)	7439-96-5					4.8E+02	4.4E+03		4.3E+02	
				9.0E-05	H					1.04	1	1.0	Yes	Mephosfolan	950-10-7					1.8E+00	2.5E+02		1.8E+00	
				3.0E-02	I					-2.82	1	1.0	No	Mepiquat Chloride	24307-26-4					6.0E+02			6.0E+02	
1.1E-02	P			4.0E-03	P					2.42	1	1.0	Yes	Mercaptobenzothiazole, 2-	149-30-4	7.1E+02	5.6E+03		6.3E+02	8.0E+01	6.9E+02		7.2E+01	
														Mercury Compounds										
				3.0E-04	I	3.0E-04	G			-0.22	0.07	1.0	Yes	~Mercuric Chloride (and other Mercury salts)	7487-94-7					6.0E+00	9.6E+01		5.7E+00	2.0E+00
						3.0E-04	I	V		0.62	1	1.0	Yes	~Mercury (elemental)	7439-97-6							6.3E-01	6.3E-01	2.0E+00
				1.0E-04	I						1	1.0	Yes	~Methyl Mercury	22967-92-6					2.0E+00	4.6E+02		2.0E+00	
				8.0E-05	I					0.71	1	1.0	Yes	~Phenylmercuric Acetate	62-38-4					1.6E+00	5.7E+02		1.6E+00	
				3.0E-05	I			V		7.67	1	0.3	No	Merphos	150-50-5					6.0E-01			6.0E-01	
				6.0E-02	I					1.65	1	1.0	Yes	Metalaxyl	57837-19-1					1.2E+03	6.4E+04		1.2E+03	
				1.0E-04	I	3.0E-02	P	V		0.68	1	1.0	Yes	Methacrylonitrile	126-98-7					2.0E+00	1.3E+02	6.3E+01	1.9E+00	
				5.0E-05	I					-0.8	1	1.0	Yes	Methamidophos	10265-92-6					1.0E+00	1.0E+03		1.0E+00	
				2.0E+00	I	2.0E+01	I	V		-0.77	1	1.0	Yes	Methanol	67-56-1					4.0E+04	1.8E+07	4.2E+04	2.0E+04	
				1.5E-03	O					2.2	1	1.0	Yes	Methidathion	950-37-8					3.0E+01	8.7E+02		2.9E+01	
				2.5E-02	I					0.6	1	1.0	Yes	Methomyl	16752-77-5					5.0E+02	6.8E+04		5.0E+02	
4.9E-02	C									1.47	1	1.0	Yes	Methoxy-5-nitroaniline, 2-	99-59-2	1.6E+02	5.4E+03		1.5E+02					
				5.0E-03	I					5.08	1	0.8	Yes	Methoxychlor	72-43-5					1.0E+02	5.9E+01		3.7E+01	4.0E+01
				8.0E-03	P	1.0E-03	P	V		0.1	1	1.0	Yes	Methoxyethanol Acetate, 2-	110-49-6					1.6E+02	3.5E+04	2.1E+00	2.1E+00	
				5.0E-03	P	7.0E-03	P	V		-0.77	1	1.0	Yes	Methoxyethanol, 2-	109-86-4					1.0E+02	6.3E+04	1.5E+01	1.3E+01	

RMLS TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1				
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	log K_{ow} (unitless)	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
				1.0E+00	X			V		0.18	1	1.0	Yes	Methyl Acetate	79-20-9					2.0E+04	2.9E+06		2.0E+04	
						2.0E-02	P	V		0.8	1	1.0	Yes	Methyl Acrylate	96-33-3							4.2E+01	4.2E+01	
				6.0E-01	I	5.0E+00	I	V		0.29	1	1.0	Yes	Methyl Ethyl Ketone (2-Butanone)	78-93-3					1.2E+04	1.5E+06	1.0E+04	5.6E+03	
		1.0E-03	X	1.0E-03	P	2.0E-05	X	V		-1.05	1	1.0	Yes	Methyl Hydrazine	60-34-4			5.6E-01	5.6E-01	2.0E+01	1.5E+04	4.2E-02	4.2E-02	
						3.0E+00	I	V		1.31	1	1.0	Yes	Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1							6.3E+03	6.3E+03	
						1.0E-03	C	V		0.79	1	1.0	Yes	Methyl Isocyanate	624-83-9							2.1E+00	2.1E+00	
				1.4E+00	I	7.0E-01	I	V		1.38	1	1.0	Yes	Methyl Methacrylate	80-62-6					2.8E+04	7.7E+05	1.5E+03	1.4E+03	
				2.5E-04	I					2.86	1	1.0	Yes	Methyl Parathion	298-00-0					5.0E+00	4.1E+01		4.5E+00	
				6.0E-02	X					-0.7	1	1.0	Yes	Methyl Phosphonic Acid	993-13-5					1.2E+03	1.2E+06		1.2E+03	
				6.0E-03	H	4.0E-02	H	V		3.44	1	0.8	Yes	Methyl Styrene (Mixed Isomers)	25013-15-4					1.2E+02	4.3E+01	8.3E+01	2.3E+01	
9.9E-02	C	2.8E-05	C							-0.66	1	1.0	Yes	Methyl methanesulfonate	66-27-3	7.9E+01	4.8E+04		7.9E+01					
1.8E-03	C	2.6E-07	C			3.0E+00	I	V		0.94	1	1.0	Yes	Methyl tert-Butyl Ether (MTBE)	1634-04-4	4.3E+03	2.0E+05	2.2E+03	1.4E+03			6.3E+03	6.3E+03	
				3.0E-04	X					-2.06	1	1.0	Yes	Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2					6.0E+00	5.9E+04		6.0E+00	
						3.0E+00	X	V		1.43	1	1.0	Yes	Methyl-2-Pentanol, 4-	108-11-2							6.3E+03	6.3E+03	
9.0E-03	P			2.0E-02	X					1.87	1	1.0	Yes	Methyl-5-Nitroaniline, 2-	99-55-8	8.7E+02	1.4E+04		8.2E+02	4.0E+02	7.3E+03		3.8E+02	
8.3E+00	C	2.4E-03	C							-0.92	1	1.0	Yes	Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	9.4E-01	1.1E+03		9.4E-01					
1.3E-01	C	3.7E-05	C							1.62	1	1.0	Yes	Methylaniline Hydrochloride, 2-	636-21-5	6.0E+01	3.9E+05		6.0E+01					
				1.0E-02	A					-1.18	1	1.0	Yes	Methylarsonic acid	124-58-3					2.0E+02	3.6E+05		2.0E+02	
				2.0E-04	X						1	0.0	No	Methylbenzene,1-4-diamine monohydrochloride, 2-	74612-12-7					4.0E+00			4.0E+00	
1.0E-01	X			3.0E-04	X						1	0.0	No	Methylbenzene-1,4-diamine sulfate, 2-	615-50-9	7.8E+01			7.8E+01	6.0E+00			6.0E+00	
2.2E+01	C	6.3E-03	C						M	6.42	1	0.8	No	Methylcholanthrene, 3-	56-49-5	1.1E-01			1.1E-01					
2.0E-03	I	1.0E-08	I	6.0E-03	I	6.0E-01	I	V	M	1.25	1	1.0	Yes	Methylene Chloride	75-09-2	1.3E+03	3.5E+04	2.0E+04	1.1E+03	1.2E+02	3.7E+03	1.3E+03	1.1E+02	5.0E+00
1.0E-01	P	4.3E-04	C	2.0E-03	P				M	3.91	1	0.9	Yes	Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	2.5E+01	4.3E+01		1.6E+01	4.0E+01	7.5E+01		2.6E+01	
4.6E-02	I	1.3E-05	C							4.37	1	1.0	Yes	Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	1.7E+02	1.2E+02		7.0E+01					
1.6E+00	C	4.6E-04	C			2.0E-02	C			1.59	1	1.0	Yes	Methylenebisbenzenamine, 4,4'-	101-77-9	4.9E+00	1.7E+02		4.7E+00					
						6.0E-04	I			5.22	1	0.9	Yes	Methylenediphenyl Diisocyanate	101-68-8									
				7.0E-02	H			V		3.48	1	1.0	Yes	Methylstyrene, Alpha-	98-83-9					1.4E+03	1.7E+03		7.8E+02	
				1.5E-01	I					3.13	1	1.0	Yes	Metolachlor	51218-45-2					3.0E+03	2.6E+04		2.7E+03	
				2.5E-02	I					1.7	1	1.0	Yes	Metribuzin	21087-64-9					5.0E+02	1.8E+04		4.9E+02	
				2.5E-01	I					2.2	1	1.0	Yes	Metsulfuron-methyl	74223-64-6					5.0E+03	2.4E+05		4.9E+03	
		4.5E-06	X	1.0E-02	X	1.0E-01	P	V		5.65	1	1.0	No	Midrange Aliphatic Hydrocarbon Streams	E1790669			1.2E+02	1.2E+02	2.0E+02		2.1E+02	1.0E+02	
				3.0E+00	P			V		6.1	1	1.0	No	Mineral oils	8012-95-1					6.0E+04			6.0E+04	
1.8E+01	C	5.1E-03	C	2.0E-04	I			V		6.89	1	0.5	No	Mirex	2385-85-5	4.3E-01		1.1E-01	8.8E-02	4.0E+00			4.0E+00	
				2.0E-03	I					3.21	1	1.0	Yes	Molinate	2212-67-1					4.0E+01	1.2E+02		3.0E+01	
				5.0E-03	I	2.0E-03	A				1	1.0	Yes	Molybdenum	7439-98-7					1.0E+02	2.3E+04		1.0E+02	
				1.0E-01	I						1	1.0	Yes	Monochloramine	10599-90-3					2.0E+03	4.6E+05		2.0E+03	4.0E+03(G)
				2.0E-03	P					1.66	1	1.0	Yes	Monomethylaniline	100-61-8					4.0E+01	7.5E+02		3.8E+01	
				2.5E-02	I					2.94	1	1.0	Yes	Myclobutanil	88671-89-0					5.0E+02	4.7E+03		4.5E+02	
				3.0E-04	X					4.04	1	0.9	Yes	N,N'-Diphenyl-1,4-benzenediamine	74-31-7					6.0E+00	8.9E+00		3.6E+00	
				2.0E-03	I			V		1.38	1	1.0	Yes	Naled	300-76-5					4.0E+01	6.8E+03		4.0E+01	
				3.0E-02	X	1.0E-01	P	V			1	0.0	No	Naphtha, High Flash Aromatic (HFAN)	64742-95-6					6.0E+02		2.1E+02	1.5E+02	
1.8E+00	C	0.0E+00	C							2.28	1	1.0	Yes	Naphthylamine, 2-	91-59-8	4.3E+00	3.6E+01		3.9E+00					
				1.2E-01	O					3.36	1	0.9	Yes	Napropamide	15299-99-7					2.4E+03	1.1E+04		2.0E+03	
		2.6E-04	C	1.1E-02	C	1.4E-05	C			-1.38	1	1.0	Yes	Nickel Acetate	373-02-4					2.2E+02	6.8E+05		2.2E+02	
		2.6E-04	C	1.1E-02	C	1.4E-05	C			-2.12	1	1.0	Yes	Nickel Carbonate	3333-67-3					2.2E+02	1.4E+06		2.2E+02	
		2.6E-04	C	1.1E-02	C	1.4E-05	C	V			1	0.0	No	Nickel Carbonyl	13463-39-3			2.2E+00	2.2E+00	2.2E+02		2.9E-02	2.9E-02	
		2.6E-04	C	1.1E-02	C	1.4E-05	C				0.04	1.0	Yes	Nickel Hydroxide	12054-48-7					2.2E+02	2.0E+03		2.0E+02	
		2.6E-04	C	1.1E-02	C	2.0E-05	C				0.04	1.0	Yes	Nickel Oxide	1313-99-1					2.2E+02	2.0E+03		2.0E+02	
		2.4E-04	I	1.1E-02	C	1.4E-05	C				0.04	0.0	Yes	Nickel Refinery Dust	E715532					2.2E+02	1.0E+04		2.2E+02	
		2.6E-04	C	2.0E-02	I	9.0E-05	A				0.04	1.0	Yes	Nickel Soluble Salts	7440-02-0					4.0E+02	1.8E+04		3.9E+02	
1.7E+00	C	4.8E-04	I	1.1E-02	C	1.4E-05	C				0.04	1.0	Yes	Nickel Subsulfide	12035-72-2	4.6E+00	1.7E+02		4.5E+00	2.2E+02	1.0E+04		2.2E+02	
9.1E-01	C	2.6E-04	C	1.1E-02	C	1.4E-05	C				1	0.0	No	Nickelocene	1271-28-9	8.6E+00			8.6E+00	2.2E+02			2.2E+02	
				1.6E+00	I						1	1.0	Yes	Nitrate (measured as nitrogen)	14797-55-8					3.2E+04	7.3E+06		3.2E+04	1.0E+04
											1	0.0	Yes	Nitrate + Nitrite (measured as nitrogen)	E701177									1.0E+04
				1.0E-01	I						1	1.0	Yes	Nitrite (measured as nitrogen)	14797-65-0					2.0E+03	4.6E+05		2.0E+03	1.0E+03
				1.0E-02	X	5.0E-05	X			1.85	1	1.0	Yes	Nitroaniline, 2-	88-74-4					2.0E+02	3.4E+03		1.9E+02	
2.0E-02	P			4.0E-03	P	6.0E-03	P			1.39	1	1.0	Yes	Nitroaniline, 4-	100-01-6	3.9E+02	1.2E+04		3.8E+02	8.0E+01	2.8E+03		7.8E+01	
		4.0E-05	I	2.0E-03	I	9.0E-03	I	V		1.85	1	1.0	Yes	Nitrobenzene	98-95-3			1.4E+01	1.4E+01	4.0E+01	6.2E+02	1.9E+01	1.3E+01	
				3.0E+03	P					-4.56	1	1.0	No	Nitrocellulose	9004-70-0					6.0E+07			6.0E+07	
				7.0E-02	H					-0.47	1	1.0	Yes	Nitrofurantoin	67-20-9					1.4E+03	1.6E+06		1.4E+03	
1.3E+00	C	3.7E-04	C							0.23	1	1.0	Yes	Nitrofurazone	59-87-0	6.0E+00	1.7E+03		6.0E+00					
1.7E-02	P			1.0E-04	P					1.62	1	1.0	Yes	Nitroglycerin	55-63-0	4.6E+02	1.8E+04		4.5E+02	2.0E+00	8.7E+01		2.0E+00	
				1.0E-01	I					-0.89	1	1.0	Yes	Nitroguanidine	556-88-7					2.0E+03	1.8E+06		2.0E+03	
		8.8E-06	P			5.0E-03	P	V		-0.35	1	1.0	Yes	Nitromethane	75-52-5			6.4E+01	6.4E+01			1.0E+01	1.0E+01	
		5.8E-04	X			2.0E-02	I	V		0.93	1	1.0	Yes	Nitropropane, 2-	79-46-9			9.7E-01	9.7E-01			4.2E+01	4.2E+01	
2.7E+01	C	7.7E-03	C						M	0.23	1	1.0	Yes	Nitroso-N-ethylurea, N-	759-73-9	9.3E-02	1.5E+01		9.2E-02					
1.2E+02	C	3.4E-02	C						M	-0.03	1	1.0	Yes	Nitroso-N-methylurea, N-	684-93-5	2.1E-02	4.6E+00		2.1E-02					
5.4E+00	I	1.6E-03	I					V		2.63	1	1.0	Yes	Nitroso-di-N-butylamine, N-	924-16-3	1.4E+00	7.9E+00	3.5E-01	2.7E-01					
7.0E+00	I	2.0E-03	C							1.36	1	1.0	Yes	Nitroso-di-N-propylamine, N-	621-64-7	1.1E+00	3.5E+01		1.1E+00					
2.8E+00	I	8.0E-04	C							-1.28	1	1.0	Yes	Nitrosodiethanolamine, N-	1116-54-7	2.8E+00	5.1E+03		2.8E+00					
1.5E+02	I	4.3E-02	I						M	0.48	1	1.0	Yes	Nitrosodiethylamine, N-	55-18-5	1.7E-02	1.7E+00		1.7E-02					
5.1E+01	I	1.4E-02	I	8.0E-06	P	4.0E-05	X	V	M	-0.57	1	1.0	Yes	Nitrosodimethylamine, N-	62-75-9	4.9E-02	2.0E+01	1.4E-02	1.1E-02	1.6E-01	7.4E+01	8.3E-02	5.5E-02	
4.9E-03	I	2.6E-06	C							3.13	1	1.0	Yes	Nitrosodiphenylamine, N-	86-30-6	1.6E+03	5.2E+03		1.2E+03					
2.2E+01	I	6.3E-03	C					V		0.04	1	1.0	Yes	Nitrosomethylethylamine, N-	10595-95-6	3.5E-01	6.4E+01	8.9E-02	7.1E-02					
6.7E+00	C	1.9E-03	C							-0.44	1	1.0	Yes	Nitrosomorpholine [N-]	59-89-2	1.2E+00	5.3E+02		1.2E+00					
9.4E+00	C	2.7E-03	C							0.36	1	1.0	Yes	Nitrosopiperidine [N-]	100-75-4	8.3E-01	1.1E+02		8.2E-01					

RMLS TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1				
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	log K_{ow} (unitless)	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
2.1E+00	I	6.1E-04	I							-0.19	1	1.0	Yes	Nitrosopyrrolidine, N-	930-55-2	3.7E+00	1.0E+03		3.7E+00					
				1.0E-04	X					2.45	1	1.0	Yes	Nitrotoluene, m-	99-08-1					2.0E+00	1.4E+01		1.7E+00	
2.2E-01	P			9.0E-04	P			V		2.3	1	1.0	Yes	Nitrotoluene, o-	88-72-2	3.5E+01	2.8E+02		3.1E+01	1.8E+01	1.5E+02		1.6E+01	
1.6E-02	P			4.0E-03	P					2.37	1	1.0	Yes	Nitrotoluene, p-	99-99-0	4.9E+02	3.4E+03		4.3E+02	8.0E+01	6.2E+02		7.1E+01	
				3.0E-04	X	2.0E-02	P	V		5.65	1	1.0	No	Nonane, n-	111-84-2					6.0E+00		4.2E+01	5.3E+00	
				1.5E-03	O					2.3	1	1.0	Yes	Norflurazon	27314-13-2					3.0E+01	7.5E+02		2.9E+01	
				3.0E-03	I					8.71	1	0.3	No	Octabromodiphenyl Ether	32536-52-0					6.0E+01			6.0E+01	
				5.0E-02	I					0.16	1	1.0	Yes	Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0					1.0E+03	6.3E+05		1.0E+03	
				2.0E-03	H					-1.01	1	1.0	Yes	Octamethylpyrophosphoramide	152-16-9					4.0E+01	1.4E+05		4.0E+01	
7.8E-03	O			1.9E-01	O					3.73	1	0.9	Yes	Oryzalin	19044-88-3	1.0E+03	3.8E+03		7.9E+02	3.8E+03	1.6E+04		3.1E+03	
				5.0E-03	I					4.8	1	0.8	Yes	Oxadiazon	19666-30-9					1.0E+02	9.0E+01		4.7E+01	
				2.5E-02	I					-0.47	1	1.0	Yes	Oxamyl	23135-22-0					5.0E+02	5.1E+05		5.0E+02	2.0E+02
7.3E-02	O			4.0E-02	O					4.73	1	0.8	Yes	Oxyfluorfen	42874-03-3	1.1E+02	1.1E+02		5.4E+01	8.0E+02	8.9E+02		4.2E+02	
				1.3E-02	I					3.2	1	0.9	Yes	Paclobutrazol	76738-62-0					2.6E+02	1.7E+03		2.3E+02	
				4.5E-03	I					-4.5	1	1.0	No	Paraquat Dichloride	1910-42-5					9.0E+01			9.0E+01	
				6.0E-03	H					3.83	1	0.9	Yes	Parathion	56-38-2					1.2E+02	3.0E+02		8.6E+01	
				5.0E-02	H			V		3.83	1	1.0	Yes	Pebulate	1114-71-2					1.0E+03	1.3E+03		5.6E+02	
				3.0E-01	O					5.2	1	0.9	Yes	Pendimethalin	40487-42-1					6.0E+03	1.8E+03		1.4E+03	
				2.0E-03	I			V		6.84	1	0.6	No	Pentabromodiphenyl Ether	32534-81-9					4.0E+01			4.0E+01	
				1.0E-04	I					7.66	1	0.6	No	Pentabromodiphenyl ether, 2,2',4,4',5- (BDE-99)	60348-60-9					2.0E+00			2.0E+00	
				8.0E-04	I			V		5.17	1	0.9	Yes	Pentachlorobenzene	608-93-5					1.6E+01	3.9E+00		3.2E+00	
9.0E-02	P							V		3.22	1	1.0	Yes	Pentachloroethane	76-01-7	8.7E+01	2.5E+02		6.5E+01					
2.6E-01	H			3.0E-03	I			V		4.64	1	0.9	Yes	Pentachloronitrobenzene	82-68-8	3.0E+01	2.0E+01		1.2E+01	6.0E+01	4.4E+01		2.6E+01	
4.0E-01	I	5.1E-06	C	5.0E-03	I					5.12	1	0.9	Yes	Pentachlorophenol	87-86-5	1.9E+01	5.2E+00		4.1E+00	1.0E+02	2.9E+01		2.3E+01	1.0E+00
4.3E-03	X			9.0E-03	P					2.38	1	1.0	Yes	Pentaerythritol tetranitrate (PETN)	78-11-5	1.8E+03	4.0E+04		1.7E+03	1.8E+02	4.3E+03		1.7E+02	
				1.0E-04	X					-0.77	1	1.0	Yes	Pentamethylphosphoramide (PMPA)	10159-46-3					2.0E+00	2.3E+03		2.0E+00	
						1.0E+00	P	V		3.39	1	1.0	Yes	Pentane, n-	109-66-0							2.1E+03	2.1E+03	
														Per- and Polyfluoroalkyl Substances (PFAS)										
				3.0E-06	D						1	0.0	No	~Ammonium perfluoro-2-methyl-3-oxahexanoate	62037-80-3					6.0E-02			6.0E-02	
				3.0E-06	D			V			1	0.0	No	~Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6					6.0E-02			6.0E-02	
				3.0E-04	P						1	0.0	No	~Perfluorobutanesulfonate	45187-15-3					6.0E+00			6.0E+00	
				3.0E-04	P					-0.34	1	1.0	Yes	~Perfluorobutanesulfonic acid (PFBS)	375-73-5					6.0E+00	8.4E+03		6.0E+00	
				2.0E-05	A					2.2	1	1.0	Yes	~Perfluorohexanesulfonate	108427-53-8					4.0E-01	2.2E+01		3.9E-01	
				2.0E-05	A					2.2	1	1.0	Yes	~Perfluorohexanesulfonic acid (PFHxS)	355-46-4					4.0E-01	2.2E+01		3.9E-01	
				3.0E-06	A					2.57	1	1.0	Yes	~Perfluorononanoate	72007-68-2					6.0E-02	2.8E+00		5.9E-02	
				3.0E-06	A					2.57	1	1.0	Yes	~Perfluorononanoic acid (PFNA)	375-95-1					6.0E-02	2.8E+00		5.9E-02	
				2.0E-06	A						1	0.0	No	~Perfluorooctanesulfonate	45298-90-6					4.0E-02			4.0E-02	
				2.0E-06	A					-1.08	1	1.0	No	~Perfluorooctanesulfonic acid (PFOS)	1763-23-1					4.0E-02			4.0E-02	
7.0E-02	D			3.0E-06	A						1	0.0	No	~Perfluorooctanoate	45285-51-6	1.1E+02			1.1E+02	6.0E-02			6.0E-02	
7.0E-02	D			3.0E-06	A						1	0.0	No	~Perfluorooctanoic acid (PFOA)	335-67-1	1.1E+02			1.1E+02	6.0E-02			6.0E-02	
				3.0E-04	P					-1.8	1	1.0	Yes	~Potassium perfluorobutanesulfonate	29420-49-3					6.0E+00	1.0E+05		6.0E+00	
				2.0E-06	A					-1.08	1	1.0	No	~Potassium perfluorooctanesulfonate	2795-39-3					4.0E-02			4.0E-02	
														Perchlorates										
				7.0E-04	I						1	1.0	Yes	~Ammonium Perchlorate	7790-98-9					1.4E+01	3.2E+03		1.4E+01	
				7.0E-04	I						1	1.0	Yes	~Lithium Perchlorate	7791-03-9					1.4E+01	3.2E+03		1.4E+01	
				7.0E-04	I						1	1.0	Yes	~Perchlorate and Perchlorate Salts	14797-73-0					1.4E+01	3.2E+03		1.4E+01	1.5E+01(G)
				7.0E-04	I						1	1.0	Yes	~Potassium Perchlorate	7778-74-7					1.4E+01	1.6E+03		1.4E+01	
				7.0E-04	I						1	1.0	Yes	~Sodium Perchlorate	7601-89-0					1.4E+01	3.2E+03		1.4E+01	
				5.0E-02	I					6.5	1	0.6	No	Permethrin	52645-53-1					1.0E+03			1.0E+03	
2.2E-03	C	6.3E-07	C							1.58	1	1.0	Yes	Phenacetin	62-44-2	3.5E+03	1.1E+05		3.4E+03					
				2.4E-01	O					3.59	1	0.9	Yes	Phenmedipham	13684-63-4					4.8E+03	1.8E+04		3.8E+03	
				3.0E-01	I	2.0E-01	C			1.46	1	1.0	Yes	Phenol	108-95-2					6.0E+03	1.4E+05		5.8E+03	
				4.0E-03	I					1.52	1	1.0	Yes	Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1					8.0E+01	3.6E+03		7.8E+01	
				5.0E-04	X					4.15	1	1.0	Yes	Phenothiazine	92-84-2					1.0E+01	7.6E+00		4.3E+00	
				2.0E-04	X			V		3.28	1	1.0	Yes	Phenyl Isothiocyanate	103-72-0					4.0E+00	7.6E+00		2.6E+00	
				6.0E-03	I					-0.33	1	1.0	Yes	Phenylenediamine, m-	108-45-2					1.2E+02	4.8E+04		1.2E+02	
1.2E-01	P			4.0E-03	P					0.15	1	1.0	Yes	Phenylenediamine, o-	95-54-5	6.5E+01	1.1E+04		6.5E+01	8.0E+01	1.5E+04		8.0E+01	
				1.0E-03	X					-0.3	1	1.0	Yes	Phenylenediamine, p-	106-50-3					2.0E+01	7.6E+03		2.0E+01	
1.9E-03	H									3.09	1	1.0	Yes	Phenylphenol, 2-	90-43-7	4.0E+03	1.2E+04		3.0E+03					
				2.0E-04	H					3.56	1	0.9	Yes	Phorate	298-02-2					4.0E+00	1.2E+01		3.0E+00	
						3.0E-04	I	V		-0.71	1	1.0	Yes	Phosgene	75-44-5							6.3E-01	6.3E-01	
				2.0E-02	I					2.78	1	1.0	Yes	Phosmet	732-11-6					4.0E+02	5.3E+03		3.7E+02	
				3.0E-04	I	3.0E-04	I	V		-0.27	1	1.0	Yes	Phosphine	7803-51-2					6.0E+00	1.4E+03	6.3E-01	5.7E-01	
						1.0E-02	I				1	1.0	Yes	Phosphoric Acid	7664-38-2									
				2.0E-05	I			V		3.08	1	1.0	Yes	Phosphorus, White	7723-14-0					4.0E-01	9.1E+01		4.0E-01	
														Phthalates										
1.4E-02	I	2.4E-06	C	2.0E-02	I					7.6	1	0.8	No	~Bis(2-ethylhexyl)phthalate	117-81-7	5.6E+02			5.6E+02	4.0E+02			4.0E+02	6.0E+00
1.9E-03	P			2.0E-01	I					4.73	1	0.9	Yes	~Butyl Benzyl Phthalate	85-68-7	4.1E+03	2.7E+03		1.6E+03	4.0E+03	2.9E+03		1.7E+03	
				1.0E+00	I					4.15	1	0.9	Yes	~Butylphthalyl Butylglycolate	85-70-1					2.0E+04	4.1E+04		1.3E+04	
				1.0E-01	I					4.5	1	0.9	Yes	~Dibutyl Phthalate	84-74-2					2.0E+03	1.6E+03		9.0E+02	
				8.0E-01	I					2.42	1	1.0	Yes	~Diethyl Phthalate	84-66-2					1.6E+04	2.0E+05		1.5E+04	
				1.0E-01	I			V		2.25	1	1.0	Yes	~Dimethylterephthalate	120-61-6					2.0E+03	2.7E+04		1.9E+03	
				1.0E-02	P					8.1	1	0.0	No	~Octyl Phthalate, di-N-	117-84-0					2.0E+02			2.0E+02	
				5.0E-01	X					2	1	1.0	Yes	~Phthalic Acid, p-	100-21-0					1.0E+04	1.6E+05		9.4E+03	
				2.0E+00	I	2.0E-02	C			1.6	1	1.0	Yes	~Phthalic Anhydride	85-44-9					4.0E+04	1.1E+06		3.9E+04	
				7.0E-02	I					1.9	1	1.0	Yes	Picloram	1918-02-1					1.4E+03	4.3E+04		1.4E+03	5.0E+02

TMS TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1				
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	log K_{ow} (unitless)	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
				1.0E-04	X					0.93	1	1.0	Yes	Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3					2.0E+00	2.1E+02		2.0E+00	
				2.0E-03	X					1.44	1	1.0	Yes	Picric Acid (2,4,6-Trinitrophenol)	88-89-1					4.0E+01	2.7E+03		4.0E+01	
				7.3E-04	O					4.2	1	0.9	Yes	Pirimiphos, Methyl	29232-93-7					1.5E+01	2.3E+01		8.9E+00	
3.0E+01	C	8.6E-03	C	7.0E-06	H						1	0.0	No	Polybrominated Biphenyls	36355-01-8	2.6E-01			2.6E-01	1.4E-01			1.4E-01	
														Polychlorinated Biphenyls (PCBs)										
7.0E-02	G	2.0E-05	G	7.0E-05	I			V		5.69	1	0.9	No	~Aroclor 1016	12674-11-2	1.1E+02		2.8E+01	2.2E+01	1.4E+00			1.4E+00	
2.0E+00	G	5.7E-04	G					V		4.65	1	1.0	Yes	~Aroclor 1221	11104-28-2	3.9E+00	1.2E+00	9.8E-01	4.7E-01					
2.0E+00	G	5.7E-04	G					V		4.4	1	1.0	Yes	~Aroclor 1232	11141-16-5	3.9E+00	1.2E+00	9.8E-01	4.7E-01					
2.0E+00	G	5.7E-04	G					V		6.34	1	0.7	No	~Aroclor 1242	53469-21-9	3.9E+00		9.8E-01	7.8E-01					
2.0E+00	G	5.7E-04	G					V		6.2	1	0.7	No	~Aroclor 1248	12672-29-6	3.9E+00		9.8E-01	7.8E-01					
2.0E+00	G	5.7E-04	G	2.0E-05	I			V		6.5	1	0.5	No	~Aroclor 1254	11097-69-1	3.9E+00		9.8E-01	7.8E-01	4.0E-01			4.0E-01	
2.0E+00	G	5.7E-04	G					V		7.55	1	0.0	No	~Aroclor 1260	11096-82-5	3.9E+00		9.8E-01	7.8E-01					
				6.0E-04	X			V		6.34	1	0.0	No	~Aroclor 5460	11126-42-4					1.2E+01			1.2E+01	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V		8.27	1	0.0	No	~Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V		7.5	1	0.0	No	~Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167)	52663-72-6	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V		7.6	1	0.0	No	~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V		7.6	1	0.0	No	~Hexachlorobiphenyl, 2,3,3',4,4',5- (PCB 156)	38380-08-4	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01	
3.9E+03	W	1.1E+00	W	2.3E-08	W	1.3E-06	W	V		7.41	1	0.1	No	~Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	2.0E-03		4.9E-04	4.0E-04	4.7E-04		2.8E-03	4.0E-04	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V		6.98	1	0.4	No	~Pentachlorobiphenyl, 2',3,4,4',5- (PCB 123)	65510-44-3	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V		7.12	1	0.3	No	~Pentachlorobiphenyl, 2,3',4,4',5- (PCB 118)	31508-00-6	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V		6.79	1	0.5	No	~Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V		6.98	1	0.4	No	~Pentachlorobiphenyl, 2,3,4,4',5- (PCB 114)	74472-37-0	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01	
1.3E+04	W	3.8E+00	W	7.0E-09	W	4.0E-07	W	V		6.98	1	0.4	No	~Pentachlorobiphenyl, 3,3',4,4',5- (PCB 126)	57465-28-8	6.0E-04		1.5E-04	1.2E-04	1.4E-04		8.3E-04	1.2E-04	
2.0E+00	I	5.7E-04	I					V		7.1	1	0.7		~Polychlorinated Biphenyls (high risk)	1336-36-3									5.0E-01
4.0E-01	I	1.0E-04	I					V		7.1	1	0.7	No	~Polychlorinated Biphenyls (low risk)	1336-36-3	1.9E+01		5.6E+00	4.4E+00					5.0E-01
7.0E-02	I	2.0E-05	I					V		7.1	1	0.7		~Polychlorinated Biphenyls (lowest risk)	1336-36-3									5.0E-01
1.3E+01	W	3.8E-03	W	7.0E-06	W	4.0E-04	W			6.63	1	0.6	No	~Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	6.0E-01			6.0E-01	1.4E-01			1.4E-01	
3.9E+01	W	1.1E-02	W	2.3E-06	W	1.3E-04	W	V		6.34	1	0.7	No	~Tetrachlorobiphenyl, 3,4,4',5- (PCB 81)	70362-50-4	2.0E-01		4.9E-02	4.0E-02	4.7E-02		2.8E-01	4.0E-02	
						6.0E-04	I			10.46	1	0.0	No	Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9									
														Polynuclear Aromatic Hydrocarbons (PAHs)										
				6.0E-02	I			V		3.92	1	1.0	Yes	~Acenaphthene	83-32-9					1.2E+03	9.6E+02		5.3E+02	
				3.0E-01	I			V		4.45	1	1.0	Yes	~Anthracene	120-12-7					6.0E+03	2.5E+03		1.8E+03	
1.0E-01	E	6.0E-05	E					V	M	5.76	1	1.0	No	~Benz[a]anthracene	56-55-3	2.5E+01		3.4E+00	3.0E+00					
				9.0E-05	X	2.0E-06	X			6.44	1	0.9	No	~Benzo(e)pyrene	192-97-2					1.8E+00			1.8E+00	
1.2E+00	C	1.1E-04	C							6.11	1	0.9	No	~Benzo(j)fluoranthene	205-82-3	6.5E+00			6.5E+00					
1.0E+00	I	6.0E-04	I	3.0E-04	I	2.0E-06	I		M	6.13	1	1.0	No	~Benzo[a]pyrene	50-32-8	2.5E+00			2.5E+00	6.0E+00			6.0E+00	2.0E-01
1.0E-01	E	6.0E-05	E						M	5.78	1	1.0	No	~Benzo[b]fluoranthene	205-99-2	2.5E+01			2.5E+01					
1.0E-02	E	6.0E-06	E						M	6.11	1	0.9	No	~Benzo[k]fluoranthene	207-08-9	2.5E+02			2.5E+02					
				8.0E-02	I			V		3.9	1	1.0	Yes	~Chloronaphthalene, Beta-	91-58-7					1.6E+03	1.4E+03		7.5E+02	
1.0E-03	E	6.0E-07	E						M	5.81	1	1.0	No	~Chrysene	218-01-9	2.5E+03			2.5E+03					
1.0E+00	E	6.0E-04	E						M	6.75	1	0.6	No	~Dibenz[a,h]anthracene	53-70-3	2.5E+00			2.5E+00					
1.2E+01	C	1.1E-03	C							7.71	1	0.3	No	~Dibenzo(a,e)pyrene	192-65-4	6.5E-01			6.5E-01					
2.5E+02	C	7.1E-02	C						M	5.8	1	0.9	No	~Dimethylbenz(a)anthracene, 7,12-	57-97-6	1.0E-02			1.0E-02					
				4.0E-02	I					5.16	1	1.0	No	~Fluoranthene	206-44-0					8.0E+02			8.0E+02	
				4.0E-02	I			V		4.18	1	1.0	Yes	~Fluorene	86-73-7					8.0E+02	4.6E+02		2.9E+02	
1.0E-01	E	6.0E-05	E						M	6.7	1	0.6	No	~Indeno[1,2,3-cd]pyrene	193-39-5	2.5E+01			2.5E+01					
2.9E-02	P			7.0E-02	A			V		3.87	1	1.0	Yes	~Methylnaphthalene, 1-	90-12-0	2.7E+02	2.0E+02		1.1E+02	1.4E+03	1.1E+03		6.2E+02	
				4.0E-03	I			V		3.86	1	1.0	Yes	~Methylnaphthalene, 2-	91-57-6					8.0E+01	6.5E+01		3.6E+01	
1.2E-01	C	3.4E-05	C	2.0E-02	I	3.0E-03	I	V		3.3	1	1.0	Yes	~Naphthalene	91-20-3	6.5E+01	1.0E+02	1.7E+01	1.2E+01	4.0E+02	7.0E+02	6.3E+00	6.1E+00	
1.2E+00	C	1.1E-04	C							4.75	1	0.9	Yes	~Nitropyrene, 4-	57835-92-4	6.5E+00	2.7E+00		1.9E+00					
				3.0E-02	I			V		4.88	1	1.0	Yes	~Pyrene	129-00-0					6.0E+02	1.5E+02		1.2E+02	
1.5E-01	I			9.0E-03	I					4.1	1	0.9	Yes	Prochloraz	67747-09-5	5.2E+01	1.4E+02		3.8E+01	1.8E+02	5.1E+02		1.3E+02	
				6.0E-03	H			V		5.58	1	0.8	Yes	Profluralin	26399-36-0					1.2E+02	3.3E+01		2.6E+01	
				1.5E-02	I					2.99	1	1.0	Yes	Prometon	1610-18-0					3.0E+02	1.6E+03		2.5E+02	
				4.0E-02	O					3.51	1	0.9	Yes	Prometryn	7287-19-6					8.0E+02	2.3E+03		6.0E+02	
				7.5E-02	I					3.43	1	0.9	Yes	Pronamide	23950-58-5					1.5E+03	5.5E+03		1.2E+03	
				1.3E-02	I					2.18	1	1.0	Yes	Propachlor	1918-16-7					2.6E+02	4.3E+03		2.5E+02	
				5.0E-03	I					3.07	1	1.0	Yes	Propanil	709-98-8					1.0E+02	4.4E+02		8.2E+01	
1.9E-01	O			4.0E-02	O					5	1	0.8	Yes	Propargite	2312-35-8	4.1E+01	2.5E+01		1.6E+01	8.0E+02	5.5E+02		3.3E+02	
				2.0E-03	I			V		-0.38	1	1.0	Yes	Propargyl Alcohol	107-19-7					4.0E+01	1.2E+04		4.0E+01	
				2.0E-02	I					2.93	1	1.0	Yes	Propazine	139-40-2					4.0E+02	2.4E+03		3.4E+02	
				2.0E-02	I					2.6	1	1.0	Yes	Propham	122-42-9					4.0E+02	2.8E+03		3.5E+02	
				1.0E-01	O					3.72	1	0.9	Yes	Propiconazole	60207-90-1					2.0E+03	8.2E+03		1.6E+03	
						8.0E-03	I	V		0.59	1	1.0	Yes	Propionaldehyde	123-38-6							1.7E+01	1.7E+01	
				1.0E-01	X	1.0E+00	X	V		3.69	1	1.0	Yes	Propyl benzene	103-65-1					2.0E+03	1.8E+03	2.1E+03	6.6E+02	
						3.0E+00	C	V		1.77	1	1.0	Yes	Propylene	115-07-1							6.3E+03	6.3E+03	
				2.0E+01	P					-0.92	1	1.0	Yes	Propylene Glycol	57-55-6					4.0E+05	3.2E+08		4.0E+05	
						2.7E-04	A			1.83	1	1.0	Yes	Propylene Glycol Dinitrate	6423-43-4									
				7.0E-01	H	2.0E+00	I	V		-0.49	1	1.0	Yes	Propylene Glycol Monomethyl Ether	107-98-2					1.4E+04	6.3E+06	4.2E+03	3.2E+03	
2.4E-01	I	3.7E-06	I			3.0E-02	I	V		0.03	1	1.0	Yes	Propylene Oxide	75-56-9	3.2E+01	4.7E+03	1.5E+02	2.7E+01			6.3E+01	6.3E+01	
				1.0E-03	I			V		0.65	1	1.0	Yes	Pyridine	110-86-1					2.0E+01	1.5E+03		2.0E+01	
				5.0E-04	I					4.44	1	0.9	Yes	Quinalphos	13593-03-8					1.0E+01	1.0E+01		5.1E+00	
3.0E+00	I									2.03	1	1.0	Yes	Quinoline	91-22-5	2.6E+00	2.9E+01		2.4E+00					
				9.0E-03	I					4.28	1	0.9	Yes	Quizalofop-ethyl	76578-14-8					1.8E+02	3.8E+02		1.2E+02	
						3.0E+04	A				1	0.0		Refractory Ceramic Fibers (units in fibers)	E715557									
				3.0E-02	I					6.14	1	0.7	Yes	Resmethrin	10453-86-8					6.0E+02	7.6E+01		6.7E+01	

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1				
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	log K_{ow} (unitless)	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
				5.0E-02	H			V		4.88	1	0.8	Yes	Ronnel	299-84-3					1.0E+03	6.8E+02		4.1E+02	
				4.0E-03	I					4.1	1	0.9	Yes	Rotenone	83-79-4					8.0E+01	2.6E+02		6.1E+01	
2.2E-01	C	6.3E-05	C						M	3.45	1	1.0	Yes	Safrole	94-59-7	1.1E+01	6.0E+01		9.6E+00					
				5.0E-03	I						1	1.0	Yes	Selenious Acid	7783-00-8					1.0E+02	2.3E+04		1.0E+02	
				5.0E-03	I	2.0E-02	C				1	1.0	Yes	Selenium	7782-49-2					1.0E+02	2.3E+04		1.0E+02	5.0E+01
				5.0E-03	C	2.0E-02	C				1	1.0	Yes	Selenium Sulfide	7446-34-6					1.0E+02	2.3E+04		1.0E+02	
				1.4E-01	O					4.38	1	0.9	Yes	Sethoxydim	74051-80-2					2.8E+03	3.8E+03		1.6E+03	
						3.0E-03	C				1	1.0	Yes	Silica (crystalline, respirable)	7631-86-9									
				5.0E-03	I						0.04	1.0	Yes	Silver	7440-22-4					1.0E+02	1.5E+03		9.4E+01	
1.2E-01	H			5.0E-03	I					2.18	1	1.0	Yes	Simazine	122-34-9	6.5E+01	9.3E+02		6.1E+01	1.0E+02	1.6E+03		9.4E+01	4.0E+00
				1.3E-02	I					0.37	1	1.0	Yes	Sodium Acifluorfen	62476-59-9					2.6E+02	2.1E+05		2.6E+02	
				4.0E-03	I						1	1.0	Yes	Sodium Azide	26628-22-8					8.0E+01	1.8E+04		8.0E+01	
2.7E-01	H			3.0E-02	I					-1.43	1	1.0	Yes	Sodium Diethyldithiocarbamate	148-18-5	2.9E+01	8.5E+04		2.9E+01	6.0E+02	1.9E+06		6.0E+02	
				5.0E-02	A	1.3E-02	C				1	1.0	Yes	Sodium Fluoride	7681-49-4					1.0E+03	2.3E+05		1.0E+03	4.0E+03
				2.0E-05	I					-3.78	1	1.0	No	Sodium Fluoroacetate	62-74-8					4.0E-01			4.0E-01	
				1.0E-03	H						1	1.0	Yes	Sodium Metavanadate	13718-26-8					2.0E+01	4.6E+03		2.0E+01	
				8.0E-04	P						1	1.0	Yes	Sodium Tungstate	13472-45-2					1.6E+01	3.6E+03		1.6E+01	
				8.0E-04	P						1	1.0	Yes	Sodium Tungstate Dihydrate	10213-10-2					1.6E+01	3.6E+03		1.6E+01	
2.4E-02	H			3.0E-02	I					3.53	1	0.9	Yes	Stirofos (Tetrachlorovinphos)	961-11-5	3.2E+02	1.9E+03		2.8E+02	6.0E+02	3.8E+03		5.2E+02	
				6.0E-01	I						1	1.0	Yes	Strontium, Stable	7440-24-6					1.2E+04	2.7E+06		1.2E+04	
				3.0E-04	I					1.93	1	1.0	Yes	Strychnine	57-24-9					6.0E+00	3.2E+02		5.9E+00	
				2.0E-01	I	1.0E+00	I	V		2.95	1	1.0	Yes	Styrene	100-42-5					4.0E+03	1.0E+04	2.1E+03	1.2E+03	1.0E+02
				3.0E-03	P					2.76	1	1.0	Yes	Styrene-Acrylonitrile (SAN) Trimer (THNA isomer)	57964-39-3					6.0E+01	2.4E+02		4.8E+01	
				3.0E-03	P					3.1	1	1.0	Yes	Styrene-Acrylonitrile (SAN) Trimer (THNP isomer)	57964-40-6					6.0E+01	2.4E+02		4.8E+01	
				1.0E-03	P	2.0E-03	X			-0.77	1	1.0	Yes	Sulfolane	126-33-0					2.0E+01	1.7E+04		2.0E+01	
				8.0E-04	P					3.9	1	0.9	Yes	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9					1.6E+01	3.5E+01		1.1E+01	
						1.0E-03	C	V			1	1.0	Yes	Sulfur Trioxide	7446-11-9							2.1E+00	2.1E+00	
						1.0E-03	C				1	1.0	Yes	Sulfuric Acid	7664-93-9									
2.5E-02	I	7.1E-06	I	5.0E-02	H					4.82	1	0.8	Yes	Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl ester	140-57-8	3.1E+02	2.3E+02		1.3E+02	1.0E+03	8.2E+02		4.5E+02	
				3.0E-02	H					3.3	1	0.9	Yes	TCMTB	21564-17-0					6.0E+02	2.4E+03		4.8E+02	
				7.0E-02	I					1.79	1	1.0	Yes	Tebuthiuron	34014-18-1					1.4E+03	4.7E+04		1.4E+03	
				2.0E-02	H					5.96	1	0.7	No	Temephos	3383-96-8					4.0E+02			4.0E+02	
				1.3E-02	I					1.89	1	1.0	Yes	Terbacil	5902-51-2					2.6E+02	7.0E+03		2.5E+02	
				2.5E-05	H			V		4.48	1	0.9	Yes	Terbufos	13071-79-9					5.0E-01	4.5E-01		2.4E-01	
				1.0E-03	I					3.74	1	0.9	Yes	Terbutryn	886-50-0					2.0E+01	4.1E+01		1.3E+01	
5.0E-03	C	1.3E-06	C					V		1.76	1	1.0	Yes	Tert-Butyl Acetate	540-88-5	1.6E+03	2.4E+04	4.3E+02	3.3E+02					
				1.0E-04	I					6.77	1	0.6	No	Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1					2.0E+00			2.0E+00	
				3.0E-05	P			V		4.64	1	1.0	Yes	Tetrachlorobenzene, 1,2,4,5-	95-94-3					6.0E-01	2.4E-01		1.7E-01	
2.6E-02	I	7.4E-06	I	3.0E-02	I			V		2.93	1	1.0	Yes	Tetrachloroethane, 1,1,1,2-	630-20-6	3.0E+02	1.1E+03	7.6E+01	5.7E+01	6.0E+02	2.4E+03		4.8E+02	
2.0E-01	I	5.8E-05	C	2.0E-02	I			V		2.39	1	1.0	Yes	Tetrachloroethane, 1,1,2,2-	79-34-5	3.9E+01	3.3E+02	9.7E+00	7.6E+00	4.0E+02	3.6E+03		3.6E+02	
2.1E-03	I	2.6E-07	I	6.0E-03	I	4.0E-02	I	V		3.4	1	1.0	Yes	Tetrachloroethylene	127-18-4	3.7E+03	6.5E+03	2.2E+03	1.1E+03	1.2E+02	2.3E+02	8.3E+01	4.1E+01	5.0E+00
				3.0E-02	I					4.45	1	0.9	Yes	Tetrachlorophenol, 2,3,4,6-	58-90-2					6.0E+02	3.9E+02		2.4E+02	
1.6E+01	X			6.0E-05	X			V		4.54	1	0.9	Yes	Tetrachlorotoluene, p- alpha, alpha, alpha-	5216-25-1	4.9E-01	2.5E-01		1.7E-01	1.2E+00	6.8E-01		4.3E-01	
				5.0E-04	I					3.99	1	0.9	Yes	Tetraethyl Dithiopyrophosphate	3689-24-5					1.0E+01	2.4E+01		7.1E+00	
						8.0E+01	I	V		1.68	1	1.0	Yes	Tetrafluoroethane, 1,1,1,2-	811-97-2							1.7E+05	1.7E+05	
				1.0E-04	X					-1.32	1	1.0	Yes	Tetramethylphosphoramide, -N,N,N',N'- (TMPA)	16853-36-4					2.0E+00	4.8E+03		2.0E+00	
				2.0E-03	P					1.64	1	1.0	Yes	Tetryl (Trinitrophenylmethylnitramine)	479-45-8					4.0E+01	2.5E+03		3.9E+01	
				2.0E-05	G						1	0.9	Yes	Thallic Oxide	1314-32-5					4.0E-01	9.1E+01		4.0E-01	
				1.0E-05	X						1	1.0	Yes	Thallium (I) Nitrate	10102-45-1					2.0E-01	4.6E+01		2.0E-01	
				1.0E-05	X						1	1.0	Yes	Thallium (Soluble Salts)	7440-28-0					2.0E-01	4.6E+01		2.0E-01	2.0E+00
				1.0E-05	X			V		-0.17	1	1.0	Yes	Thallium Acetate	563-68-8					2.0E-01	1.7E+02		2.0E-01	
				2.0E-05	X					-0.86	1	1.0	Yes	Thallium Carbonate	6533-73-9					4.0E-01	9.3E+04		4.0E-01	
				1.0E-05	X						1	1.0	Yes	Thallium Chloride	7791-12-0					2.0E-01	4.6E+01		2.0E-01	
				1.0E-05	G						1	1.0	Yes	Thallium Selenite	12039-52-0					2.0E-01	4.6E+01		2.0E-01	
				2.0E-05	X						1	0.9	Yes	Thallium Sulfate	7446-18-6					4.0E-01	9.1E+01		4.0E-01	
				4.3E-02	O					1.56	1	1.0	Yes	Thifensulfuron-methyl	79277-27-3					8.6E+02	1.2E+05		8.6E+02	
				1.0E-02	I					3.4	1	0.9	Yes	Thiobencarb	28249-77-6					2.0E+02	7.7E+02		1.6E+02	
				7.0E-02	X					-0.63	1	1.0	Yes	Thiodiglycol	111-48-8					1.4E+03	9.7E+05		1.4E+03	
				3.0E-04	H					2.16	1	1.0	Yes	Thiofanox	39196-18-4					6.0E+00	4.4E+01		5.3E+00	
1.2E-02	O			1.6E-01	O					1.4	1	1.0	Yes	Thiophanate, Methyl	23564-05-8	6.7E+02	7.9E+04		6.7E+02	3.2E+03	4.1E+05		3.2E+03	
				1.5E-02	O					1.73	1	1.0	Yes	Thiram	137-26-8					3.0E+02	1.2E+04		2.9E+02	
				6.0E-01	H						1	1.0	Yes	Tin	7440-31-5					1.2E+04	2.7E+06		1.2E+04	
						1.0E-04	A	V			1	1.0	Yes	Titanium Tetrachloride	7550-45-0							2.1E-01	2.1E-01	
				8.0E-02	I	5.0E+00	I	V		2.73	1	1.0	Yes	Toluene	108-88-3					1.6E+03	5.3E+03	1.0E+04	1.1E+03	1.0E+03
3.9E-02	C	1.1E-05	C			8.0E-06	C	V		3.74	1	1.0	Yes	Toluene-2,4-diisocyanate	584-84-9	2.0E+02	2.2E+01	5.1E+01	1.4E+01			1.7E-02	1.7E-02	
1.8E-01	X			2.0E-04	X					0.16	1	1.0	Yes	Toluene-2,5-diamine	95-70-5	4.3E+01	8.2E+03		4.3E+01	4.0E+00	8.3E+02		4.0E+00	
3.9E-02	C	1.1E-05	C			8.0E-06	C	V		3.74	1	1.0	Yes	Toluene-2,6-diisocyanate	91-08-7	2.0E+02	2.2E+02	5.1E+01	3.4E+01			1.7E-02	1.7E-02	
				1.0E-04	X					0.71	1	1.0	Yes	Toluenediamine, 2,3-	2687-25-4					2.0E+00	1.8E+02		2.0E+00	
				1.0E-04	X					0.66	1	1.0	Yes	Toluenediamine, 3,4-	496-72-0					2.0E+00	1.9E+02		2.0E+00	
				5.0E-03	P					2.27	1	1.0	Yes	Toluic Acid, p-	99-94-5					1.0E+02	8.9E+02		9.0E+01	
1.6E-02	P	5.1E-05	C							1.32	1	1.0	Yes	Toluidine, o- (Methylaniline, 2-)	95-53-4	4.9E+02	1.4E+04		4.7E+02					
3.0E-02	P			4.0E-03	X					1.39	1	1.0	Yes	Toluidine, p-	106-49-0	2.6E+02	6.8E+03		2.5E+02	8.0E+01	2.3E+03		7.7E+01	
				3.0E+00	P			V		6.1	1	1.0	No	Total Petroleum Hydrocarbons (Aliphatic High)	E1790670					6.0E+04			6.0E+04	
						6.0E-01	P	V		3.9	1	1.0	Yes	Total Petroleum Hydrocarbons (Aliphatic Low)	E1790666							1.3E+03	1.3E+03	
				1.0E-02	X	1.0E-01	P	V		5.65	1	1.0	No	Total Petroleum Hydrocarbons (Aliphatic Medium)	E1790668					2.0E+02		2.1E+02	1.0E+02	
				4.0E-02	P					5.16	1	1.0	No	Total Petroleum Hydrocarbons (Aromatic High)	E1790676					8.0E+02			8.0E+02	

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1				
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	log K_{ow} (unitless)	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
				4.0E-03	P	3.0E-02	P	V		2.13	1	1.0	Yes	Total Petroleum Hydrocarbons (Aromatic Low)	E1790672					8.0E+01	6.1E+02	6.3E+01	3.3E+01	
				4.0E-03	P	3.0E-03	P	V		3.58	1	1.0	Yes	Total Petroleum Hydrocarbons (Aromatic Medium)	E1790674					8.0E+01	9.0E+01	6.3E+00	5.5E+00	
1.1E+00	I	3.2E-04	I	9.0E-05	P					5.9	1	0.8	No	Toxaphene	8001-35-2	7.1E+00			7.1E+00	1.8E+00			1.8E+00	3.0E+00
				3.0E-05	X					5.9	1	0.8	No	Toxaphene, Weathered	E1841606					6.0E-01			6.0E-01	
				7.5E-03	I					7.56	1	0.5	No	Tralomethrin	66841-25-6					1.5E+02			1.5E+02	
				3.0E-04	A			V		4.1	1	0.9	Yes	Tri-n-butyltin	688-73-3					6.0E+00	9.9E+00		3.7E+00	
				8.0E+01	X					0.25	1	1.0	Yes	Triacetin	102-76-1					1.6E+06	5.3E+08		1.6E+06	
				3.4E-02	O					2.77	1	1.0	Yes	Triadimefon	43121-43-3					6.8E+02	7.8E+03		6.3E+02	
7.2E-02	O			2.5E-02	O			V		4.6	1	0.9	Yes	Triallate	2303-17-5	1.1E+02	8.3E+01		4.7E+01	5.0E+02	4.2E+02		2.3E+02	
				1.0E-02	I					1.1	1	1.0	Yes	Triasulfuron	82097-50-5					2.0E+02	6.0E+04		2.0E+02	
				8.0E-03	I					0.78	1	1.0	Yes	Tribenuron-methyl	101200-48-0					1.6E+02	5.0E+03		1.6E+02	
				5.0E-03	I			V		4.66	1	0.9	Yes	Tribromobenzene, 1,2,4-	615-54-3					1.0E+02	8.1E+01		4.5E+01	
				9.0E-03	X					4.13	1	0.9	Yes	Tribromophenol, 2,4,6-	118-79-6					1.8E+02	3.7E+02		1.2E+02	
				2.0E-04	O					5.7	1	0.9	Yes	Tribufos	78-48-8					4.0E+00	6.6E-01		5.7E-01	
9.0E-03	P			1.0E-02	P					4	1	0.9	Yes	Tributyl Phosphate	126-73-8	8.7E+02	1.3E+03		5.2E+02	2.0E+02	3.3E+02		1.2E+02	
				3.0E-04	P						1	0.0	No	Tributyltin Compounds	E1790679					6.0E+00			6.0E+00	
				3.0E-04	I					4.05	1	1.0	Yes	Tributyltin Oxide	56-35-9					6.0E+00	9.5E+01		5.7E+00	
											1	0.0	No	Trichloramine	10025-85-1									4.0E+03(G)
				3.0E+01	I	5.0E+00	P	V		3.16	1	1.0	Yes	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1					6.0E+05	1.9E+06	1.0E+04	1.0E+04	
7.0E-02	I			2.0E-02	I					1.33	1	1.0	Yes	Trichloroacetic Acid	76-03-9	1.1E+02	4.6E+03		1.1E+02	4.0E+02	1.8E+04		3.9E+02	6.0E+01(G)
2.9E-02	H									-0.67	1	1.0	Yes	Trichloroaniline HCl, 2,4,6-	33663-50-2	2.7E+02	3.7E+05		2.7E+02					
7.0E-03	X			3.0E-05	X					3.52	1	1.0	Yes	Trichloroaniline, 2,4,6-	634-93-5	1.1E+03	2.0E+03		7.1E+02	6.0E-01	1.2E+00		4.0E-01	
				8.0E-04	X			V		4.05	1	1.0	Yes	Trichlorobenzene, 1,2,3-	87-61-6					1.6E+01	1.3E+01		7.0E+00	
2.9E-02	P			1.0E-02	I	2.0E-03	P	V		4.02	1	1.0	Yes	Trichlorobenzene, 1,2,4-	120-82-1	2.7E+02	2.0E+02		1.2E+02	2.0E+02	1.6E+02	4.2E+00	4.0E+00	7.0E+01
				2.0E+00	I	5.0E+00	I	V		2.49	1	1.0	Yes	Trichloroethane, 1,1,1-	71-55-6					4.0E+04	2.5E+05	1.0E+04	8.0E+03	2.0E+02
5.7E-02	I	1.6E-05	I	4.0E-03	I	2.0E-04	X	V		1.89	1	1.0	Yes	Trichloroethane, 1,1,2-	79-00-5	1.4E+02	2.0E+03	3.5E+01	2.8E+01	8.0E+01	1.3E+03	4.2E-01	4.1E-01	5.0E+00
4.6E-02	I	4.1E-06	I	5.0E-04	I	2.0E-03	I	V	M	2.42	1	1.0	Yes	Trichloroethylene	79-01-6	1.2E+02	7.4E+02	9.6E+01	4.9E+01	1.0E+01	6.9E+01	4.2E+00	2.8E+00	5.0E+00
				3.0E-01	I			V		2.53	1	1.0	Yes	Trichlorofluoromethane	75-69-4					6.0E+03	3.6E+04		5.2E+03	
				1.0E-01	I					3.72	1	1.0	Yes	Trichlorophenol, 2,4,5-	95-95-4					2.0E+03	2.9E+03		1.2E+03	
1.1E-02	I	3.1E-06	I	1.0E-03	P					3.69	1	1.0	Yes	Trichlorophenol, 2,4,6-	88-06-2	7.1E+02	9.8E+02		4.1E+02	2.0E+01	3.0E+01		1.2E+01	
				1.0E-02	I					3.31	1	0.9	Yes	Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5					2.0E+02	8.7E+02		1.6E+02	
				8.0E-03	I					3.8	1	0.9	Yes	Trichlorophenoxypropionic acid, -2,4,5	93-72-1					1.6E+02	3.6E+02		1.1E+02	5.0E+01
				5.0E-03	I			V		2.43	1	1.0	Yes	Trichloropropane, 1,1,2-	598-77-6					1.0E+02	7.5E+02		8.8E+01	
3.0E+01	I			4.0E-03	I	3.0E-04	I	V	M	2.27	1	1.0	Yes	Trichloropropane, 1,2,3-	96-18-4	8.4E-02	7.3E-01		7.5E-02	8.0E+01	7.7E+02	6.3E-01	6.2E-01	
				3.0E-03	X	3.0E-04	P	V		2.78	1	1.0	Yes	Trichloropropene, 1,2,3-	96-19-5					6.0E+01	2.6E+02	6.3E-01	6.2E-01	
				2.0E-02	A					5.11	1	0.8	Yes	Tricresyl Phosphate (TCP)	1330-78-5					4.0E+02	2.6E+02		1.6E+02	
				3.0E-03	I					5.18	1	0.8	Yes	Tridiphane	58138-08-2					6.0E+01	2.6E+01		1.8E+01	
						7.0E-03	I	V		1.45	1	1.0	Yes	Triethylamine	121-44-8							1.5E+01	1.5E+01	
				2.0E+00	P					-1.75	1	1.0	Yes	Triethylene Glycol	112-27-6					4.0E+04	1.8E+08		4.0E+04	
						2.0E+01	P	V		1.74	1	1.0	Yes	Trifluoroethane, 1,1,1-	420-46-2							4.2E+04	4.2E+04	
7.7E-03	I			7.5E-03	I			V		5.34	1	0.8	Yes	Trifluralin	1582-09-8	1.0E+03	3.4E+02		2.6E+02	1.5E+02	5.5E+01		4.0E+01	
2.0E-02	P			1.0E-02	P					-0.65	1	1.0	Yes	Trimethyl Phosphate	512-56-1	3.9E+02	2.8E+05		3.9E+02	2.0E+02	1.6E+05		2.0E+02	
				1.0E-02	I	6.0E-02	I	V		3.66	1	1.0	Yes	Trimethylbenzene, 1,2,3-	526-73-8					2.0E+02	1.9E+02	1.3E+02	5.5E+01	
				1.0E-02	I	6.0E-02	I	V		3.63	1	1.0	Yes	Trimethylbenzene, 1,2,4-	95-63-6					2.0E+02	2.0E+02	1.3E+02	5.6E+01	
				1.0E-02	I	6.0E-02	I	V		3.42	1	1.0	Yes	Trimethylbenzene, 1,3,5-	108-67-8					2.0E+02	2.8E+02	1.3E+02	6.0E+01	
				1.0E-02	X			V		4.08	1	1.0	Yes	Trimethylpentene, 2,4,4-	25167-70-8					2.0E+02	4.7E+01		3.8E+01	
				3.0E-02	I					1.18	1	1.0	Yes	Trinitrobenzene, 1,3,5-	99-35-4					6.0E+02	4.7E+04		5.9E+02	
3.0E-02	I			5.0E-04	I					1.6	1	1.0	Yes	Trinitrotoluene, 2,4,6-	118-96-7	2.6E+02	1.1E+04		2.5E+02	1.0E+01	4.5E+02		9.8E+00	
				2.0E-02	P					2.83	1	1.0	Yes	Triphenylphosphine Oxide	791-28-6					4.0E+02	3.8E+03		3.6E+02	
				2.0E-02	A					3.65	1	0.9	Yes	Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8					4.0E+02	3.2E+03		3.6E+02	
				1.0E-02	X					2.59	1	1.0	Yes	Tris(1-chloro-2-propyl)phosphate	13674-84-5					2.0E+02	3.8E+03		1.9E+02	
2.3E+00	C	6.6E-04	C					V		4.29	1	1.0	No	Tris(2,3-dibromopropyl)phosphate	126-72-7	3.4E+00		8.5E-01	6.8E-01					
2.0E-02	P			7.0E-03	P					1.44	1	1.0	Yes	Tris(2-chloroethyl)phosphate	115-96-8	3.9E+02	3.0E+04		3.8E+02	1.4E+02	1.2E+04		1.4E+02	
3.2E-03	P			1.0E-01	P					9.49	1	0.0	No	Tris(2-ethylhexyl)phosphate	78-42-2	2.4E+03			2.4E+03	2.0E+03			2.0E+03	
				8.0E-04	P						1	1.0	Yes	Tungsten	7440-33-7					1.6E+01	3.6E+03		1.6E+01	
				2.0E-04	A	4.0E-05	A				1	1.0	Yes	Uranium	7440-61-1					4.0E+00	9.1E+02		4.0E+00	3.0E+01
1.0E+00	C	2.9E-04	C						M	-0.15	1	1.0	Yes	Urethane	51-79-6	2.5E+00	6.1E+02		2.5E+00					
		8.3E-03	P	9.0E-03	I	7.0E-06	P				0.026	1.0	Yes	Vanadium Pentoxide	1314-62-1					1.8E+02	1.1E+03		1.5E+02	
				5.0E-03	G	1.0E-04	A				0.026	1.0	Yes	Vanadium and Compounds	7440-62-2					1.0E+02	6.0E+02		8.6E+01	
				1.0E-03	I			V		3.84	1	1.0	Yes	Vernolate	1929-77-7					2.0E+01	2.5E+01		1.1E+01	
				1.2E-03	O					3.1	1	0.9	Yes	Vinclozolin	50471-44-8					2.4E+01	1.8E+02		2.1E+01	
				1.0E+00	H	2.0E-01	I	V		0.73	1	1.0	Yes	Vinyl Acetate	108-05-4					2.0E+04	1.4E+06	4.2E+02	4.1E+02	
		1.5E-05	P			3.0E-03	I	V		1.57	1	1.0	Yes	Vinyl Bromide	593-60-2			3.7E+01	3.7E+01			6.3E+00	6.3E+00	
7.2E-01	I	4.4E-06	I	3.0E-03	I	8.0E-02	A	V	M	1.38	1	1.0	Yes	Vinyl Chloride	75-01-4	2.1E+00	2.8E+01	3.4E+01	1.9E+00	6.0E+01	8.9E+02	1.7E+02	4.2E+01	2.0E+00
				3.0E-04	I					2.7	1	1.0	Yes	Warfarin	81-81-2					6.0E+00	8.4E+01		5.6E+00	
				2.0E-01	G	1.0E-01	G	V		3.2	1	1.0	Yes	Xylene, m-	108-38-3					4.0E+03	7.1E+03	2.1E+02	1.9E+02	
				2.0E-01	G	1.0E-01	G	V		3.12	1	1.0	Yes	Xylene, o-	95-47-6					4.0E+03	8.0E+03	2.1E+02	1.9E+02	
				2.0E-01	G	1.0E-01	G	V		3.15	1	1.0	Yes	Xylene, p-	106-42-3					4.0E+03	7.6E+03	2.1E+02	1.9E+02	
				2.0E-01	I	1.0E-01	I	V		3.16	1	1.0	Yes	Xylenes	1330-20-7					4.0E+03	7.5E+03	2.1E+02	1.9E+02	1.0E+04
				3.0E-04	I						1	1.0	Yes	Zinc Phosphide	1314-84-7					6.0E+00	2.3E+03		6.0E+00	
				3.0E-01	I						1	1.0	Yes	Zinc and Compounds	7440-66-6					6.0E+03	2.3E+06		6.0E+03	
				5.0E-02	I					1.3	1	1.0	Yes	Zineb	12122-67-7					1.0E+03	9.7E+04		9.9E+02	
				8.0E-05	X						1	1.0	Yes	Zirconium	7440-67-7					1.6E+00	3.6E+02		1.6E+00	

OMLS TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
						Acephate	30560-19-1		
2.2E-06	I	9.0E-03	I	V		Acetaldehyde	75-07-0	1.3E+02	9.4E+00
						Acetochlor	34256-82-1		
				V		Acetone	67-64-1		
		2.0E-03	X			Acetone Cyanohydrin	75-86-5		2.1E+00
		6.0E-02	I	V		Acetonitrile	75-05-8		6.3E+01
				V		Acetophenone	98-86-2		
1.3E-03	C					Acetylaminofluorene, 2-	53-96-3	2.2E-01	
		2.0E-05	I	V		Acrolein	107-02-8		2.1E-02
1.0E-04	I	6.0E-03	I		M	Acrylamide	79-06-1	1.0E+00	6.3E+00
		2.0E-04	P	V		Acrylic Acid	79-10-7		2.1E-01
6.8E-05	I	2.0E-03	I	V		Acrylonitrile	107-13-1	4.1E+00	2.1E+00
		6.0E-03	P			Adiponitrile	111-69-3		6.3E+00
						Alachlor	15972-60-8		
						Aldicarb	116-06-3		
						Aldicarb Sulfone	1646-88-4		
						Aldicarb sulfoxide	1646-87-3		
4.9E-03	I			V		Aldrin	309-00-2	5.7E-02	
		1.0E-04	X	V		Allyl Alcohol	107-18-6		1.0E-01
6.0E-06	C	1.0E-03	I	V		Allyl Chloride	107-05-1	4.7E+01	1.0E+00
		5.0E-03	P			Aluminum	7429-90-5		5.2E+00
						Aluminum Phosphide	20859-73-8		
						Ametryn	834-12-8		
6.0E-03	C					Aminobiphenyl, 4-	92-67-1	4.7E-02	
						Aminophenol, m-	591-27-5		
						Aminophenol, o-	95-55-6		
						Aminophenol, p-	123-30-8		
						Amitraz	33089-61-1		
		5.0E-01	I	V		Ammonia	7664-41-7		5.2E+02
						Ammonium Picrate	131-74-8		
						Ammonium Sulfamate	7773-06-0		
		3.0E-03	X	V		Amyl Alcohol, tert-	75-85-4		3.1E+00
1.6E-06	C	1.0E-03	I			Aniline	62-53-3	1.8E+02	1.0E+00
						Anthraquinone, 9,10-	84-65-1		
		3.0E-04	A			Antimony (metallic)	7440-36-0		3.1E-01
						Antimony Pentoxide	1314-60-9		
						Antimony Tetroxide	1332-81-6		
		2.0E-04	I			Antimony Trioxide	1309-64-4		2.1E-01
4.3E-03	I	1.5E-05	C			Arsenic, Inorganic	7440-38-2	6.5E-02	1.6E-02
		5.0E-05	I			Arsine	7784-42-1		5.2E-02
						Asbestos (units in fibers)	1332-21-4		
						Asulam	3337-71-1		
						Atrazine	1912-24-9		
2.5E-04	C					Auramine	492-80-8	1.1E+00	
						Avermectin B1	65195-55-3		
		1.0E-02	A			Azinphos-methyl	86-50-0		1.0E+01
3.1E-05	I			V		Azobenzene	103-33-3	9.1E+00	
		7.0E-06	P			Azodicarbonamide	123-77-3		7.3E-03
		5.0E-04	H			Barium	7440-39-3		5.2E-01
				V		Benfluralin	1861-40-1		
						Benomyl	17804-35-2		
						Bensulfuron-methyl	83055-99-6		
						Bentazon	25057-89-0		
				V		Benzaldehyde	100-52-7		
7.8E-06	I	3.0E-02	I	V		Benzene	71-43-2	3.6E+01	3.1E+01
						Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1		
				V		Benzenethiol	108-98-5		
6.7E-02	I				M	Benzidine	92-87-5	1.5E-03	
						Benzoic Acid	65-85-0		
				V		Benzotrichloride	98-07-7		
						Benzyl Alcohol	100-51-6		
4.9E-05	C	1.0E-03	P	V		Benzyl Chloride	100-44-7	5.7E+00	1.0E+00
2.4E-03	I	2.0E-05	I			Beryllium and compounds	7440-41-7	1.2E-01	2.1E-02
						Bifenox	42576-02-3		
						Biphenthrin	82657-04-3		
		4.0E-04	X	V		Biphenyl, 1,1'-	92-52-4		4.2E-01
				V		Bis(2-chloro-1-methylethyl) ether	108-60-1		
						Bis(2-chloroethoxy)methane	111-91-1		
3.3E-04	I			V		Bis(2-chloroethyl)ether	111-44-4	8.5E-01	
6.2E-02	I			V		Bis(chloromethyl)ether	542-88-1	4.5E-03	
						Bisphenol A	80-05-7		
		2.0E-02	H			Boron And Borates Only	7440-42-8		2.1E+01
		2.0E-02	P	V		Boron Trichloride	10294-34-5		2.1E+01
		1.3E-02	C	V		Boron Trifluoride	7637-07-2		1.4E+01
						Bromate	15541-45-4		
		6.0E-05	X	V		Bromo-2-chloroethane, 1-	107-04-0		6.3E-02
				V		Bromo-3-fluorobenzene, 1-	1073-06-9		
				V		Bromo-4-fluorobenzene, 1-	460-00-4		
						Bromoacetic acid	79-08-3		
		6.0E-02	I	V		Bromobenzene	108-86-1		6.3E+01
		4.0E-02	X	V		Bromochloromethane	74-97-5		4.2E+01
3.7E-05	C			V		Bromodichloromethane	75-27-4	7.6E+00	
1.1E-06	I			V		Bromoform	75-25-2	2.6E+02	
		5.0E-03	I	V		Bromomethane	74-83-9		5.2E+00
				V		Bromophos	2104-96-3		
		1.0E-01	A	V		Bromopropane, 1-	106-94-5		1.0E+02
						Bromoxynil	1689-84-5		
				V		Bromoxynil Octanoate	1689-99-2		
3.0E-05	I	2.0E-03	I	V		Butadiene, 1,3-	106-99-0	9.4E+00	2.1E+00
				V		Butanol, N-	71-36-3		

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
		5.0E+00	I	V		Butyl Alcohol, t-	75-65-0		5.2E+03
		3.0E+01	P	V		Butyl alcohol, sec-	78-92-2		3.1E+04
				V		Butylate	2008-41-5		
5.7E-08	C					Butylated hydroxyanisole	25013-16-5	4.9E+03	
						Butylated hydroxytoluene	128-37-0		
				V		Butylbenzene, n-	104-51-8		
				V		Butylbenzene, sec-	135-98-8		
				V		Butylbenzene, tert-	98-06-6		
						Cacodylic Acid	75-60-5		
1.8E-03	I	1.0E-05	A			Cadmium (Diet)	7440-43-9	1.6E-01	1.0E-02
1.8E-03	I	1.0E-05	A			Cadmium (Water)	7440-43-9	1.6E-01	1.0E-02
		2.2E-03	C			Caprolactam	105-60-2		2.3E+00
4.3E-05	C					Captafol	2425-06-1	6.5E+00	
6.6E-07	C					Captan	133-06-2	4.3E+02	
						Carbaryl	63-25-2		
						Carbofuran	1563-66-2		
		7.0E-01	I	V		Carbon Disulfide	75-15-0		7.3E+02
6.0E-06	I	1.0E-01	I	V		Carbon Tetrachloride	56-23-5	4.7E+01	1.0E+02
		1.0E-01	P	V		Carbonyl Sulfide	463-58-1		1.0E+02
						Carbosulfan	55285-14-8		
						Carboxin	5234-68-4		
		9.0E-04	I			Ceric oxide	1306-38-3		9.4E-01
				V		Chloral Hydrate	302-17-0		
						Chloramben	133-90-4		
						Chloramines, Organic	E701235		
						Chloranil	118-75-2		
				V		Chlordane (alpha)	5103-71-9		
				V		Chlordane (gamma)	5103-74-2		
1.0E-04	I	7.0E-04	I	V		Chlordane (technical mixture)	12789-03-6	2.8E+00	7.3E-01
4.6E-03	C					Chlordecone (Kepone)	143-50-0	6.1E-02	
						Chlorfenvinphos	470-90-6		
						Chlorimuron, Ethyl-	90982-32-4		
		1.5E-04	A	V		Chlorine	7782-50-5		1.5E-01
		2.0E-04	I	V		Chlorine Dioxide	10049-04-4		2.1E-01
						Chlorite (Sodium Salt)	7758-19-2		
		5.0E+01	I	V		Chloro-1,1-difluoroethane, 1-	75-68-3		5.2E+04
3.0E-04	I	2.0E-02	I	V		Chloro-1,3-butadiene, 2-	126-99-8	9.4E-01	2.1E+01
						Chloro-2-methylaniline HCl, 4-	3165-93-3		
7.7E-05	C					Chloro-2-methylaniline, 4-	95-69-2	3.6E+00	
				V		Chloroacetaldehyde, 2-	107-20-0		
						Chloroacetic Acid	79-11-8		
		3.0E-05	I			Chloroacetophenone, 2-	532-27-4		3.1E-02
						Chloroaniline, p-	106-47-8		
		5.0E-02	P	V		Chlorobenzene	108-90-7		5.2E+01
						Chlorobenzene sulfonic acid, p-	98-66-8		
3.1E-05	C					Chlorobenzilate	510-15-6	9.1E+00	
						Chlorobenzoic Acid, p-	74-11-3		
8.6E-06	C	3.0E-01	P	V		Chlorobenzotrifluoride, 4-	98-56-6	3.3E+01	3.1E+02
				V		Chlorobutane, 1-	109-69-3		
		5.0E+01	I	V		Chlorodifluoromethane	75-45-6		5.2E+04
				V		Chloroethanol, 2-	107-07-3		
2.3E-05	I	9.8E-02	A	V		Chloroform	67-66-3	1.2E+01	1.0E+02
		9.0E-02	I	V		Chloromethane	74-87-3		9.4E+01
6.9E-04	C			V		Chloromethyl Methyl Ether	107-30-2	4.1E-01	
		1.0E-05	X			Chloronitrobenzene, o-	88-73-3		1.0E-02
		2.0E-03	P			Chloronitrobenzene, p-	100-00-5		2.1E+00
				V		Chlorophenol, 2-	95-57-8		
		4.0E-04	C	V		Chloropicrin	76-06-2		4.2E-01
						Chlorothalonil	1897-45-6		
				V		Chlorotoluene, o-	95-49-8		
				V		Chlorotoluene, p-	106-43-4		
6.9E-02	C					Chlorozotocin	54749-90-5	4.1E-03	
						Chlorpropham	101-21-3		
						Chlorpyrifos	2921-88-2		
						Chlorpyrifos Methyl	5598-13-0		
						Chlorsulfuron	64902-72-3		
						Chlorthal-dimethyl	1861-32-1		
						Chlorthiophos	60238-56-4		
						Chromium(III), Insoluble Salts	16065-83-1		
8.4E-02	G	1.0E-04	I		M	Chromium(VI)	18540-29-9	1.2E-03	1.0E-01
						Chromium, Total	7440-47-3		
						Clofentezine	74115-24-5		
9.0E-03	P	6.0E-06	P			Cobalt	7440-48-4	3.1E-02	6.3E-03
6.2E-04	I			V	M	Coke Oven Emissions	E649830	1.6E-01	
						Copper	7440-50-8		
		6.0E-01	C			Cresol, m-	108-39-4		6.3E+02
		6.0E-01	C			Cresol, o-	95-48-7		6.3E+02
		6.0E-01	C			Cresol, p-	106-44-5		6.3E+02
						Cresol, p-chloro-m-	59-50-7		
		6.0E-01	C			Cresols	1319-77-3		6.3E+02
				V		Crotonaldehyde, trans-	123-73-9		
		4.0E-01	I	V		Cumene	98-82-8		4.2E+02
6.3E-05	C					Cupferron	135-20-6	4.5E+00	
						Cyanazine	21725-46-2		
						Cyanides			
						~Calcium Cyanide	592-01-8		
						~Copper Cyanide	544-92-3		
		8.0E-04	G	V		~Cyanide (CN-)	57-12-5		8.3E-01
				V		~Cyanogen	460-19-5		
				V		~Cyanogen Bromide	506-68-3		

RMLs TR=1E-04 THQ=1.0

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Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
				V		~Cyanogen Chloride	506-77-4		
		8.0E-04	I	V		~Hydrogen Cyanide	74-90-8		8.3E-01
						~Potassium Cyanide	151-50-8		
						~Potassium Silver Cyanide	506-61-6		
						~Silver Cyanide	506-64-9		
						~Sodium Cyanide	143-33-9		
						~Thiocyanates	E1790665		
				V		~Thiocyanic Acid	463-56-9		
						~Zinc Cyanide	557-21-1		
		6.0E+00	I	V		Cyclohexane	110-82-7		6.3E+03
						Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3		
		7.0E-01	P	V		Cyclohexanone	108-94-1		7.3E+02
		1.0E+00	X	V		Cyclohexene	110-83-8		1.0E+03
				V		Cyclohexylamine	108-91-8		
						Cyfluthrin	68359-37-5		
						Cypermethrin	52315-07-8		
						Cyromazine	66215-27-8		
6.9E-05	C					DDD, p,p`- (DDD)	72-54-8	4.1E+00	
9.7E-05	C			V		DDE, p,p'-	72-55-9	2.9E+00	
9.7E-05	I					DDT	50-29-3	2.9E+00	
						Dalapon	75-99-0		
5.1E-06	C					Daminozide	1596-84-5	5.5E+01	
						Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5		
						Demeton	8065-48-3		
						Di(2-ethylhexyl)adipate	103-23-1		
						Diallate	2303-16-4		
						Diazinon	333-41-5		
				V		Dibenzothiophene	132-65-0		
6.0E-03	P	2.0E-04	I	V	M	Dibromo-3-chloropropane, 1,2-	96-12-8	1.7E-02	2.1E-01
						Dibromoacetic acid	631-64-1		
				V		Dibromobenzene, 1,3-	108-36-1		
				V		Dibromobenzene, 1,4-	106-37-6		
				V		Dibromochloromethane	124-48-1		
6.0E-04	I	9.0E-03	I	V		Dibromoethane, 1,2-	106-93-4	4.7E-01	9.4E+00
		4.0E-03	X	V		Dibromomethane (Methylene Bromide)	74-95-3		4.2E+00
						Dibutyltin Compounds	E1790661		
						Dicamba	1918-00-9		
						Dichloramine	3400-09-7		
4.2E-03	P			V		Dichloro-2-butene, 1,4-	764-41-0	6.7E-02	
4.2E-03	P			V		Dichloro-2-butene, cis-1,4-	1476-11-5	6.7E-02	
4.2E-03	P			V		Dichloro-2-butene, trans-1,4-	110-57-6	6.7E-02	
						Dichloroacetic Acid	79-43-6		
		2.0E-01	H	V		Dichlorobenzene, 1,2-	95-50-1		2.1E+02
1.1E-05	C	8.0E-01	I	V		Dichlorobenzene, 1,4-	106-46-7	2.6E+01	8.3E+02
3.4E-04	C					Dichlorobenzidine, 3,3'-	91-94-1	8.3E-01	
						Dichlorobenzophenone, 4,4'-	90-98-2		
		1.0E-01	X	V		Dichlorodifluoromethane	75-71-8		1.0E+02
1.6E-06	C			V		Dichloroethane, 1,1-	75-34-3	1.8E+02	
2.6E-05	I	7.0E-03	P	V		Dichloroethane, 1,2-	107-06-2	1.1E+01	7.3E+00
		2.0E-01	I	V		Dichloroethylene, 1,1-	75-35-4		2.1E+02
				V		Dichloroethylene, cis-1,2-	156-59-2		
		4.0E-02	X	V		Dichloroethylene, trans-1,2-	156-60-5		4.2E+01
						Dichlorophenol, 2,4-	120-83-2		
						Dichlorophenoxy Acetic Acid, 2,4-	94-75-7		
3.7E-06	P	4.0E-03	I	V		Dichloropropane, 1,2-	78-87-5	7.6E+01	4.2E+00
				V		Dichloropropane, 1,3-	142-28-9		
						Dichloropropanol, 2,3-	616-23-9		
4.0E-06	I	2.0E-02	I	V		Dichloropropene, 1,3-	542-75-6	7.0E+01	2.1E+01
8.3E-05	C	5.0E-04	I			Dichlorvos	62-73-7	3.4E+00	5.2E-01
						Dicrotophos	141-66-2		
		3.0E-04	X	V		Dicyclopentadiene	77-73-6		3.1E-01
4.6E-03	I					Dieldrin	60-57-1	6.1E-02	
3.0E-04	C	5.0E-03	I			Diesel Engine Exhaust	E17136615	9.4E-01	5.2E+00
		2.0E-04	P			Diethanolamine	111-42-2		2.1E-01
		1.0E-04	P			Diethylene Glycol Monobutyl Ether	112-34-5		1.0E-01
		3.0E-04	P			Diethylene Glycol Monoethyl Ether	111-90-0		3.1E-01
				V		Diethylformamide	617-84-5		
1.0E-01	C					Diethylstilbestrol	56-53-1	2.8E-03	
						Difenzoquat	43222-48-6		
						Diflubenzuron	35367-38-5		
		4.0E+01	I	V		Difluoroethane, 1,1-	75-37-6		4.2E+04
		3.0E+01	X	V		Difluoropropane, 2,2-	420-45-1		3.1E+04
1.3E-05	C			V		Dihydrosafrole	94-58-6	2.2E+01	
		7.0E-01	P	V		Diisopropyl Ether	108-20-3		7.3E+02
				V		Diisopropyl Methylphosphonate	1445-75-6		
						Dimethipin	55290-64-7		
						Dimethoate	60-51-5		
						Dimethoxybenzidine, 3,3'-	119-90-4		
						Dimethyl methylphosphonate	756-79-6		
1.3E-03	C					Dimethylamino azobenzene [p-]	60-11-7	2.2E-01	
						Dimethylaniline HCl, 2,4-	21436-96-4		
						Dimethylaniline, 2,4-	95-68-1		
				V		Dimethylaniline, N,N-	121-69-7		
						Dimethylbenzidine, 3,3'-	119-93-7		
		3.0E-02	I	V		Dimethylformamide	68-12-2		3.1E+01
		2.0E-06	X	V		Dimethylhydrazine, 1,1-	57-14-7		2.1E-03
1.6E-01	C			V		Dimethylhydrazine, 1,2-	540-73-8	1.8E-03	
						Dimethylphenol, 2,4-	105-67-9		
						Dimethylphenol, 2,6-	576-26-1		
						Dimethylphenol, 3,4-	95-65-8		

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
1.3E-05	C			V		Dimethylvinylchloride	513-37-1	2.2E+01	
						Dinitro-o-cresol, 4,6-	534-52-1		
						Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5		
		2.0E-03	X			Dinitroaniline, 3,5-	618-87-1		2.1E+00
						Dinitrobenzene, 1,2-	528-29-0		
						Dinitrobenzene, 1,3-	99-65-0		
						Dinitrobenzene, 1,4-	100-25-4		
						Dinitrophenol, 2,4-	51-28-5		
						Dinitrotoluene Mixture, 2,4/2,6-	E1615210		
8.9E-05	C					Dinitrotoluene, 2,4-	121-14-2	3.2E+00	
						Dinitrotoluene, 2,6-	606-20-2		
						Dinitrotoluene, 2-Amino-4,6-	35572-78-2		
						Dinitrotoluene, 4-Amino-2,6-	19406-51-0		
						Dinitrotoluene, Technical grade	25321-14-6		
						Dinoseb	88-85-7		
5.0E-06	I	3.0E-02	I	V		Dioxane, 1,4-	123-91-1	5.6E+01	3.1E+01
						Dioxins			
1.3E+00	I					~Hexachlorodibenzo-p-dioxin, Mixture	34465-46-8	2.2E-04	
3.8E+01	C	4.0E-08	C	V		~TCDD, 2,3,7,8-	1746-01-6	7.4E-06	4.2E-05
						Diphenamid	957-51-7		
		4.0E-04	X	V		Diphenyl Ether	101-84-8		4.2E-01
						Diphenyl Sulfone	127-63-9		
						Diphenylamine	122-39-4		
2.2E-04	I					Diphenylhydrazine, 1,2-	122-66-7	1.3E+00	
						Diquat	2764-72-9		
2.1E-03	C					Direct Black 38	1937-37-7	1.3E-01	
2.1E-03	C					Direct Blue 6	2602-46-2	1.3E-01	
1.9E-03	C					Direct Brown 95	16071-86-6	1.5E-01	
						Disulfoton	298-04-4		
				V		Dithiane, 1,4-	505-29-3		
						Diuron	330-54-1		
						Dodine	2439-10-3		
				V		EPTC	759-94-4		
				V		Endosulfan	115-29-7		
						Endosulfan Sulfate	1031-07-8		
						Endothall	145-73-3		
						Endrin	72-20-8		
1.2E-06	I	1.0E-03	I	V		Epichlorohydrin	106-89-8	2.3E+02	1.0E+00
		2.0E-02	I	V		Epoxybutane, 1,2-	106-88-7		2.1E+01
						Ethanol, 2-(2-methoxyethoxy)-	111-77-3		
						Ethephon	16672-87-0		
						Ethion	563-12-2		
		6.0E-02	P	V		Ethoxyethanol Acetate, 2-	111-15-9		6.3E+01
		4.0E-02	P	V		Ethoxyethanol, 2-	110-80-5		4.2E+01
		7.0E-02	P	V		Ethyl Acetate	141-78-6		7.3E+01
		8.0E-03	P	V		Ethyl Acrylate	140-88-5		8.3E+00
		4.0E+00	P	V		Ethyl Chloride (Chloroethane)	75-00-3		4.2E+03
				V		Ethyl Ether	60-29-7		
		3.0E-01	P	V		Ethyl Methacrylate	97-63-2		3.1E+02
8.0E-08	I	4.0E+01	I	V		Ethyl Tertiary Butyl Ether (ETBE)	637-92-3	3.5E+03	4.2E+04
						Ethyl-p-nitrophenyl Phosphonate	2104-64-5		
2.5E-06	C	1.0E+00	I	V		Ethylbenzene	100-41-4	1.1E+02	1.0E+03
						Ethylene Cyanohydrin	109-78-4		
				V		Ethylene Diamine	107-15-3		
		4.0E-01	C			Ethylene Glycol	107-21-1		4.2E+02
		1.6E+00	I			Ethylene Glycol Monobutyl Ether	111-76-2		1.7E+03
3.0E-03	I	3.0E-02	C	V	M	Ethylene Oxide	75-21-8	3.4E-02	3.1E+01
1.3E-05	C					Ethylene Thiourea	96-45-7	2.2E+01	
1.9E-02	C			V		Ethyleneimine	151-56-4	1.5E-02	
						Ethylphthalyl Ethyl Glycolate	84-72-0		
						Fenamiphos	22224-92-6		
						Fenpropathrin	39515-41-8		
						Fenvalerate	51630-58-1		
						Fluometuron	2164-17-2		
		1.3E-02	C			Fluoride	16984-48-8		1.4E+01
		1.3E-02	C			Fluorine (Soluble Fluoride)	7782-41-4		1.4E+01
						Fluridone	59756-60-4		
						Flurprimidol	56425-91-3		
						Flusilazole	85509-19-9		
						Flutolanil	66332-96-5		
						Fluvalinate	69409-94-5		
						Folpet	133-07-3		
						Fomesafen	72178-02-0		
						Fonofos	944-22-9		
1.3E-05	I	9.8E-03	A	V		Formaldehyde	50-00-0	2.2E+01	1.0E+01
		3.0E-04	X	V		Formic Acid	64-18-6		3.1E-01
						Fosetyl-AL	39148-24-8		
						Furans			
				V		~Dibenzofuran	132-64-9		
				V		~Furan	110-00-9		
		2.0E+00	I	V		~Tetrahydrofuran	109-99-9		2.1E+03
						Furazolidone	67-45-8		
		5.0E-02	H	V		Furfural	98-01-1		5.2E+01
4.3E-04	C					Furium	531-82-8	6.5E-01	
8.6E-06	C					Furmecyclox	60568-05-0	3.3E+01	
						Glufosinate, Ammonium	77182-82-2		
		8.0E-05	C			Glutaraldehyde	111-30-8		8.3E-02
		1.0E-03	X	V		Glycidaldehyde	765-34-4		1.0E+00
						Glyphosate	1071-83-6		
				V		Guanidine	113-00-8		

RMLs TR=1E-04 THQ=1.0

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Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
						Guanidine Chloride	50-01-1		
						Guanidine Nitrate	506-93-4		
						Haloxyfop, Methyl	69806-40-2		
1.3E-03	I			V		Heptachlor	76-44-8	2.2E-01	
2.6E-03	I			V		Heptachlor Epoxide	1024-57-3	1.1E-01	
		3.0E-03	X	V		Heptanal, n-	111-71-7		3.1E+00
		4.0E-01	P	V		Heptane, N-	142-82-5		4.2E+02
				V		Hexabromobenzene	87-82-1		
						Hexabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-153)	68631-49-2		
4.6E-04	I			V		Hexachlorobenzene	118-74-1	6.1E-01	
2.2E-05	I			V		Hexachlorobutadiene	87-68-3	1.3E+01	
1.8E-03	I					Hexachlorocyclohexane, Alpha-	319-84-6	1.6E-01	
5.3E-04	I					Hexachlorocyclohexane, Beta-	319-85-7	5.3E-01	
3.1E-04	C					Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	9.1E-01	
5.1E-04	I					Hexachlorocyclohexane, Technical	608-73-1	5.5E-01	
		2.0E-04	I	V		Hexachlorocyclopentadiene	77-47-4		2.1E-01
1.1E-05	C	3.0E-02	I	V		Hexachloroethane	67-72-1	2.6E+01	3.1E+01
						Hexachlorophene	70-30-4		
						Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4		
		1.0E-05	I	V		Hexamethylene Diisocyanate, 1,6-	822-06-0		1.0E-02
		4.0E-04	C			Hexamethylene diisocyanate biuret	4035-89-6		4.2E-01
		4.0E-04	C			Hexamethylene diisocyanate isocyanurate	3779-63-3		4.2E-01
						Hexamethylphosphoramide	680-31-9		
2.0E-07	X	6.0E-01	P	V		Hexane, Commercial	E5241997	1.4E+03	6.3E+02
		7.0E-01	I	V		Hexane, N-	110-54-3		7.3E+02
						Hexanedioic Acid	124-04-9		
		4.0E-04	P	V		Hexanol, 1-,2-ethyl- (2-Ethyl-1-hexanol)	104-76-7		4.2E-01
		3.0E-02	I	V		Hexanone, 2-	591-78-6		3.1E+01
						Hexazinone	51235-04-2		
						Hexythiazox	78587-05-0		
						Hydramethylnon	67485-29-4		
4.9E-03	I	3.0E-05	P	V		Hydrazine	302-01-2	5.7E-02	3.1E-02
4.9E-03	I					Hydrazine Sulfate	10034-93-2	5.7E-02	
		2.0E-02	I	V		Hydrogen Chloride	7647-01-0		2.1E+01
		1.4E-02	C	V		Hydrogen Fluoride	7664-39-3		1.5E+01
		2.0E-03	I	V		Hydrogen Sulfide	7783-06-4		2.1E+00
						Hydroquinone	123-31-9		
						Imazalil	35554-44-0		
						Imazaquin	81335-37-7		
						Imazethapyr	81335-77-5		
						Iodine	7553-56-2		
						Iprodione	36734-19-7		
						Iron	7439-89-6		
				V		Isobutyl Alcohol	78-83-1		
		2.0E+00	C			Isophorone	78-59-1		2.1E+03
				V		Isopropalin	33820-53-0		
		2.0E-01	P	V		Isopropanol	67-63-0		2.1E+02
						Isopropyl Methyl Phosphonic Acid	1832-54-8		
						Isoxaben	82558-50-7		
		3.0E-01	A	V		JP-7	E1737665		3.1E+02
						Lactofen	77501-63-4		
						Lactonitrile	78-97-7		
						Lanthanum	7439-91-0		
						Lanthanum Acetate Hydrate	100587-90-4		
						Lanthanum Chloride Heptahydrate	10025-84-0		
						Lanthanum Chloride, Anhydrous	10099-58-8		
						Lanthanum Nitrate Hexahydrate	10277-43-7		
						Lead Compounds			
1.2E-05	C					~Lead Phosphate	7446-27-7	2.3E+01	
8.0E-05	C					~Lead acetate	301-04-2	3.5E+00	
						~Lead and Compounds	7439-92-1		1.5E-01
1.1E-05	C					~Lead subacetate	1335-32-6	2.6E+01	
				V		~Tetraethyl Lead	78-00-2		
				V		Lewisite	541-25-3		
						Linuron	330-55-2		
						Lithium	7439-93-2		
						MCPA	94-74-6		
						MCPB	94-81-5		
						MCPP	93-65-2		
						Malathion	121-75-5		
		7.0E-04	C			Maleic Anhydride	108-31-6		7.3E-01
						Maleic Hydrazide	123-33-1		
						Malononitrile	109-77-3		
						Mancozeb	8018-01-7		
						Maneb	12427-38-2		
		5.0E-05	I			Manganese (Diet)	7439-96-5		5.2E-02
		5.0E-05	I			Manganese (Non-diet)	7439-96-5		5.2E-02
						Mephosfolan	950-10-7		
						Mepiquat Chloride	24307-26-4		
						Mercaptobenzothiazole, 2-	149-30-4		
						Mercury Compounds			
		3.0E-04	G			~Mercuric Chloride (and other Mercury salts)	7487-94-7		3.1E-01
		3.0E-04	I	V		~Mercury (elemental)	7439-97-6		3.1E-01
						~Methyl Mercury	22967-92-6		
						~Phenylmercuric Acetate	62-38-4		
				V		Merphos	150-50-5		
						Metalaxyl	57837-19-1		
		3.0E-02	P	V		Methacrylonitrile	126-98-7		3.1E+01
						Methamidophos	10265-92-6		
		2.0E+01	I	V		Methanol	67-56-1		2.1E+04

RMLs TR=1E-04 THQ=1.0

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Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
						Methidathion	950-37-8		
						Methomyl	16752-77-5		
						Methoxy-5-nitroaniline, 2-	99-59-2		
						Methoxychlor	72-43-5		
		1.0E-03	P	V		Methoxyethanol Acetate, 2-	110-49-6		1.0E+00
		7.0E-03	P	V		Methoxyethanol, 2-	109-86-4		7.3E+00
				V		Methyl Acetate	79-20-9		
		2.0E-02	P	V		Methyl Acrylate	96-33-3		2.1E+01
		5.0E+00	I	V		Methyl Ethyl Ketone (2-Butanone)	78-93-3		5.2E+03
1.0E-03	X	2.0E-05	X	V		Methyl Hydrazine	60-34-4	2.8E-01	2.1E-02
		3.0E+00	I	V		Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1		3.1E+03
		1.0E-03	C	V		Methyl Isocyanate	624-83-9		1.0E+00
		7.0E-01	I	V		Methyl Methacrylate	80-62-6		7.3E+02
						Methyl Parathion	298-00-0		
						Methyl Phosphonic Acid	993-13-5		
		4.0E-02	H	V		Methyl Styrene (Mixed Isomers)	25013-15-4		4.2E+01
2.8E-05	C					Methyl methanesulfonate	66-27-3	1.0E+01	
2.6E-07	C	3.0E+00	I	V		Methyl tert-Butyl Ether (MTBE)	1634-04-4	1.1E+03	3.1E+03
						Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2		
		3.0E+00	X	V		Methyl-2-Pentanol, 4-	108-11-2		3.1E+03
						Methyl-5-Nitroaniline, 2-	99-55-8		
2.4E-03	C					Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	1.2E-01	
3.7E-05	C					Methylaniline Hydrochloride, 2-	636-21-5	7.6E+00	
						Methylarsonic acid	124-58-3		
						Methylbenzene,1-4-diamine monohydrochloride, 2-	74612-12-7		
						Methylbenzene-1,4-diamine sulfate, 2-	615-50-9		
6.3E-03	C				M	Methylcholanthrene, 3-	56-49-5	1.6E-02	
1.0E-08	I	6.0E-01	I	V	M	Methylene Chloride	75-09-2	1.0E+04	6.3E+02
4.3E-04	C				M	Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	2.4E-01	
1.3E-05	C					Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	2.2E+01	
4.6E-04	C	2.0E-02	C			Methylenebisbenzenamine, 4,4'-	101-77-9	6.1E-01	2.1E+01
		6.0E-04	I			Methylenediphenyl Diisocyanate	101-68-8		6.3E-01
				V		Methylstyrene, Alpha-	98-83-9		
						Metolachlor	51218-45-2		
						Metribuzin	21087-64-9		
						Metsulfuron-methyl	74223-64-6		
4.5E-06	X	1.0E-01	P	V		Midrange Aliphatic Hydrocarbon Streams	E1790669	6.2E+01	1.0E+02
				V		Mineral oils	8012-95-1		
5.1E-03	C			V		Mirex	2385-85-5	5.5E-02	
						Molinate	2212-67-1		
		2.0E-03	A			Molybdenum	7439-98-7		2.1E+00
						Monochloramine	10599-90-3		
						Monomethylaniline	100-61-8		
						Myclobutanil	88671-89-0		
						N,N'-Diphenyl-1,4-benzenediamine	74-31-7		
				V		Naled	300-76-5		
		1.0E-01	P	V		Naphtha, High Flash Aromatic (HFAN)	64742-95-6		1.0E+02
0.0E+00	C					Naphthylamine, 2-	91-59-8		
						Napropamide	15299-99-7		
2.6E-04	C	1.4E-05	C			Nickel Acetate	373-02-4	1.1E+00	1.5E-02
2.6E-04	C	1.4E-05	C			Nickel Carbonate	3333-67-3	1.1E+00	1.5E-02
2.6E-04	C	1.4E-05	C	V		Nickel Carbonyl	13463-39-3	1.1E+00	1.5E-02
2.6E-04	C	1.4E-05	C			Nickel Hydroxide	12054-48-7	1.1E+00	1.5E-02
2.6E-04	C	2.0E-05	C			Nickel Oxide	1313-99-1	1.1E+00	2.1E-02
2.4E-04	I	1.4E-05	C			Nickel Refinery Dust	E715532	1.2E+00	1.5E-02
2.6E-04	C	9.0E-05	A			Nickel Soluble Salts	7440-02-0	1.1E+00	9.4E-02
4.8E-04	I	1.4E-05	C			Nickel Subsulfide	12035-72-2	5.8E-01	1.5E-02
2.6E-04	C	1.4E-05	C			Nickelocene	1271-28-9	1.1E+00	1.5E-02
						Nitrate (measured as nitrogen)	14797-55-8		
						Nitrate + Nitrite (measured as nitrogen)	E701177		
						Nitrite (measured as nitrogen)	14797-65-0		
		5.0E-05	X			Nitroaniline, 2-	88-74-4		5.2E-02
		6.0E-03	P			Nitroaniline, 4-	100-01-6		6.3E+00
4.0E-05	I	9.0E-03	I	V		Nitrobenzene	98-95-3	7.0E+00	9.4E+00
						Nitrocellulose	9004-70-0		
						Nitrofurantoin	67-20-9		
3.7E-04	C					Nitrofurazone	59-87-0	7.6E-01	
						Nitroglycerin	55-63-0		
						Nitroguanidine	556-88-7		
8.8E-06	P	5.0E-03	P	V		Nitromethane	75-52-5	3.2E+01	5.2E+00
5.8E-04	X	2.0E-02	I	V		Nitropropane, 2-	79-46-9	4.8E-01	2.1E+01
7.7E-03	C				M	Nitroso-N-ethylurea, N-	759-73-9	1.3E-02	
3.4E-02	C				M	Nitroso-N-methylurea, N-	684-93-5	3.0E-03	
1.6E-03	I			V		Nitroso-di-N-butylamine, N-	924-16-3	1.8E-01	
2.0E-03	C					Nitroso-di-N-propylamine, N-	621-64-7	1.4E-01	
8.0E-04	C					Nitrosodiethanolamine, N-	1116-54-7	3.5E-01	
4.3E-02	I				M	Nitrosodiethylamine, N-	55-18-5	2.4E-03	
1.4E-02	I	4.0E-05	X	V	M	Nitrosodimethylamine, N-	62-75-9	7.2E-03	4.2E-02
2.6E-06	C					Nitrosodiphenylamine, N-	86-30-6	1.1E+02	
6.3E-03	C			V		Nitrosomethylethylamine, N-	10595-95-6	4.5E-02	
1.9E-03	C					Nitrosomorpholine [N-]	59-89-2	1.5E-01	
2.7E-03	C					Nitrosopiperidine [N-]	100-75-4	1.0E-01	
6.1E-04	I					Nitrosopyrrolidine, N-	930-55-2	4.6E-01	
						Nitrotoluene, m-	99-08-1		
				V		Nitrotoluene, o-	88-72-2		
						Nitrotoluene, p-	99-99-0		
		2.0E-02	P	V		Nonane, n-	111-84-2		2.1E+01
						Norflurazon	27314-13-2		
						Octabromodiphenyl Ether	32536-52-0		
						Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0		

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
						Octamethylpyrophosphoramide	152-16-9		
						Oryzalin	19044-88-3		
						Oxadiazon	19666-30-9		
						Oxamyl	23135-22-0		
						Oxyfluorfen	42874-03-3		
						Paclobutrazol	76738-62-0		
						Paraquat Dichloride	1910-42-5		
						Parathion	56-38-2		
				V		Pebulate	1114-71-2		
						Pendimethalin	40487-42-1		
				V		Pentabromodiphenyl Ether	32534-81-9		
						Pentabromodiphenyl ether, 2,2',4,4',5- (BDE-99)	60348-60-9		
				V		Pentachlorobenzene	608-93-5		
				V		Pentachloroethane	76-01-7		
				V		Pentachloronitrobenzene	82-68-8		
5.1E-06	C					Pentachlorophenol	87-86-5	5.5E+01	
						Pentaerythritol tetranitrate (PETN)	78-11-5		
						Pentamethylphosphoramide (PMPA)	10159-46-3		
		1.0E+00	P	V		Pentane, n-	109-66-0		1.0E+03
						Per- and Polyfluoroalkyl Substances (PFAS)			
						~Ammonium perfluoro-2-methyl-3-oxahexanoate	62037-80-3		
				V		~Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6		
						~Perfluorobutanesulfonate	45187-15-3		
						~Perfluorobutanesulfonic acid (PFBS)	375-73-5		
						~Perfluorohexanesulfonate	108427-53-8		
						~Perfluorohexanesulfonic acid (PFHxS)	355-46-4		
						~Perfluorononanoate	72007-68-2		
						~Perfluorononanoic acid (PFNA)	375-95-1		
						~Perfluorooctanesulfonate	45298-90-6		
						~Perfluorooctanesulfonic acid (PFOS)	1763-23-1		
						~Perfluorooctanoate	45285-51-6		
						~Perfluorooctanoic acid (PFOA)	335-67-1		
						~Potassium perfluorobutanesulfonate	29420-49-3		
						~Potassium perfluorooctanesulfonate	2795-39-3		
						Perchlorates			
						~Ammonium Perchlorate	7790-98-9		
						~Lithium Perchlorate	7791-03-9		
						~Perchlorate and Perchlorate Salts	14797-73-0		
						~Potassium Perchlorate	7778-74-7		
						~Sodium Perchlorate	7601-89-0		
						Permethrin	52645-53-1		
6.3E-07	C					Phenacetin	62-44-2	4.5E+02	
						Phenmedipham	13684-63-4		
		2.0E-01	C			Phenol	108-95-2		2.1E+02
						Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1		
						Phenothiazine	92-84-2		
				V		Phenyl Isothiocyanate	103-72-0		
						Phenylenediamine, m-	108-45-2		
						Phenylenediamine, o-	95-54-5		
						Phenylenediamine, p-	106-50-3		
						Phenylphenol, 2-	90-43-7		
						Phorate	298-02-2		
		3.0E-04	I	V		Phosgene	75-44-5		3.1E-01
						Phosmet	732-11-6		
		3.0E-04	I	V		Phosphine	7803-51-2		3.1E-01
		1.0E-02	I			Phosphoric Acid	7664-38-2		1.0E+01
				V		Phosphorus, White	7723-14-0		
						Phthalates			
2.4E-06	C					~Bis(2-ethylhexyl)phthalate	117-81-7	1.2E+02	
						~Butyl Benzyl Phthalate	85-68-7		
						~Butylphthalyl Butylglycolate	85-70-1		
						~Dibutyl Phthalate	84-74-2		
						~Diethyl Phthalate	84-66-2		
				V		~Dimethylterephthalate	120-61-6		
						~Octyl Phthalate, di-N-	117-84-0		
						~Phthalic Acid, p-	100-21-0		
		2.0E-02	C			~Phthalic Anhydride	85-44-9		2.1E+01
						Picloram	1918-02-1		
						Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3		
						Picric Acid (2,4,6-Trinitrophenol)	88-89-1		
						Pirimiphos, Methyl	29232-93-7		
8.6E-03	C					Polybrominated Biphenyls	36355-01-8	3.3E-02	
						Polychlorinated Biphenyls (PCBs)			
2.0E-05	G			V		~Aroclor 1016	12674-11-2	1.4E+01	
5.7E-04	G			V		~Aroclor 1221	11104-28-2	4.9E-01	
5.7E-04	G			V		~Aroclor 1232	11141-16-5	4.9E-01	
5.7E-04	G			V		~Aroclor 1242	53469-21-9	4.9E-01	
5.7E-04	G			V		~Aroclor 1248	12672-29-6	4.9E-01	
5.7E-04	G			V		~Aroclor 1254	11097-69-1	4.9E-01	
5.7E-04	G			V		~Aroclor 1260	11096-82-5	4.9E-01	
				V		~Aroclor 5460	11126-42-4		
1.1E-03	W	1.3E-03	W	V		~Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	2.5E-01	1.4E+00
1.1E-03	W	1.3E-03	W	V		~Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167)	52663-72-6	2.5E-01	1.4E+00
1.1E-03	W	1.3E-03	W	V		~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	2.5E-01	1.4E+00
1.1E-03	W	1.3E-03	W	V		~Hexachlorobiphenyl, 2,3,3',4,4',5- (PCB 156)	38380-08-4	2.5E-01	1.4E+00
1.1E+00	W	1.3E-06	W	V		~Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	2.5E-04	1.4E-03
1.1E-03	W	1.3E-03	W	V		~Pentachlorobiphenyl, 2',3,4,4',5- (PCB 123)	65510-44-3	2.5E-01	1.4E+00
1.1E-03	W	1.3E-03	W	V		~Pentachlorobiphenyl, 2,3',4,4',5- (PCB 118)	31508-00-6	2.5E-01	1.4E+00
1.1E-03	W	1.3E-03	W	V		~Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	2.5E-01	1.4E+00
1.1E-03	W	1.3E-03	W	V		~Pentachlorobiphenyl, 2,3,4,4',5- (PCB 114)	74472-37-0	2.5E-01	1.4E+00

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
3.8E+00	W	4.0E-07	W	V		~Pentachlorobiphenyl, 3,3',4,4',5- (PCB 126)	57465-28-8	7.4E-05	4.2E-04
5.7E-04	I			V		~Polychlorinated Biphenyls (high risk)	1336-36-3	4.9E-01	
1.0E-04	I			V		~Polychlorinated Biphenyls (low risk)	1336-36-3	2.8E+00	
2.0E-05	I			V		~Polychlorinated Biphenyls (lowest risk)	1336-36-3	1.4E+01	
3.8E-03	W	4.0E-04	W			~Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	7.4E-02	4.2E-01
1.1E-02	W	1.3E-04	W	V		~Tetrachlorobiphenyl, 3,4,4',5- (PCB 81)	70362-50-4	2.5E-02	1.4E-01
		6.0E-04	I			Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9		6.3E-01
						Polynuclear Aromatic Hydrocarbons (PAHs)			
				V		~Acenaphthene	83-32-9		
				V		~Anthracene	120-12-7		
6.0E-05	E			V	M	~Benz[a]anthracene	56-55-3	1.7E+00	
		2.0E-06	X			~Benzo(e)pyrene	192-97-2		2.1E-03
1.1E-04	C					~Benzo(j)fluoranthene	205-82-3	2.6E+00	
6.0E-04	I	2.0E-06	I		M	~Benzo[a]pyrene	50-32-8	1.7E-01	2.1E-03
6.0E-05	E				M	~Benzo[b]fluoranthene	205-99-2	1.7E+00	
6.0E-06	E				M	~Benzo[k]fluoranthene	207-08-9	1.7E+01	
				V		~Chloronaphthalene, Beta-	91-58-7		
6.0E-07	E				M	~Chrysene	218-01-9	1.7E+02	
6.0E-04	E				M	~Dibenz[a,h]anthracene	53-70-3	1.7E-01	
1.1E-03	C					~Dibenzo(a,e)pyrene	192-65-4	2.6E-01	
7.1E-02	C				M	~Dimethylbenz(a)anthracene, 7,12-	57-97-6	1.4E-03	
						~Fluoranthene	206-44-0		
				V		~Fluorene	86-73-7		
6.0E-05	E				M	~Indeno[1,2,3-cd]pyrene	193-39-5	1.7E+00	
				V		~Methylnaphthalene, 1-	90-12-0		
				V		~Methylnaphthalene, 2-	91-57-6		
3.4E-05	C	3.0E-03	I	V		~Naphthalene	91-20-3	8.3E+00	3.1E+00
1.1E-04	C					~Nitropyrene, 4-	57835-92-4	2.6E+00	
				V		~Pyrene	129-00-0		
						Prochloraz	67747-09-5		
				V		Profluralin	26399-36-0		
						Prometon	1610-18-0		
						Prometryn	7287-19-6		
						Pronamide	23950-58-5		
						Propachlor	1918-16-7		
						Propanil	709-98-8		
						Propargite	2312-35-8		
				V		Propargyl Alcohol	107-19-7		
						Propazine	139-40-2		
						Propham	122-42-9		
						Propiconazole	60207-90-1		
		8.0E-03	I	V		Propionaldehyde	123-38-6		8.3E+00
		1.0E+00	X	V		Propyl benzene	103-65-1		1.0E+03
		3.0E+00	C	V		Propylene	115-07-1		3.1E+03
						Propylene Glycol	57-55-6		
		2.7E-04	A			Propylene Glycol Dinitrate	6423-43-4		2.8E-01
		2.0E+00	I	V		Propylene Glycol Monomethyl Ether	107-98-2		2.1E+03
3.7E-06	I	3.0E-02	I	V		Propylene Oxide	75-56-9	7.6E+01	3.1E+01
				V		Pyridine	110-86-1		
						Quinalphos	13593-03-8		
						Quinoline	91-22-5		
						Quizalofop-ethyl	76578-14-8		
		3.0E+04	A			Refractory Ceramic Fibers (units in fibers)	E715557		3.1E+04
						Resmethrin	10453-86-8		
				V		Ronnel	299-84-3		
						Rotenone	83-79-4		
6.3E-05	C				M	Safrole	94-59-7	1.6E+00	
						Selenious Acid	7783-00-8		
		2.0E-02	C			Selenium	7782-49-2		2.1E+01
		2.0E-02	C			Selenium Sulfide	7446-34-6		2.1E+01
						Sethoxydim	74051-80-2		
		3.0E-03	C			Silica (crystalline, respirable)	7631-86-9		3.1E+00
						Silver	7440-22-4		
						Simazine	122-34-9		
						Sodium Acifluorfen	62476-59-9		
						Sodium Azide	26628-22-8		
						Sodium Diethyldithiocarbamate	148-18-5		
		1.3E-02	C			Sodium Fluoride	7681-49-4		1.4E+01
						Sodium Fluoroacetate	62-74-8		
						Sodium Metavanadate	13718-26-8		
						Sodium Tungstate	13472-45-2		
						Sodium Tungstate Dihydrate	10213-10-2		
						Stirofos (Tetrachlorovinphos)	961-11-5		
						Strontium, Stable	7440-24-6		
						Strychnine	57-24-9		
		1.0E+00	I	V		Styrene	100-42-5		1.0E+03
						Styrene-Acrylonitrile (SAN) Trimer (THNA isomer)	57964-39-3		
						Styrene-Acrylonitrile (SAN) Trimer (THNP isomer)	57964-40-6		
		2.0E-03	X			Sulfolane	126-33-0		2.1E+00
						Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9		
		1.0E-03	C	V		Sulfur Trioxide	7446-11-9		1.0E+00
		1.0E-03	C			Sulfuric Acid	7664-93-9		1.0E+00
7.1E-06	I					Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl ester	140-57-8	4.0E+01	
						TCMTB	21564-17-0		
						Tebuthiuron	34014-18-1		
						Temephos	3383-96-8		
						Terbacil	5902-51-2		
				V		Terbufos	13071-79-9		
						Terbutryn	886-50-0		
1.3E-06	C			V		Tert-Butyl Acetate	540-88-5	2.2E+02	

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
						Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1		
				V		Tetrachlorobenzene, 1,2,4,5-	95-94-3		
7.4E-06	I			V		Tetrachloroethane, 1,1,1,2-	630-20-6	3.8E+01	
5.8E-05	C			V		Tetrachloroethane, 1,1,2,2-	79-34-5	4.8E+00	
2.6E-07	I	4.0E-02	I	V		Tetrachloroethylene	127-18-4	1.1E+03	4.2E+01
						Tetrachlorophenol, 2,3,4,6-	58-90-2		
				V		Tetrachlorotoluene, p- alpha, alpha, alpha-	5216-25-1		
						Tetraethyl Dithiopyrophosphate	3689-24-5		
		8.0E+01	I	V		Tetrafluoroethane, 1,1,1,2-	811-97-2		8.3E+04
						Tetramethylphosphoramide, -N,N,N',N" (TMPA)	16853-36-4		
						Tetryl (Trinitrophenylmethylnitramine)	479-45-8		
						Thallic Oxide	1314-32-5		
						Thallium (I) Nitrate	10102-45-1		
						Thallium (Soluble Salts)	7440-28-0		
				V		Thallium Acetate	563-68-8		
						Thallium Carbonate	6533-73-9		
						Thallium Chloride	7791-12-0		
						Thallium Selenite	12039-52-0		
						Thallium Sulfate	7446-18-6		
						Thifensulfuron-methyl	79277-27-3		
						Thiobencarb	28249-77-6		
						Thiodiglycol	111-48-8		
						Thiofanox	39196-18-4		
						Thiophanate, Methyl	23564-05-8		
						Thiram	137-26-8		
						Tin	7440-31-5		
		1.0E-04	A	V		Titanium Tetrachloride	7550-45-0		1.0E-01
		5.0E+00	I	V		Toluene	108-88-3		5.2E+03
1.1E-05	C	8.0E-06	C	V		Toluene-2,4-diisocyanate	584-84-9	2.6E+01	8.3E-03
						Toluene-2,5-diamine	95-70-5		
1.1E-05	C	8.0E-06	C	V		Toluene-2,6-diisocyanate	91-08-7	2.6E+01	8.3E-03
						Toluenediamine, 2,3-	2687-25-4		
						Toluenediamine, 3,4-	496-72-0		
						Toluic Acid, p-	99-94-5		
5.1E-05	C					Toluidine, o- (Methylaniline, 2-)	95-53-4	5.5E+00	
						Toluidine, p-	106-49-0		
				V		Total Petroleum Hydrocarbons (Aliphatic High)	E1790670		
		6.0E-01	P	V		Total Petroleum Hydrocarbons (Aliphatic Low)	E1790666		6.3E+02
		1.0E-01	P	V		Total Petroleum Hydrocarbons (Aliphatic Medium)	E1790668		1.0E+02
						Total Petroleum Hydrocarbons (Aromatic High)	E1790676		
		3.0E-02	P	V		Total Petroleum Hydrocarbons (Aromatic Low)	E1790672		3.1E+01
		3.0E-03	P	V		Total Petroleum Hydrocarbons (Aromatic Medium)	E1790674		3.1E+00
3.2E-04	I					Toxaphene	8001-35-2	8.8E-01	
						Toxaphene, Weathered	E1841606		
						Tralomethrin	66841-25-6		
				V		Tri-n-butyltin	688-73-3		
						Triacetin	102-76-1		
						Triadimefon	43121-43-3		
				V		Triallate	2303-17-5		
						Triasulfuron	82097-50-5		
						Tribenuron-methyl	101200-48-0		
				V		Tribromobenzene, 1,2,4-	615-54-3		
						Tribromophenol, 2,4,6-	118-79-6		
						Tribufos	78-48-8		
						Tributyl Phosphate	126-73-8		
						Tributyltin Compounds	E1790679		
						Tributyltin Oxide	56-35-9		
						Trichloramine	10025-85-1		
		5.0E+00	P	V		Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1		5.2E+03
						Trichloroacetic Acid	76-03-9		
						Trichloroaniline HCl, 2,4,6-	33663-50-2		
						Trichloroaniline, 2,4,6-	634-93-5		
				V		Trichlorobenzene, 1,2,3-	87-61-6		
		2.0E-03	P	V		Trichlorobenzene, 1,2,4-	120-82-1		2.1E+00
		5.0E+00	I	V		Trichloroethane, 1,1,1-	71-55-6		5.2E+03
1.6E-05	I	2.0E-04	X	V		Trichloroethane, 1,1,2-	79-00-5	1.8E+01	2.1E-01
4.1E-06	I	2.0E-03	I	V	M	Trichloroethylene	79-01-6	4.8E+01	2.1E+00
				V		Trichlorofluoromethane	75-69-4		
						Trichlorophenol, 2,4,5-	95-95-4		
3.1E-06	I					Trichlorophenol, 2,4,6-	88-06-2	9.1E+01	
						Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5		
						Trichlorophenoxypropionic acid, -2,4,5	93-72-1		
				V		Trichloropropane, 1,1,2-	598-77-6		
		3.0E-04	I	V	M	Trichloropropane, 1,2,3-	96-18-4		3.1E-01
		3.0E-04	P	V		Trichloropropene, 1,2,3-	96-19-5		3.1E-01
						Tricresyl Phosphate (TCP)	1330-78-5		
						Tridiphane	58138-08-2		
		7.0E-03	I	V		Triethylamine	121-44-8		7.3E+00
						Triethylene Glycol	112-27-6		
		2.0E+01	P	V		Trifluoroethane, 1,1,1-	420-46-2		2.1E+04
				V		Trifluralin	1582-09-8		
						Trimethyl Phosphate	512-56-1		
		6.0E-02	I	V		Trimethylbenzene, 1,2,3-	526-73-8		6.3E+01
		6.0E-02	I	V		Trimethylbenzene, 1,2,4-	95-63-6		6.3E+01
		6.0E-02	I	V		Trimethylbenzene, 1,3,5-	108-67-8		6.3E+01
				V		Trimethylpentene, 2,4,4-	25167-70-8		
						Trinitrobenzene, 1,3,5-	99-35-4		
						Trinitrotoluene, 2,4,6-	118-96-7		
						Triphenylphosphine Oxide	791-28-6		
						Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8		

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
6.6E-04	C			V		Tris(1-chloro-2-propyl)phosphate	13674-84-5		
						Tris(2,3-dibromopropyl)phosphate	126-72-7	4.3E-01	
						Tris(2-chloroethyl)phosphate	115-96-8		
						Tris(2-ethylhexyl)phosphate	78-42-2		
						Tungsten	7440-33-7		
		4.0E-05	A			Uranium	7440-61-1		4.2E-02
2.9E-04	C				M	Urethane	51-79-6	3.5E-01	
8.3E-03	P	7.0E-06	P			Vanadium Pentoxide	1314-62-1	3.4E-02	7.3E-03
		1.0E-04	A			Vanadium and Compounds	7440-62-2		1.0E-01
				V		Vernolate	1929-77-7		
						Vinclozolin	50471-44-8		
		2.0E-01	I	V		Vinyl Acetate	108-05-4		2.1E+02
1.5E-05	P	3.0E-03	I	V		Vinyl Bromide	593-60-2	1.9E+01	3.1E+00
4.4E-06	I	8.0E-02	A	V	M	Vinyl Chloride	75-01-4	1.7E+01	8.3E+01
						Warfarin	81-81-2		
		1.0E-01	G	V		Xylene, m-	108-38-3		1.0E+02
		1.0E-01	G	V		Xylene, o-	95-47-6		1.0E+02
		1.0E-01	G	V		Xylene, p-	106-42-3		1.0E+02
		1.0E-01	I	V		Xylenes	1330-20-7		1.0E+02
						Zinc Phosphide	1314-84-7		
						Zinc and Compounds	7440-66-6		
						Zineb	12122-67-7		
						Zirconium	7440-67-7		

RMLs
TR=1E-04
THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
				3.0E-04	O						1.36E+09		1	0.1	Acephate	30560-19-1					3.5E+02	8.3E+02		2.5E+02
		2.2E-06	I			9.0E-03	I	V		1.07E+05	1.36E+09	8.72E+03	1		Acetaldehyde	75-07-0			4.9E+03	4.9E+03			3.4E+02	3.4E+02
				2.0E-02	I						1.36E+09		1	0.1	Acetochlor	34256-82-1					2.3E+04	5.5E+04		1.6E+04
				9.0E-01	I			V		1.14E+05	1.36E+09	1.37E+04	1		Acetone	67-64-1					1.1E+06			1.1E+06
						2.0E-03	X				1.36E+09		1	0.1	Acetone Cyanohydrin	75-86-5							1.2E+07	1.2E+07
						6.0E-02	I	V		1.28E+05	1.36E+09	1.30E+04	1		Acetonitrile	75-05-8							3.4E+03	3.4E+03
				1.0E-01	I			V		2.52E+03	1.36E+09	5.97E+04	1		Acetophenone	98-86-2					1.2E+05			1.2E+05
3.8E+00	C	1.3E-03	C								1.36E+09		1	0.1	Acetylaminofluorene, 2-	53-96-3	8.6E+01	2.0E+02	1.3E+06	6.0E+01				
				5.0E-04	I	2.0E-05	I	V		2.27E+04	1.36E+09	6.91E+03	1		Acrolein	107-02-8					5.8E+02		6.1E-01	6.0E-01
5.0E-01	I	1.0E-04	I	2.0E-03	I	6.0E-03	I		M		1.36E+09		1	0.1	Acrylamide	79-06-1	6.5E+02	1.5E+03	1.7E+07	4.6E+02	2.3E+03	5.5E+03	3.6E+07	1.6E+03
				5.0E-01	I	2.0E-04	P	V		1.09E+05	1.36E+09	9.53E+04	1		Acrylic Acid	79-10-7					5.8E+05		8.3E+01	8.3E+01
5.4E-01	I	6.8E-05	I	1.0E-02	A	2.0E-03	I	V		1.13E+04	1.36E+09	7.69E+03	1		Acrylonitrile	107-13-1	6.1E+02		1.4E+02	1.1E+02	1.2E+04		6.7E+01	6.7E+01
						6.0E-03	P				1.36E+09		1	0.1	Adiponitrile	111-69-3							3.6E+07	3.6E+07
5.6E-02	C			1.0E-02	I						1.36E+09		1	0.1	Alachlor	15972-60-8	5.8E+03	1.4E+04		4.1E+03	1.2E+04	2.8E+04		8.2E+03
				1.0E-03	I						1.36E+09		1	0.1	Aldicarb	116-06-3					1.2E+03	2.8E+03		8.2E+02
				1.0E-03	I						1.36E+09		1	0.1	Aldicarb Sulfone	1646-88-4					1.2E+03	2.8E+03		8.2E+02
											1.36E+09		1	0.1	Aldicarb sulfoxide	1646-87-3								
1.7E+01	I	4.9E-03	I	3.0E-05	I			V			1.36E+09	1.72E+06	1		Aldrin	309-00-2	1.9E+01		4.3E+02	1.8E+01	3.5E+01			3.5E+01
				4.0E-03	P	1.0E-04	X	V		1.11E+05	1.36E+09	3.42E+04	1		Allyl Alcohol	107-18-6					4.7E+03		1.5E+01	1.5E+01
2.1E-02	C	6.0E-06	C			1.0E-03	I	V		1.42E+03	1.36E+09	1.58E+03	1		Allyl Chloride	107-05-1	1.6E+04		3.2E+02	3.2E+02			6.9E+00	6.9E+00
				1.0E+00	P	5.0E-03	P				1.36E+09		1		Aluminum	7429-90-5					1.2E+06		3.0E+07	1.1E+06
				4.0E-04	I						1.36E+09		1		Aluminum Phosphide	20859-73-8					4.7E+02			4.7E+02
				9.0E-03	I						1.36E+09		1	0.1	Ametryn	834-12-8					1.1E+04	2.5E+04		7.4E+03
2.1E+01	C	6.0E-03	C								1.36E+09		1	0.1	Aminobiphenyl, 4-	92-67-1	1.6E+01	3.7E+01	2.8E+05	1.1E+01				
				8.0E-02	P						1.36E+09		1	0.1	Aminophenol, m-	591-27-5					9.3E+04	2.2E+05		6.6E+04
				4.0E-03	X						1.36E+09		1	0.1	Aminophenol, o-	95-55-6					4.7E+03	1.1E+04		3.3E+03
				2.0E-02	P						1.36E+09		1	0.1	Aminophenol, p-	123-30-8					2.3E+04	5.5E+04		1.6E+04
				2.5E-03	I						1.36E+09		1	0.1	Amitraz	33089-61-1					2.9E+03	6.9E+03		2.1E+03
						5.0E-01	I	V					1		Ammonia	7664-41-7								
				2.0E-03	X						1.36E+09		1	0.1	Ammonium Picrate	131-74-8					2.3E+03	5.5E+03		1.6E+03
				2.0E-01	I						1.36E+09		1		Ammonium Sulfamate	7773-06-0					2.3E+05			2.3E+05
						3.0E-03	X	V		1.37E+04	1.36E+09	2.62E+04	1		Amyl Alcohol, tert-	75-85-4							3.4E+02	3.4E+02
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I				1.36E+09		1	0.1	Aniline	62-53-3	5.7E+04	1.4E+05	1.0E+09	4.0E+04	8.2E+03	1.9E+04	6.0E+06	5.7E+03
4.0E-02	P			2.0E-03	X						1.36E+09		1	0.1	Anthraquinone, 9,10-	84-65-1	8.2E+03	1.9E+04		5.7E+03	2.3E+03	5.5E+03		1.6E+03
				4.0E-04	I	3.0E-04	A				1.36E+09		0.15		Antimony (metallic)	7440-36-0					4.7E+02		1.8E+06	4.7E+02
				5.0E-04	H						1.36E+09		0.15		Antimony Pentoxide	1314-60-9					5.8E+02			5.8E+02
				4.0E-04	H						1.36E+09		0.15		Antimony Tetroxide	1332-81-6					4.7E+02			4.7E+02
						2.0E-04	I				1.36E+09		0.15		Antimony Trioxide	1309-64-4							1.2E+06	1.2E+06
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C				1.36E+09		1	0.03	Arsenic, Inorganic	7440-38-2	3.6E+02	1.7E+03	3.9E+05	3.0E+02	5.8E+02	2.8E+03	8.9E+04	4.8E+02
				3.5E-06	C	5.0E-05	I				1.36E+09		1		Arsine	7784-42-1					4.1E+00		3.0E+05	4.1E+00
													1		Asbestos (units in fibers)	1332-21-4								
				3.6E-02	O						1.36E+09		1	0.1	Asulam	3337-71-1					4.2E+04	9.9E+04		3.0E+04
2.3E-01	C			3.0E-03	A						1.36E+09		1	0.1	Atrazine	1912-24-9	1.4E+03	3.4E+03		1.0E+03	3.5E+03	8.3E+03		2.5E+03
8.8E-01	C	2.5E-04	C								1.36E+09		1	0.1	Auramine	492-80-8	3.7E+02	8.8E+02	6.7E+06	2.6E+02				
				4.0E-04	I						1.36E+09		1	0.1	Avermectin B1	65195-55-3					4.7E+02	1.1E+03		3.3E+02
				3.0E-03	A	1.0E-02	A				1.36E+09		1	0.1	Azinphos-methyl	86-50-0					3.5E+03	8.3E+03	6.0E+07	2.5E+03
1.1E-01	I	3.1E-05	I					V			1.36E+09	5.23E+05	1		Azobenzene	103-33-3	3.0E+03		2.1E+04	2.6E+03				
				1.0E+00	P	7.0E-06	P				1.36E+09		1	0.1	Azodicarbonamide	123-77-3					1.2E+06	2.8E+06	4.2E+04	4.0E+04
				2.0E-01	I	5.0E-04	H				1.36E+09		0.07		Barium	7440-39-3					2.3E+05		3.0E+06	2.2E+05
				5.0E-03	O			V			1.36E+09	3.07E+05	1		Benfluralin	1861-40-1					5.8E+03			5.8E+03
				5.0E-02	I						1.36E+09		1	0.1	Benomyl	17804-35-2					5.8E+04	1.4E+05		4.1E+04
				2.0E-01	I						1.36E+09		1	0.1	Bensulfuron-methyl	83055-99-6					2.3E+05	5.5E+05		1.6E+05
				3.0E-02	I						1.36E+09		1	0.1	Bentazon	25057-89-0					3.5E+04	8.3E+04		2.5E+04
4.0E-03	P			1.0E-01	I			V		1.16E+03	1.36E+09	2.25E+04	1		Benzaldehyde	100-52-7	8.2E+04			8.2E+04	1.2E+05			1.2E+05
5.5E-02	I	7.8E-06	I	4.0E-03	I	3.0E-02	I	V		1.82E+03	1.36E+09	3.54E+03	1		Benzene	71-43-2	5.9E+03		5.6E+02	5.1E+02	4.7E+03		4.6E+02	4.2E+02
1.0E-01	X			3.0E-04	X						1.36E+09		1	0.1	Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1	3.3E+03	7.7E+03		2.3E+03	3.5E+02	8.3E+02		2.5E+02
				1.0E-03	P			V		1.26E+03	1.36E+09	1.94E+04	1		Benzenethiol	108-98-5					1.2E+03			1.2E+03
2.3E+02	I	6.7E-02	I	3.0E-03	I				M		1.36E+09		1	0.1	Benzidine	92-87-5	1.4E+00	3.4E+00	2.5E+04	1.0E+00	3.5E+03	8.3E+03		2.5E+03
				4.0E+00	I						1.36E+09		1	0.1	Benzoic Acid	65-85-0					4.7E+06	1.1E+07		3.3E+06
1.3E+01	I							V		3.24E+02	1.36E+09	6.76E+04	1		Benzotrichloride	98-07-7	2.5E+01			2.5E+01				
				1.0E-01	P						1.36E+09		1	0.1	Benzyl Alcohol	100-51-6					1.2E+05	2.8E+05		8.2E+04
1.7E-01	I	4.9E-05	C	2.0E-03	P	1.0E-03	P	V		1.46E+03	1.36E+09	2.55E+04	1		Benzyl Chloride	100-44-7	1.9E+03		6.4E+02	4.8E+02	2.3E+03		1.1E+02	1.1E+02
		2.4E-03	I	2.0E-03	I	2.0E-05	I				1.36E+09		0.007		Beryllium and compounds	7440-41-7			6.9E+05	6.9E+05	2.3E+03		1.2E+05	2.3E+03
				9.0E-03	P						1.36E+09		1	0.1	Bifenox	42576-02-3					1.1E+04	2.5E+04		7.4E+03
				1.5E-02	I						1.36E+09		1	0.1	Biphenthrin	82657-04-3					1.8E+04	4.1E+04		1.2E+04
8.0E-03	I			5.0E-01	I	4.0E-04	X	V			1.36E+09	1.14E+05	1		Biphenyl, 1,1'-	92-52-4	4.1E+04			4.1E+04	5.8E+05		2.0E+02	2.0E+02
				4.0E-02	I			V		1.02E+03	1.36E+09	3.50E+04	1		Bis(2-chloro-1-methylethyl) ether	108-60-1					4.7E+04			4.7E+04
				3.0E-03	P						1.36E+09		1	0.1	Bis(2-chloroethoxy)methane	111-91-1					3.5E+03	8.3E+03		2.5E+03
1.1E+00	I	3.3E-04	I					V		5.05E+03	1.36E+09	4.25E+04	1		Bis(2-chloroethyl)ether	111-44-4	3.0E+02		1.6E+02	1.0E+02				
2.2E+02	I	6.2E-02	I					V		4.22E+03	1.36E+09	1.88E+03	1		Bis(chloromethyl)ether	542-88-1	1.5E+00		3.7E-02	3.6E-02				
				5.0E-02	I						1.36E+09		1	0.1	Bisphenol A	80-05-7					5.8E+04	1.4E+05		4.1E+04
				2.0E-01	I	2.0E-02	H				1.36E+09		1		Boron And Borates Only	7440-42-8					2.3E+05		1.2E+08	2.3E+05
				2.0E+00	P	2.0E-02	P	V			1.36E+09		1		Boron Trichloride	10294-34-5					2.3E+06		1.2E+08	2.3E+06
				4.0E-02	C	1.3E-02	C	V			1.36E+09		1		Boron Trifluoride	7637-07-2					4.7E+04		7.7E+07	4.7E+04
7.0E-01	I			4.0E-03	I						1.36E+09		1		Bromate	15541-45-4	4.7E+02			4.7E+02	4.7E+03			4.7E+03
				1.0E-04	X	6.0E-05	X	V		2.38E+03	1.36E+09	5.92E+03	1		Bromo-2-chloroethane, 1-	107-04-0					1.2E+02		1.6E+00	1.5E+00
				3.0E-04	X			V		8.96E+02	1.36E+09	1.12E+04	1		Bromo-3-fluorobenzene, 1-	1073-06-9					3.5E+02			3.5E+02
				3.0E-04	X			V		3.23E+02	1.36E+09	1.14E+04	1		Bromo-4-fluorobenzene, 1-	460-00-4					3.5E+02			3.5E+02

RMLS
TR=1E-04
THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																		Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
											1.36E+09		1	0.1	Bromoacetic acid	79-08-3								
				8.0E-03	I	6.0E-02	I	V		6.79E+02	1.36E+09	8.37E+03	1		Bromobenzene	108-86-1					9.3E+03		2.2E+03	1.8E+03
						4.0E-02	X	V		4.04E+03	1.36E+09	3.58E+03	1		Bromochloromethane	74-97-5							6.3E+02	6.3E+02
6.2E-02	I	3.7E-05	C	8.0E-03	P			V		9.32E+02	1.36E+09	3.97E+03	1		Bromodichloromethane	75-27-4	5.3E+03		1.3E+02	1.3E+02	9.3E+03			9.3E+03
7.9E-03	I	1.1E-06	I	2.0E-02	I			V		9.15E+02	1.36E+09	9.70E+03	1		Bromoform	75-25-2	4.1E+04		1.1E+04	8.6E+03	2.3E+04			2.3E+04
				1.4E-03	I	5.0E-03	I	V		3.59E+03	1.36E+09	1.40E+03	1		Bromomethane	74-83-9					1.6E+03		3.1E+01	3.0E+01
				5.0E-03	H			V			1.36E+09	1.24E+05	1		Bromophos	2104-96-3					5.8E+03			5.8E+03
						1.0E-01	A	V		9.66E+02	1.36E+09	2.14E+03	1		Bromopropane, 1-	106-94-5							9.4E+02	9.4E+02
1.0E-01	O			1.5E-02	O						1.36E+09		1	0.1	Bromoxynil	1689-84-5	3.2E+03	7.5E+03		2.2E+03	1.8E+04	4.1E+04		1.2E+04
1.0E-01	O			1.5E-02	O			V			1.36E+09	4.74E+05	1		Bromoxynil Octanoate	1689-99-2	3.2E+03			3.2E+03	1.8E+04			1.8E+04
6.0E-01	C	3.0E-05	I			2.0E-03	I	V		6.67E+02	1.36E+09	8.66E+02	1		Butadiene, 1,3-	106-99-0	5.5E+02		3.5E+01	3.3E+01			7.6E+00	7.6E+00
				1.0E-01	I			V		7.64E+03	1.36E+09	3.00E+04	1		Butanol, N-	71-36-3					1.2E+05			1.2E+05
5.0E-04	I			4.0E-01	I	5.0E+00	I	V			1.36E+09	2.87E+04	1		Butyl Alcohol, t-	75-65-0	6.5E+05			6.5E+05	4.7E+05		6.3E+05	2.7E+05
				2.0E+00	P	3.0E+01	P	V		2.13E+04	1.36E+09	2.92E+04	1		Butyl alcohol, sec-	78-92-2					2.3E+06		3.8E+06	1.5E+06
				5.0E-02	I			V			1.36E+09	8.63E+04	1		Butylate	2008-41-5					5.8E+04			5.8E+04
2.0E-04	C	5.7E-08	C								1.36E+09		1	0.1	Butylated hydroxyanisole	25013-16-5	1.6E+06	3.9E+06	2.9E+10	1.1E+06				
3.6E-03	P			3.0E-01	P						1.36E+09		1	0.1	Butylated hydroxytoluene	128-37-0	9.1E+04	2.1E+05		6.4E+04	3.5E+05	8.3E+05		2.5E+05
				5.0E-02	P			V		1.08E+02	1.36E+09	8.14E+03	1		Butylbenzene, n-	104-51-8					5.8E+04			5.8E+04
				1.0E-01	X			V		1.45E+02	1.36E+09	7.35E+03	1		Butylbenzene, sec-	135-98-8					1.2E+05			1.2E+05
				1.0E-01	X			V		1.83E+02	1.36E+09	7.36E+03	1		Butylbenzene, tert-	98-06-6					1.2E+05			1.2E+05
				2.0E-02	A						1.36E+09		1	0.1	Cacodylic Acid	75-60-5					2.3E+04	5.5E+04		1.6E+04
		1.8E-03	I	1.0E-04	A	1.0E-05	A				1.36E+09		0.025	0.001	Cadmium (Diet)	7440-43-9			9.3E+05	9.3E+05	1.2E+02	6.9E+02	6.0E+04	1.0E+02
		1.8E-03	I	1.0E-04	A	1.0E-05	A						0.05	0.001	Cadmium (Water)	7440-43-9								
				5.0E-01	I	2.2E-03	C				1.36E+09		1	0.1	Caprolactam	105-60-2					5.8E+05	1.4E+06	1.3E+07	4.0E+05
1.5E-01	C	4.3E-05	C	2.0E-03	I						1.36E+09		1	0.1	Captafol	2425-06-1	2.2E+03	5.2E+03	3.9E+07	1.5E+03	2.3E+03	5.5E+03		1.6E+03
2.3E-03	C	6.6E-07	C	1.3E-01	I						1.36E+09		1	0.1	Captan	133-06-2	1.4E+05	3.4E+05	2.5E+09	1.0E+05	1.5E+05	3.6E+05		1.1E+05
				1.0E-01	I						1.36E+09		1	0.1	Carbaryl	63-25-2					1.2E+05	2.8E+05		8.2E+04
				5.0E-03	I						1.36E+09		1	0.1	Carbofuran	1563-66-2					5.8E+03	1.4E+04		4.1E+03
				1.0E-01	I	7.0E-01	I	V		7.38E+02	1.36E+09	1.17E+03	1		Carbon Disulfide	75-15-0					1.2E+05		3.6E+03	3.5E+03
7.0E-02	I	6.0E-06	I	4.0E-03	I	1.0E-01	I	V		4.58E+02	1.36E+09	1.49E+03	1		Carbon Tetrachloride	56-23-5	4.7E+03		3.1E+02	2.9E+02	4.7E+03		6.5E+02	5.7E+02
						1.0E-01	P	V		5.89E+03	1.36E+09	6.46E+02	1		Carbonyl Sulfide	463-58-1							2.8E+02	2.8E+02
				1.0E-02	I						1.36E+09		1	0.1	Carbosulfan	55285-14-8					1.2E+04	2.8E+04		8.2E+03
				1.0E-01	I						1.36E+09		1	0.1	Carboxin	5234-68-4					1.2E+05	2.8E+05		8.2E+04
						9.0E-04	I				1.36E+09		1		Ceric oxide	1306-38-3							5.4E+06	5.4E+06
				1.0E-01	I			V			1.36E+09	1.45E+05	1		Chloral Hydrate	302-17-0					1.2E+05			1.2E+05
				1.5E-02	I						1.36E+09		1	0.1	Chloramben	133-90-4					1.8E+04	4.1E+04		1.2E+04
											1.36E+09		1	0.1	Chloramines, Organic	E701235								
4.0E-01	H										1.36E+09		1	0.1	Chloranil	118-75-2	8.1E+02	1.9E+03		5.7E+02				
				5.0E-04	G			V			1.36E+09	1.49E+06	1	0.04	Chlordane (alpha)	5103-71-9					5.8E+02	3.4E+03		5.0E+02
				5.0E-04	G			V			1.36E+09	1.49E+06	1	0.04	Chlordane (gamma)	5103-74-2					5.8E+02	3.4E+03		5.0E+02
3.5E-01	I	1.0E-04	I	5.0E-04	I	7.0E-04	I	V			1.36E+09	1.53E+06	1	0.04	Chlordane (technical mixture)	12789-03-6	9.3E+02	5.5E+03	1.9E+04	7.7E+02	5.8E+02	3.4E+03	4.7E+03	4.5E+02
1.0E+01	I	4.6E-03	C	3.0E-04	I						1.36E+09		1	0.1	Chlordecone (Kepone)	143-50-0	3.3E+01	7.7E+01	3.6E+05	2.3E+01	3.5E+02	8.3E+02		2.5E+02
				7.0E-04	A						1.36E+09		1	0.1	Chlorfenvinphos	470-90-6					8.2E+02	1.9E+03		5.7E+02
				9.0E-02	O						1.36E+09		1	0.1	Chlorimuron, Ethyl-	90982-32-4					1.1E+05	2.5E+05		7.4E+04
				1.0E-01	I	1.5E-04	A	V		2.78E+03	1.36E+09	1.22E+03	1		Chlorine	7782-50-5					1.2E+05		7.8E-01	7.8E-01
				3.0E-02	I	2.0E-04	I	V			1.36E+09		1		Chlorine Dioxide	10049-04-4					3.5E+04		1.2E+06	3.4E+04
				3.0E-02	I						1.36E+09		1		Chlorite (Sodium Salt)	7758-19-2					3.5E+04			3.5E+04
						5.0E+01	I	V		1.15E+03	1.36E+09	1.03E+03	1		Chloro-1,1-difluoroethane, 1-	75-68-3							2.3E+05	2.3E+05
		3.0E-04	I	2.0E-02	H	2.0E-02	I	V		7.86E+02	1.36E+09	1.08E+03	1		Chloro-1,3-butadiene, 2-	126-99-8			4.4E+00	4.4E+00	2.3E+04		9.4E+01	9.4E+01
4.6E-01	H										1.36E+09		1	0.1	Chloro-2-methylaniline HCl, 4-	3165-93-3	7.1E+02	1.7E+03		5.0E+02				
1.0E-01	P	7.7E-05	C	3.0E-03	X						1.36E+09		1	0.1	Chloro-2-methylaniline, 4-	95-69-2	3.3E+03	7.7E+03	2.2E+07	2.3E+03	3.5E+03	8.3E+03		2.5E+03
2.7E-01	X							V		1.18E+04	1.36E+09	1.62E+04	1		Chloroacetaldehyde, 2-	107-20-0	1.2E+03			1.2E+03				
											1.36E+09		1	0.1	Chloroacetic Acid	79-11-8								
						3.0E-05	I				1.36E+09		1	0.1	Chloroacetophenone, 2-	532-27-4							1.8E+05	1.8E+05
2.0E-01	P			5.0E-04	P						1.36E+09		1	0.1	Chloroaniline, p-	106-47-8	1.6E+03	3.9E+03		1.1E+03	5.8E+02	1.4E+03		4.1E+02
				2.0E-02	I	5.0E-02	P	V		7.61E+02	1.36E+09	6.45E+03	1		Chlorobenzene	108-90-7					2.3E+04		1.4E+03	1.3E+03
				1.0E-01	X						1.36E+09		1	0.1	Chlorobenzene sulfonic acid, p-	98-66-8					1.2E+05	2.8E+05		8.2E+04
1.1E-01	C	3.1E-05	C	2.0E-02	I						1.36E+09		1	0.1	Chlorobenzilate	510-15-6	3.0E+03	7.0E+03	5.4E+07	2.1E+03	2.3E+04	5.5E+04		1.6E+04
				3.0E-02	X						1.36E+09		1	0.1	Chlorobenzoic Acid, p-	74-11-3					3.5E+04	8.3E+04		2.5E+04
		8.6E-06	C	3.0E-03	P	3.0E-01	P	V		2.90E+02	1.36E+09	6.76E+03	1		Chlorobenzotrifluoride, 4-	98-56-6			9.6E+02	9.6E+02	3.5E+03		8.9E+03	2.5E+03
				4.0E-02	P			V		7.28E+02	1.36E+09	1.76E+03	1		Chlorobutane, 1-	109-69-3					4.7E+04			4.7E+04
						5.0E+01	I	V		1.68E+03	1.36E+09	9.38E+02	1		Chlorodifluoromethane	75-45-6							2.1E+05	2.1E+05
				2.0E-02	P			V		1.11E+05	1.36E+09	7.81E+04	1		Chloroethanol, 2-	107-07-3					2.3E+04			2.3E+04
3.1E-02	C	2.3E-05	I	1.0E-02	I	9.8E-02	A	V		2.54E+03	1.36E+09	2.63E+03	1		Chloroform	67-66-3	1.1E+04		1.4E+02	1.4E+02	1.2E+04		1.1E+03	1.0E+03
						9.0E-02	I	V		1.32E+03	1.36E+09	1.18E+03	1		Chloromethane	74-87-3							4.6E+02	4.6E+02
2.4E+00	C	6.9E-04	C					V		9.32E+03	1.36E+09	5.33E+03	1		Chloromethyl Methyl Ether	107-30-2	1.4E+02		9.5E+00	8.9E+00				
3.0E-01	P			3.0E-03	P	1.0E-05	X				1.36E+09		1	0.1	Chloronitrobenzene, o-	88-73-3	1.1E+03	2.6E+03		7.7E+02	3.5E+03	8.3E+03	6.0E+04	2.4E+03
6.0E-02	P			7.0E-04	P	2.0E-03	P				1.36E+09		1	0.1	Chloronitrobenzene, p-	100-00-5	5.5E+03	1.3E+04		3.8E+03	8.2E+02	1.9E+03	1.2E+07	5.7E+02
				5.0E-03	I			V		2.74E+04	1.36E+09	1.39E+05	1		Chlorophenol, 2-	95-57-8					5.8E+03			5.8E+03
						4.0E-04	C	V		6.17E+02	1.36E+09	4.68E+03	1		Chloropicrin	76-06-2							8.2E+00	8.2E+00
1.7E-02	C			1.5E-02	I						1.36E+09		1	0.1	Chlorothalonil	1897-45-6	1.9E+04	4.5E+04		1.4E+04	1.8E+04	4.1E+04		1.2E+04
				2.0E-02	I			V		9.07E+02	1.36E+09	8.12E+03	1		Chlorotoluene, o-	95-49-8					2.3E+04			2.3E+04
				2.0E-02	X			V		2.53E+02	1.36E+09	7.29E+03	1		Chlorotoluene, p-	106-43-4					2.3E+04			2.3E+04
2.4E+02	C	6.9E-02	C								1.36E+09		1	0.1	Chlorozotocin	54749-90-5	1.4E+00	3.2E+00	2.4E+04	9.6E-01				
				5.0E-03	O						1.36E+09		1	0.1	Chlorpropham	101-21-3					5.8E+03	1.4E+04		4.1E+03
				1.0E-03	A						1.36E+09		1	0.1	Chlorpyrifos	2921-88-2					1.2E+03	2.8E+03		8.2E+02
				1.0E-02	H						1.36E+09		1	0.1	Chlorpyrifos Methyl	5598-13-0					1.2E+04	2.8E+04		8.2E+03
				5.0E-02	O						1.36E+09		1	0.1	Chlorsulfuron	64902-72-3					5.8E+04	1.4E+05		4.1E+04

RMLS
TR=1E-04
THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																		Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
				1.0E-02	I						1.36E+09		1	0.1	Chlorthal-dimethyl	1861-32-1					1.2E+04	2.8E+04		8.2E+03
				8.0E-04	H						1.36E+09		1	0.1	Chlorthiophos	60238-56-4					9.3E+02	2.2E+03		6.6E+02
				1.5E+00	I						1.36E+09		0.013		Chromium(III), Insoluble Salts	16065-83-1					1.8E+06			1.8E+06
5.0E-01	C	8.4E-02	G	3.0E-03	I	1.0E-04	I		M		1.36E+09		0.025		Chromium(VI)	18540-29-9	6.5E+02		2.0E+04	6.3E+02	3.5E+03		6.0E+05	3.5E+03
											1.36E+09		0.013		Chromium, Total	7440-47-3								
				1.3E-02	I						1.36E+09		1	0.1	Clofentezine	74115-24-5					1.5E+04	3.6E+04		1.1E+04
		9.0E-03	P	3.0E-04	P	6.0E-06	P				1.36E+09		1		Cobalt	7440-48-4			1.9E+05	1.9E+05	3.5E+02		3.6E+04	3.5E+02
		6.2E-04	I					V	M				1		Coke Oven Emissions	E649830								
				4.0E-02	H						1.36E+09		1		Copper	7440-50-8					4.7E+04			4.7E+04
				5.0E-02	I	6.0E-01	C				1.36E+09		1	0.1	Cresol, m-	108-39-4					5.8E+04	1.4E+05	3.6E+09	4.1E+04
				5.0E-02	I	6.0E-01	C				1.36E+09		1	0.1	Cresol, o-	95-48-7					5.8E+04	1.4E+05	3.6E+09	4.1E+04
				2.0E-02	P	6.0E-01	C				1.36E+09		1	0.1	Cresol, p-	106-44-5					2.3E+04	5.5E+04	3.6E+09	1.6E+04
				1.0E-01	A						1.36E+09		1	0.1	Cresol, p-chloro-m-	59-50-7					1.2E+05	2.8E+05		8.2E+04
				1.0E-01	A	6.0E-01	C				1.36E+09		1	0.1	Cresols	1319-77-3					1.2E+05	2.8E+05	3.6E+09	8.2E+04
1.9E+00	H			1.0E-03	P			V		1.66E+04	1.36E+09	1.89E+04	1		Crotonaldehyde, trans-	123-73-9	1.7E+02			1.7E+02	1.2E+03			1.2E+03
				1.0E-01	I	4.0E-01	I	V		2.68E+02	1.36E+09	6.21E+03	1		Cumene	98-82-8					1.2E+05		1.1E+04	9.9E+03
2.2E-01	C	6.3E-05	C								1.36E+09		1	0.1	Cupferron	135-20-6	1.5E+03	3.5E+03	2.6E+07	1.0E+03				
8.4E-01	H			2.0E-03	H						1.36E+09		1	0.1	Cyanazine	21725-46-2	3.9E+02	9.2E+02		2.7E+02	2.3E+03	5.5E+03		1.6E+03
															Cyanides									
				1.0E-03	I						1.36E+09		1		~Calcium Cyanide	592-01-8					1.2E+03			1.2E+03
				5.0E-03	I						1.36E+09		1		~Copper Cyanide	544-92-3					5.8E+03			5.8E+03
				6.0E-04	I	8.0E-04	G	V		9.54E+05	1.36E+09	5.33E+04	1		~Cyanide (CN-)	57-12-5					7.0E+02		1.9E+02	1.5E+02
				1.0E-03	I			V			1.36E+09		1		~Cyanogen	460-19-5					1.2E+03			1.2E+03
				9.0E-02	I			V			1.36E+09		1		~Cyanogen Bromide	506-68-3					1.1E+05			1.1E+05
				5.0E-02	I			V			1.36E+09		1		~Cyanogen Chloride	506-77-4					5.8E+04			5.8E+04
				6.0E-04	I	8.0E-04	I	V		1.00E+07	1.36E+09	5.22E+04	1		~Hydrogen Cyanide	74-90-8					7.0E+02		1.8E+02	1.5E+02
				2.0E-03	I						1.36E+09		1		~Potassium Cyanide	151-50-8					2.3E+03			2.3E+03
				5.0E-03	I						1.36E+09		0.04		~Potassium Silver Cyanide	506-61-6					5.8E+03			5.8E+03
				1.0E-01	I						1.36E+09		0.04		~Silver Cyanide	506-64-9					1.2E+05			1.2E+05
				1.0E-03	I						1.36E+09		1		~Sodium Cyanide	143-33-9					1.2E+03			1.2E+03
				2.0E-04	P						1.36E+09		1		~Thiocyanates	E1790665					2.3E+02			2.3E+02
				2.0E-04	X			V			1.36E+09		1		~Thiocyanic Acid	463-56-9					2.3E+02			2.3E+02
				5.0E-02	I						1.36E+09		1		~Zinc Cyanide	557-21-1					5.8E+04			5.8E+04
						6.0E+00	I	V		1.17E+02	1.36E+09	1.04E+03	1		Cyclohexane	110-82-7							2.7E+04	2.7E+04
2.0E-02	X			2.0E-02	X						1.36E+09		1	0.1	Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	1.6E+04	3.9E+04		1.1E+04	2.3E+04	5.5E+04		1.6E+04
				5.0E+00	I	7.0E-01	P	V		5.11E+03	1.36E+09	4.17E+04	1		Cyclohexanone	108-94-1					5.8E+06		1.3E+05	1.3E+05
				5.0E-03	P	1.0E+00	X	V		2.83E+02	1.36E+09	1.46E+03	1		Cyclohexene	110-83-8					5.8E+03		6.4E+03	3.1E+03
				2.0E-01	I			V		2.93E+05	1.36E+09	7.46E+04	1		Cyclohexylamine	108-91-8					2.3E+05			2.3E+05
				2.5E-02	I						1.36E+09		1	0.1	Cyfluthrin	68359-37-5					2.9E+04	6.9E+04		2.1E+04
				7.2E-02	O						1.36E+09		1	0.1	Cypermethrin	52315-07-8					8.4E+04	2.0E+05		5.9E+04
				5.0E-01	O						1.36E+09		1	0.1	Cyromazine	66215-27-8					5.8E+05	1.4E+06		4.1E+05
2.4E-01	I	6.9E-05	C	3.0E-05	X						1.36E+09		1	0.1	DDD, p,p`- (DDD)	72-54-8	1.4E+03	3.2E+03	2.4E+07	9.6E+02	3.5E+01	8.3E+01		2.5E+01
3.4E-01	I	9.7E-05	C	3.0E-04	X			V			1.36E+09	2.10E+06	1		DDE, p,p'-	72-55-9	9.6E+02		2.7E+04	9.3E+02	3.5E+02			3.5E+02
3.4E-01	I	9.7E-05	I	5.0E-04	I						1.36E+09		1	0.03	DDT	50-29-3	9.6E+02	7.6E+03	1.7E+07	8.5E+02	5.8E+02	4.6E+03		5.2E+02
				3.0E-02	I						1.36E+09		1	0.1	Dalapon	75-99-0					3.5E+04	8.3E+04		2.5E+04
1.8E-02	C	5.1E-06	C	1.5E-01	I						1.36E+09		1	0.1	Daminozide	1596-84-5	1.8E+04	4.3E+04	3.3E+08	1.3E+04	1.8E+05	4.1E+05		1.2E+05
7.0E-04	I			7.0E-03	I						1.36E+09		1	0.1	Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5	4.7E+05	1.1E+06		3.3E+05	8.2E+03	1.9E+04		5.7E+03
				4.0E-05	I						1.36E+09		1	0.1	Demeton	8065-48-3					4.7E+01	1.1E+02		3.3E+01
1.2E-03	I			6.0E-01	I						1.36E+09		1	0.1	Di(2-ethylhexyl)adipate	103-23-1	2.7E+05	6.4E+05		1.9E+05	7.0E+05	1.7E+06		4.9E+05
6.1E-02	H										1.36E+09		1	0.1	Diallate	2303-16-4	5.4E+03	1.3E+04		3.8E+03				
				7.0E-04	A						1.36E+09		1	0.1	Diazinon	333-41-5					8.2E+02	1.9E+03		5.7E+02
				1.0E-02	X			V			1.36E+09	5.24E+05	1		Dibenzothiophene	132-65-0					1.2E+04			1.2E+04
8.0E-01	P	6.0E-03	P	2.0E-04	P	2.0E-04	I	V	M	9.79E+02	1.36E+09	3.20E+04	1		Dibromo-3-chloropropane, 1,2-	96-12-8	4.1E+02		6.5E+00	6.4E+00	2.3E+02		2.8E+01	2.5E+01
											1.36E+09		1	0.1	Dibromoacetic acid	631-64-1								
				4.0E-04	X			V		1.59E+02	1.36E+09	1.93E+04	1		Dibromobenzene, 1,3-	108-36-1					4.7E+02			4.7E+02
				1.0E-02	I			V			1.36E+09	2.20E+04	1		Dibromobenzene, 1,4-	106-37-6					1.2E+04			1.2E+04
8.4E-02	I			2.0E-02	I			V		8.02E+02	1.36E+09	7.95E+03	1		Dibromochloromethane	124-48-1	3.9E+03			3.9E+03	2.3E+04			2.3E+04
2.0E+00	I	6.0E-04	I	9.0E-03	I	9.0E-03	I	V		1.34E+03	1.36E+09	8.64E+03	1		Dibromoethane, 1,2-	106-93-4	1.6E+02		1.8E+01	1.6E+01	1.1E+04		3.4E+02	3.3E+02
						4.0E-03	X	V		2.82E+03	1.36E+09	5.64E+03	1		Dibromomethane (Methylene Bromide)	74-95-3							9.9E+01	9.9E+01
				3.0E-04	P						1.36E+09		1	0.1	Dibutyltin Compounds	E1790661					3.5E+02	8.3E+02		2.5E+02
				3.0E-02	I						1.36E+09		1	0.1	Dicamba	1918-00-9					3.5E+04	8.3E+04		2.5E+04
											1.36E+09		1	0.1	Dichloramine	3400-09-7								
		4.2E-03	P					V		5.54E+02	1.36E+09	3.21E+03	1		Dichloro-2-butene, 1,4-	764-41-0			9.4E-01	9.4E-01				
		4.2E-03	P					V		5.19E+02	1.36E+09	1.11E+04	1		Dichloro-2-butene, cis-1,4-	1476-11-5			3.2E+00	3.2E+00				
		4.2E-03	P					V		7.60E+02	1.36E+09	1.11E+04	1		Dichloro-2-butene, trans-1,4-	110-57-6			3.2E+00	3.2E+00				
5.0E-02	I			4.0E-03	I						1.36E+09		1	0.1	Dichloroacetic Acid	79-43-6	6.5E+03	1.5E+04		4.6E+03	4.7E+03	1.1E+04		3.3E+03
				9.0E-02	I	2.0E-01	H	V		3.76E+02	1.36E+09	1.17E+04	1		Dichlorobenzene, 1,2-	95-50-1					1.1E+05		1.0E+04	9.3E+03
5.4E-03	C	1.1E-05	C	7.0E-02	A	8.0E-01	I	V			1.36E+09	1.04E+04	1		Dichlorobenzene, 1,4-	106-46-7	6.1E+04		1.2E+03	1.1E+03	8.2E+04		3.7E+04	2.5E+04
4.5E-01	I	3.4E-04	C								1.36E+09		1	0.1	Dichlorobenzidine, 3,3'-	91-94-1	7.3E+02	1.7E+03	4.9E+06	5.1E+02				
				9.0E-03	X						1.36E+09		1	0.1	Dichlorobenzophenone, 4,4'-	90-98-2					1.1E+04	2.5E+04		7.4E+03
				2.0E-01	I	1.0E-01	X	V		8.45E+02	1.36E+09	8.41E+02	1		Dichlorodifluoromethane	75-71-8					2.3E+05		3.7E+02	3.7E+02
5.7E-03	C	1.6E-06	C	2.0E-01	P			V		1.69E+03	1.36E+09	2.08E+03	1		Dichloroethane, 1,1-	75-34-3	5.7E+04		1.6E+03	1.6E+03	2.3E+05			2.3E+05
9.1E-02	I	2.6E-05	I	6.0E-03	X	7.0E-03	P	V		2.98E+03	1.36E+09	4.57E+03	1		Dichloroethane, 1,2-	107-06-2	3.6E+03		2.2E+02	2.0E+02	7.0E+03		1.4E+02	1.4E+02
				5.0E-02	I	2.0E-01	I	V		1.19E+03	1.36E+09	1.16E+03	1		Dichloroethylene, 1,1-	75-35-4					5.8E+04		1.0E+03	1.0E+03
				2.0E-03	I			V		2.37E+03	1.36E+09	2.50E+03	1		Dichloroethylene, cis-1,2-	156-59-2					2.3E+03			2.3E+03
				2.0E-02	I	4.0E-02	X	V		1.85E+03	1.36E+09	1.75E+03	1		Dichloroethylene, trans-1,2-	156-60-5					2.3E+04		3.1E+02	3.0E+02
				3.0E-03	I						1.36E+09		1	0.1	Dichlorophenol, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03
				1.0E-02	I						1.36E+09		1	0.05	Dichlorophenoxy Acetic Acid, 2,4-	94-75-7					1.2E+04	5.5E+04		9.6E+03

RMLS
TR=1E-04
THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																		Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
3.7E-02	P	3.7E-06	P	4.0E-02	P	4.0E-03	I	V		1.36E+03	1.36E+09	3.79E+03	1		Dichloropropane, 1,2-	78-87-5	8.8E+03		1.3E+03	1.1E+03	4.7E+04		6.6E+01	6.6E+01
				2.0E-02	P			V		1.49E+03	1.36E+09	6.76E+03	1		Dichloropropane, 1,3-	142-28-9					2.3E+04			2.3E+04
				3.0E-03	I						1.36E+09		1	0.1	Dichloropropanol, 2,3-	616-23-9					3.5E+03	8.3E+03		2.5E+03
1.0E-01	I	4.0E-06	I	3.0E-02	I	2.0E-02	I	V		1.57E+03	1.36E+09	3.55E+03	1		Dichloropropene, 1,3-	542-75-6	3.3E+03		1.1E+03	8.2E+02	3.5E+04		3.1E+02	3.1E+02
2.9E-01	I	8.3E-05	C	5.0E-04	I	5.0E-04	I				1.36E+09		1	0.1	Dichlorvos	62-73-7	1.1E+03	2.7E+03	2.0E+07	7.9E+02	5.8E+02	1.4E+03	3.0E+06	4.1E+02
				3.0E-05	O						1.36E+09		1	0.1	Dicrotophos	141-66-2					3.5E+01	8.3E+01		2.5E+01
				8.0E-02	P	3.0E-04	X	V		2.56E+02	1.36E+09	4.11E+03	1		Dicyclopentadiene	77-73-6					9.3E+04		5.4E+00	5.4E+00
1.6E+01	I	4.6E-03	I	5.0E-05	I						1.36E+09		1	0.1	Dieldrin	60-57-1	2.0E+01	4.8E+01	3.6E+05	1.4E+01	5.8E+01	1.4E+02		4.1E+01
		3.0E-04	C			5.0E-03	I						1	0.1	Diesel Engine Exhaust	E17136615								
				2.0E-03	P	2.0E-04	P				1.36E+09		1	0.1	Diethanolamine	111-42-2					2.3E+03	5.5E+03	1.2E+06	1.6E+03
				3.0E-02	P	1.0E-04	P				1.36E+09		1	0.1	Diethylene Glycol Monobutyl Ether	112-34-5					3.5E+04	8.3E+04	6.0E+05	2.4E+04
				6.0E-02	P	3.0E-04	P				1.36E+09		1	0.1	Diethylene Glycol Monoethyl Ether	111-90-0					7.0E+04	1.7E+05	1.8E+06	4.8E+04
				1.0E-03	P			V		1.12E+05	1.36E+09	1.39E+05	1		Diethylformamide	617-84-5					1.2E+03			1.2E+03
3.5E+02	C	1.0E-01	C								1.36E+09		1	0.1	Diethylstilbestrol	56-53-1	9.3E-01	2.2E+00	1.7E+04	6.6E-01				
				8.3E-02	O						1.36E+09		1	0.1	Difenzoquat	43222-48-6					9.7E+04	2.3E+05		6.8E+04
				2.0E-02	I						1.36E+09		1	0.1	Diflubenzuron	35367-38-5					2.3E+04	5.5E+04		1.6E+04
						4.0E+01	I	V		1.43E+03	1.36E+09	1.15E+03	1		Difluoroethane, 1,1-	75-37-6							2.0E+05	2.0E+05
						3.0E+01	X	V		6.91E+02	1.36E+09	7.58E+02	1		Difluoropropane, 2,2-	420-45-1							1.0E+05	1.0E+05
4.4E-02	C	1.3E-05	C					V			1.36E+09	1.23E+05	1		Dihydrosafrole	94-58-6	7.4E+03		1.2E+04	4.5E+03				
						7.0E-01	P	V		2.26E+03	1.36E+09	3.06E+03	1		Diisopropyl Ether	108-20-3							9.4E+03	9.4E+03
				8.0E-02	I			V		5.30E+02	1.36E+09	3.81E+04	1		Diisopropyl Methylphosphonate	1445-75-6					9.3E+04			9.3E+04
				2.2E-02	O						1.36E+09		1	0.1	Dimethipin	55290-64-7					2.5E+04	6.0E+04		1.8E+04
				2.2E-03	O						1.36E+09		1	0.1	Dimethoate	60-51-5					2.6E+03	6.1E+03		1.8E+03
1.6E+00	P										1.36E+09		1	0.1	Dimethoxybenzidine, 3,3'-	119-90-4	2.0E+02	4.8E+02		1.4E+02				
1.7E-03	P			6.0E-02	P						1.36E+09		1	0.1	Dimethyl methylphosphonate	756-79-6	1.9E+05	4.5E+05		1.4E+05	7.0E+04	1.7E+05		4.9E+04
4.6E+00	C	1.3E-03	C								1.36E+09		1	0.1	Dimethylamino azobenzene [p-]	60-11-7	7.1E+01	1.7E+02	1.3E+06	5.0E+01				
5.8E-01	H										1.36E+09		1	0.1	Dimethylaniline HCl, 2,4-	21436-96-4	5.6E+02	1.3E+03		4.0E+02				
2.0E-01	P			2.0E-03	X						1.36E+09		1	0.1	Dimethylaniline, 2,4-	95-68-1	1.6E+03	3.9E+03		1.1E+03	2.3E+03	5.5E+03		1.6E+03
2.7E-02	P			2.0E-03	I			V		8.30E+02	1.36E+09	3.13E+04	1		Dimethylaniline, N,N-	121-69-7	1.2E+04			1.2E+04	2.3E+03			2.3E+03
1.1E+01	P										1.36E+09		1	0.1	Dimethylbenzidine, 3,3'-	119-93-7	3.0E+01	7.0E+01		2.1E+01				
				1.0E-01	P	3.0E-02	I	V		1.06E+05	1.36E+09	1.28E+05	1		Dimethylformamide	68-12-2					1.2E+05		1.7E+04	1.5E+04
				1.0E-04	X	2.0E-06	X	V		1.72E+05	1.36E+09	2.77E+04	1		Dimethylhydrazine, 1,1-	57-14-7					1.2E+02		2.4E-01	2.4E-01
5.5E+02	C	1.6E-01	C					V		1.89E+05	1.36E+09	1.68E+05	1		Dimethylhydrazine, 1,2-	540-73-8	5.9E-01		1.3E+00	4.1E-01				
				2.0E-02	I						1.36E+09		1	0.1	Dimethylphenol, 2,4-	105-67-9					2.3E+04	5.5E+04		1.6E+04
				6.0E-04	I						1.36E+09		1	0.1	Dimethylphenol, 2,6-	576-26-1					7.0E+02	1.7E+03		4.9E+02
				1.0E-03	I						1.36E+09		1	0.1	Dimethylphenol, 3,4-	95-65-8					1.2E+03	2.8E+03		8.2E+02
4.5E-02	C	1.3E-05	C					V		4.73E+02	1.36E+09	5.48E+03	1		Dimethylvinylchloride	513-37-1	7.3E+03		5.2E+02	4.8E+02				
				8.0E-05	X						1.36E+09		1	0.1	Dinitro-o-cresol, 4,6-	534-52-1					9.3E+01	2.2E+02		6.6E+01
				2.0E-03	I						1.36E+09		1	0.1	Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5					2.3E+03	5.5E+03		1.6E+03
				4.0E-04	X	2.0E-03	X				1.36E+09		1	0.1	Dinitroaniline, 3,5-	618-87-1					4.7E+02	1.1E+03	1.2E+07	3.3E+02
				1.0E-04	P						1.36E+09		1	0.1	Dinitrobenzene, 1,2-	528-29-0					1.2E+02	2.8E+02		8.2E+01
				1.0E-04	I						1.36E+09		1	0.1	Dinitrobenzene, 1,3-	99-65-0					1.2E+02	2.8E+02		8.2E+01
				1.0E-04	P						1.36E+09		1	0.1	Dinitrobenzene, 1,4-	100-25-4					1.2E+02	2.8E+02		8.2E+01
				2.0E-03	I						1.36E+09		1	0.1	Dinitrophenol, 2,4-	51-28-5					2.3E+03	5.5E+03		1.6E+03
6.8E-01	I										1.36E+09		1	0.1	Dinitrotoluene Mixture, 2,4/2,6-	E1615210	4.8E+02	1.1E+03		3.4E+02				
3.1E-01	C	8.9E-05	C	2.0E-03	I						1.36E+09		1	0.102	Dinitrotoluene, 2,4-	121-14-2	1.1E+03	2.4E+03	1.9E+07	7.4E+02	2.3E+03	5.4E+03		1.6E+03
1.5E+00	P			3.0E-04	X						1.36E+09		1	0.099	Dinitrotoluene, 2,6-	606-20-2	2.2E+02	5.2E+02		1.5E+02	3.5E+02	8.4E+02		2.5E+02
				1.0E-04	X						1.36E+09		1	0.006	Dinitrotoluene, 2-Amino-4,6-	35572-78-2					1.2E+02	4.6E+03		1.1E+02
				1.0E-04	X						1.36E+09		1	0.009	Dinitrotoluene, 4-Amino-2,6-	19406-51-0					1.2E+02	3.1E+03		1.1E+02
4.5E-01	X			9.0E-04	X						1.36E+09		1	0.1	Dinitrotoluene, Technical grade	25321-14-6	7.3E+02	1.7E+03		5.1E+02	1.1E+03	2.5E+03		7.4E+02
				1.0E-03	I						1.36E+09		1	0.1	Dinoseb	88-85-7					1.2E+03	2.8E+03		8.2E+02
1.0E-01	I	5.0E-06	I	3.0E-02	I	3.0E-02	I	V		1.16E+05	1.36E+09	3.96E+04	1		Dioxane, 1,4-	123-91-1	3.3E+03		9.7E+03	2.4E+03	3.5E+04		5.2E+03	4.5E+03
															Dioxins									
6.2E+03	I	1.3E+00	I								1.36E+09		1	0.03	~Hexachlorodibenzo-p-dioxin, Mixture	34465-46-8	5.3E-02	4.2E-01	1.3E+03	4.7E-02				
1.3E+05	C	3.8E+01	C	7.0E-10	I	4.0E-08	C	V			1.36E+09	1.96E+06	1	0.03	~TCDD, 2,3,7,8-	1746-01-6	2.5E-03	2.0E-02	6.3E-02	2.2E-03	8.2E-04	6.4E-03	3.4E-01	7.2E-04
				3.0E-02	I						1.36E+09		1	0.1	Diphenamid	957-51-7					3.5E+04	8.3E+04		2.5E+04
						4.0E-04	X	V			1.36E+09	8.06E+04	1		Diphenyl Ether	101-84-8							1.4E+02	1.4E+02
				8.0E-04	X						1.36E+09		1	0.1	Diphenyl Sulfone	127-63-9					9.3E+02	2.2E+03		6.6E+02
				1.0E-01	O						1.36E+09		1	0.1	Diphenylamine	122-39-4					1.2E+05	2.8E+05		8.2E+04
8.0E-01	I	2.2E-04	I								1.36E+09		1	0.1	Diphenylhydrazine, 1,2-	122-66-7	4.1E+02	9.7E+02	7.6E+06	2.9E+02				
				2.2E-03	I						1.36E+09		1	0.1	Diquat	2764-72-9					2.6E+03	6.1E+03		1.8E+03
7.4E+00	C	2.1E-03	C								1.36E+09		1	0.1	Direct Black 38	1937-37-7	4.4E+01	1.0E+02	7.9E+05	3.1E+01				
7.4E+00	C	2.1E-03	C								1.36E+09		1	0.1	Direct Blue 6	2602-46-2	4.4E+01	1.0E+02	7.9E+05	3.1E+01				
6.7E+00	C	1.9E-03	C								1.36E+09		1	0.1	Direct Brown 95	16071-86-6	4.9E+01	1.2E+02	8.8E+05	3.4E+01				
				4.0E-05	I						1.36E+09		1	0.1	Disulfoton	298-04-4					4.7E+01	1.1E+02		3.3E+01
				1.0E-02	I			V			1.36E+09	4.54E+04	1		Dithiane, 1,4-	505-29-3					1.2E+04			1.2E+04
				2.0E-03	I						1.36E+09		1	0.1	Diuron	330-54-1					2.3E+03	5.5E+03		1.6E+03
				2.0E-02	O						1.36E+09		1	0.1	Dodine	2439-10-3					2.3E+04	5.5E+04		1.6E+04
				5.0E-02	O			V			1.36E+09	1.17E+05	1		EPTC	759-94-4					5.8E+04			5.8E+04
				6.0E-03	I			V			1.36E+09	4.10E+05	1		Endosulfan	115-29-7					7.0E+03			7.0E+03
				6.0E-03	P						1.36E+09		1	0.1	Endosulfan Sulfate	1031-07-8					7.0E+03	1.7E+04		4.9E+03
				2.0E-02	I						1.36E+09		1	0.1	Endothall	145-73-3					2.3E+04	5.5E+04		1.6E+04
				3.0E-04	I						1.36E+09		1	0.1	Endrin	72-20-8					3.5E+02	8.3E+02		2.5E+02
9.9E-03	I	1.2E-06	I	6.0E-03	P	1.0E-03	I	V		1.05E+04	1.36E+09	1.89E+04	1		Epichlorohydrin	106-89-8	3.3E+04		1.9E+04	1.2E+04	7.0E+03		8.3E+01	8.2E+01
						2.0E-02	I	V		1.53E+04	1.36E+09	7.66E+03	1		Epoxybutane, 1,2-	106-88-7							6.7E+02	6.7E+02
				4.0E-02	P						1.36E+09		1	0.1	Ethanol, 2-(2-methoxyethoxy)-	111-77-3					4.7E+04	1.1E+05		3.3E+04
				5.0E-03	I						1.36E+09		1	0.1	Ethephon	16672-87-0					5.8E+03	1.4E+04		4.1E+03
				5.0E-04	I						1.36E+09		1	0.1	Ethion	563-12-2					5.8E+02	1.4E+03		4.1E+02

RMLs TR=1E-04 THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
				1.0E-01	P	6.0E-02	P	V		2.38E+04	1.36E+09	6.15E+04	1		Ethoxyethanol Acetate, 2-	111-15-9					1.2E+05		1.6E+04	1.4E+04
				9.0E-02	P	4.0E-02	P	V		1.06E+05	1.36E+09	9.84E+04	1		Ethoxyethanol, 2-	110-80-5					1.1E+05		1.7E+04	1.5E+04
				7.0E-01	P	7.0E-02	P	V		1.08E+04	1.36E+09	8.62E+03	1		Ethyl Acetate	141-78-6					8.2E+05		2.6E+03	2.6E+03
				5.0E-03	P	8.0E-03	P	V		2.50E+03	1.36E+09	6.34E+03	1		Ethyl Acrylate	140-88-5					5.8E+03		2.2E+02	2.1E+02
						4.0E+00	P	V		2.12E+03	1.36E+09	1.29E+03	1		Ethyl Chloride (Chloroethane)	75-00-3							2.3E+04	2.3E+04
				2.0E-01	I			V		1.01E+04	1.36E+09	3.12E+03	1		Ethyl Ether	60-29-7					2.3E+05			2.3E+05
						3.0E-01	P	V		1.10E+03	1.36E+09	5.77E+03	1		Ethyl Methacrylate	97-63-2							7.6E+03	7.6E+03
		8.0E-08	I	1.0E+00	I	4.0E+01	I	V		2.87E+03	1.36E+09	3.67E+03	1		Ethyl Tertiary Butyl Ether (ETBE)	637-92-3			5.6E+04	5.6E+04	1.2E+06		6.4E+05	4.1E+05
				1.0E-05	I						1.36E+09		1	0.1	Ethyl-p-nitrophenyl Phosphonate	2104-64-5					1.2E+01	2.8E+01		8.2E+00
1.1E-02	C	2.5E-06	C	5.0E-02	P	1.0E+00	I	V		4.80E+02	1.36E+09	5.67E+03	1		Ethylbenzene	100-41-4	3.0E+04		2.8E+03	2.5E+03	5.8E+04		2.5E+04	1.7E+04
				7.0E-02	P						1.36E+09		1	0.1	Ethylene Cyanohydrin	109-78-4					8.2E+04	1.9E+05		5.7E+04
				9.0E-02	P			V		1.89E+05	1.36E+09	1.80E+05	1		Ethylene Diamine	107-15-3					1.1E+05			1.1E+05
				8.0E-01	A	4.0E-01	C				1.36E+09		1	0.1	Ethylene Glycol	107-21-1					9.3E+05	2.2E+06	2.4E+09	6.6E+05
				1.0E-01	I	1.6E+00	I				1.36E+09		1	0.1	Ethylene Glycol Monobutyl Ether	111-76-2					1.2E+05	2.8E+05	9.5E+09	8.2E+04
3.1E-01	C	3.0E-03	I			3.0E-02	C	V	M	1.21E+05	1.36E+09	6.09E+03	1		Ethylene Oxide	75-21-8	1.1E+03		2.5E+00	2.5E+00			8.0E+02	8.0E+02
4.5E-02	C	1.3E-05	C	8.0E-05	I						1.36E+09		1	0.1	Ethylene Thiourea	96-45-7	7.3E+03	1.7E+04	1.3E+08	5.1E+03	9.3E+01	2.2E+02		6.6E+01
6.5E+01	C	1.9E-02	C					V		1.54E+05	1.36E+09	2.39E+04	1		Ethyleneimine	151-56-4	5.0E+00		1.5E+00	1.2E+00				
				3.0E+00	I						1.36E+09		1	0.1	Ethylphthalyl Ethyl Glycolate	84-72-0					3.5E+06	8.3E+06		2.5E+06
				2.5E-04	I						1.36E+09		1	0.1	Fenamiphos	22224-92-6					2.9E+02	6.9E+02		2.1E+02
				2.5E-02	I						1.36E+09		1	0.1	Fenpropathrin	39515-41-8					2.9E+04	6.9E+04		2.1E+04
				2.5E-02	I						1.36E+09		1	0.1	Fenvalerate	51630-58-1					2.9E+04	6.9E+04		2.1E+04
				1.3E-02	I						1.36E+09		1	0.1	Fluometuron	2164-17-2					1.5E+04	3.6E+04		1.1E+04
				4.0E-02	C	1.3E-02	C				1.36E+09		1		Fluoride	16984-48-8					4.7E+04		7.7E+07	4.7E+04
				6.0E-02	I	1.3E-02	C				1.36E+09		1		Fluorine (Soluble Fluoride)	7782-41-4					7.0E+04		7.7E+07	7.0E+04
				8.0E-02	I						1.36E+09		1	0.1	Fluridone	59756-60-4					9.3E+04	2.2E+05		6.6E+04
				4.0E-02	O						1.36E+09		1	0.1	Flurprimidol	56425-91-3					4.7E+04	1.1E+05		3.3E+04
				2.0E-03	O						1.36E+09		1	0.1	Flusilazole	85509-19-9					2.3E+03	5.5E+03		1.6E+03
				5.0E-01	O						1.36E+09		1	0.1	Flutolanil	66332-96-5					5.8E+05	1.4E+06		4.1E+05
				1.0E-02	I						1.36E+09		1	0.1	Fluvalinate	69409-94-5					1.2E+04	2.8E+04		8.2E+03
				9.0E-02	O						1.36E+09		1	0.1	Folpet	133-07-3					1.1E+05	2.5E+05		7.4E+04
				1.0E-02	O						1.36E+09		1	0.1	Fomesafen	72178-02-0					1.2E+04	2.8E+04		8.2E+03
				2.0E-03	I						1.36E+09		1	0.1	Fonofos	944-22-9					2.3E+03	5.5E+03		1.6E+03
2.1E-02	C	1.3E-05	I	2.0E-01	I	9.8E-03	A	V		4.24E+04	1.36E+09	7.77E+04	1		Formaldehyde	50-00-0	1.6E+04		7.3E+03	5.0E+03	2.3E+05		3.3E+03	3.3E+03
				9.0E-01	P	3.0E-04	X	V		1.06E+05	1.36E+09	9.30E+04	1		Formic Acid	64-18-6					1.1E+06		1.2E+02	1.2E+02
				2.5E+00	O						1.36E+09		1	0.1	Fosetyl-AL	39148-24-8					2.9E+06	6.9E+06		2.1E+06
															Furans									
				1.0E-03	X			V			1.36E+09	1.56E+05	1		~Dibenzofuran	132-64-9					1.2E+03			1.2E+03
				1.0E-03	I			V		6.22E+03	1.36E+09	2.62E+03	1		~Furan	110-00-9					1.2E+03			1.2E+03
				9.0E-01	I	2.0E+00	I	V		1.65E+05	1.36E+09	1.20E+04	1		~Tetrahydrofuran	109-99-9					1.1E+06		1.0E+05	9.5E+04
3.8E+00	H										1.36E+09		1	0.1	Furazolidone	67-45-8	8.6E+01	2.0E+02		6.0E+01				
				3.0E-03	I	5.0E-02	H	V		1.01E+04	1.36E+09	4.86E+04	1		Furfural	98-01-1					3.5E+03		1.1E+04	2.6E+03
1.5E+00	C	4.3E-04	C								1.36E+09		1	0.1	Furium	531-82-8	2.2E+02	5.2E+02	3.9E+06	1.5E+02				
3.0E-02	I	8.6E-06	C								1.36E+09		1	0.1	Furmecyclox	60568-05-0	1.1E+04	2.6E+04	1.9E+08	7.7E+03				
				6.0E-03	O						1.36E+09		1	0.1	Glufosinate, Ammonium	77182-82-2					7.0E+03	1.7E+04		4.9E+03
				1.0E-01	A	8.0E-05	C				1.36E+09		1	0.1	Glutaraldehyde	111-30-8					1.2E+05	2.8E+05	4.8E+05	7.0E+04
				4.0E-04	I	1.0E-03	X	V		1.06E+05	1.36E+09	8.43E+04	1		Glycidaldehyde	765-34-4					4.7E+02		3.7E+02	2.1E+02
				1.0E-01	I						1.36E+09		1	0.1	Glyphosate	1071-83-6					1.2E+05	2.8E+05		8.2E+04
				1.0E-02	X			V			1.36E+09	1.45E+05	1		Guanidine	113-00-8					1.2E+04			1.2E+04
				2.0E-02	P						1.36E+09		1	0.1	Guanidine Chloride	50-01-1					2.3E+04	5.5E+04		1.6E+04
				3.0E-02	X						1.36E+09		1	0.1	Guanidine Nitrate	506-93-4					3.5E+04	8.3E+04		2.5E+04
				5.0E-05	I						1.36E+09		1	0.1	Haloxyfop, Methyl	69806-40-2					5.8E+01	1.4E+02		4.1E+01
4.5E+00	I	1.3E-03	I	1.0E-04	A			V			1.36E+09	4.79E+05	1		Heptachlor	76-44-8	7.3E+01		4.5E+02	6.3E+01	1.2E+02			1.2E+02
9.1E+00	I	2.6E-03	I	1.3E-05	I			V			1.36E+09	8.43E+05	1		Heptachlor Epoxide	1024-57-3	3.6E+01		4.0E+02	3.3E+01	1.5E+01			1.5E+01
						3.0E-03	X	V		2.09E+02	1.36E+09	7.80E+03	1		Heptanal, n-	111-71-7							1.0E+02	1.0E+02
				3.0E-04	X	4.0E-01	P	V		5.79E+01	1.36E+09	8.95E+02	1		Heptane, N-	142-82-5					3.5E+02		1.6E+03	2.9E+02
				2.0E-03	I			V			1.36E+09	3.80E+05	1		Hexabromobenzene	87-82-1					2.3E+03			2.3E+03
				2.0E-04	I						1.36E+09		1	0.1	Hexabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-153)	68631-49-2					2.3E+02	5.5E+02		1.6E+02
1.6E+00	I	4.6E-04	I	1.0E-05	P			V			1.36E+09	6.80E+04	1		Hexachlorobenzene	118-74-1	2.0E+02		1.8E+02	9.6E+01	1.2E+01			1.2E+01
7.8E-02	I	2.2E-05	I	1.0E-03	P			V		1.68E+01	1.36E+09	1.08E+04	1		Hexachlorobutadiene	87-68-3	4.2E+03		6.0E+02	5.3E+02	1.2E+03			1.2E+03
6.3E+00	I	1.8E-03	I	8.0E-03	A						1.36E+09		1	0.1	Hexachlorocyclohexane, Alpha-	319-84-6	5.2E+01	1.2E+02	9.3E+05	3.6E+01	9.3E+03	2.2E+04		6.6E+03
1.8E+00	I	5.3E-04	I								1.36E+09		1	0.1	Hexachlorocyclohexane, Beta-	319-85-7	1.8E+02	4.3E+02	3.1E+06	1.3E+02				
1.1E+00	C	3.1E-04	C	1.0E-05	A						1.36E+09		1	0.04	Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	3.0E+02	1.8E+03	5.4E+06	2.5E+02	1.2E+01	6.9E+01		1.0E+01
1.8E+00	I	5.1E-04	I								1.36E+09		1	0.1	Hexachlorocyclohexane, Technical	608-73-1	1.8E+02	4.3E+02	3.3E+06	1.3E+02				
				6.0E-03	I	2.0E-04	I	V		1.57E+01	1.36E+09	8.51E+03	1		Hexachlorocyclopentadiene	77-47-4					7.0E+03		7.5E+00	7.5E+00
4.0E-02	I	1.1E-05	C	7.0E-04	I	3.0E-02	I	V			1.36E+09	8.01E+03	1		Hexachloroethane	67-72-1	8.2E+03		8.9E+02	8.0E+02	8.2E+02		1.1E+03	4.6E+02
				3.0E-04	I						1.36E+09		1	0.1	Hexachlorophene	70-30-4					3.5E+02	8.3E+02		2.5E+02
8.0E-02	I			4.0E-03	I						1.36E+09		1	0.015	Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	4.1E+03	6.4E+04		3.8E+03	4.7E+03	7.4E+04		4.4E+03
						1.0E-05	I	V		3.39E+03	1.36E+09	3.00E+05	1		Hexamethylene Diisocyanate, 1,6-	822-06-0							1.3E+01	1.3E+01
						4.0E-04	C				1.36E+09		1	0.1	Hexamethylene diisocyanate biuret	4035-89-6							2.4E+06	2.4E+06
						4.0E-04	C				1.36E+09		1	0.1	Hexamethylene diisocyanate isocyanurate	3779-63-3							2.4E+06	2.4E+06
				4.0E-04	P						1.36E+09		1	0.1	Hexamethylphosphoramide	680-31-9					4.7E+02	1.1E+03		3.3E+02
		2.0E-07	X			6.0E-01	P	V		1.41E+02	1.36E+09	8.29E+02	1		Hexane, Commercial	E5241997			5.1E+03	5.1E+03			2.2E+03	2.2E+03
						7.0E-01	I	V		1.41E+02	1.36E+09	8.29E+02	1		Hexane, N-	110-54-3							2.5E+03	2.5E+03
				2.0E+00	P						1.36E+09		1	0.1	Hexanedioic Acid	124-04-9					2.3E+06	5.5E+06		1.6E+06
9.5E-03	P			7.0E-02	P	4.0E-04	P	V		2.74E+02	1.36E+09	3.62E+04	1		Hexanol, 1-,2-ethyl- (2-Ethyl-1-hexanol)	104-76-7	3.4E+04			3.4E+04	8.2E+04		6.3E+01	6.3E+01
				5.0E-03	I	3.0E-02	I	V		3.28E+03	1.36E+09	1.33E+04	1		Hexanone, 2-	591-78-6					5.8E+03		1.7E+03	1.3E+03
				3.3E-02	I						1.36E+09		1	0.1	Hexazinone	51235-04-2					3.9E+04	9.1E+04		2.7E+04
				2.5E-02	I						1.36E+09		1	0.1	Hexythiazox	78587-05-0					2.9E+04	6.9E+04		2.1E+04

RMLS
TR=1E-04
THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
				1.7E-02	O						1.36E+09		1	0.1	Hydramethylnon	67485-29-4					2.0E+04	4.7E+04		1.4E+04
3.0E+00	I	4.9E-03	I			3.0E-05	P	V		1.12E+05	1.36E+09	6.52E+04	1		Hydrazine	302-01-2	1.1E+02		1.6E+01	1.4E+01			8.6E+00	8.6E+00
3.0E+00	I	4.9E-03	I								1.36E+09		1		Hydrazine Sulfate	10034-93-2	1.1E+02		3.4E+05	1.1E+02				
						2.0E-02	I	V			1.36E+09		1		Hydrogen Chloride	7647-01-0							1.2E+08	1.2E+08
				4.0E-02	C	1.4E-02	C	V			1.36E+09		1		Hydrogen Fluoride	7664-39-3					4.7E+04		8.3E+07	4.7E+04
						2.0E-03	I	V			1.36E+09		1		Hydrogen Sulfide	7783-06-4							1.2E+07	1.2E+07
6.0E-02	P			4.0E-02	P						1.36E+09		1	0.1	Hydroquinone	123-31-9	5.5E+03	1.3E+04		3.8E+03	4.7E+04	1.1E+05		3.3E+04
6.1E-02	O			1.1E-01	O						1.36E+09		1	0.1	Imazalil	35554-44-0	5.4E+03	1.3E+04		3.8E+03	1.3E+05	3.0E+05		8.9E+04
				2.5E-01	I						1.36E+09		1	0.1	Imazaquin	81335-37-7					2.9E+05	6.9E+05		2.1E+05
				2.5E+00	O						1.36E+09		1	0.1	Imazethapyr	81335-77-5					2.9E+06	6.9E+06		2.1E+06
				1.0E-02	A						1.36E+09		1		Iodine	7553-56-2					1.2E+04			1.2E+04
				4.0E-02	I						1.36E+09		1	0.1	Iprodione	36734-19-7					4.7E+04	1.1E+05		3.3E+04
				7.0E-01	P						1.36E+09		1		Iron	7439-89-6					8.2E+05			8.2E+05
				3.0E-01	I			V		1.00E+04	1.36E+09	2.81E+04	1		Isobutyl Alcohol	78-83-1					3.5E+05			3.5E+05
9.5E-04	I			2.0E-01	I	2.0E+00	C				1.36E+09		1	0.1	Isophorone	78-59-1	3.4E+05	8.1E+05		2.4E+05	2.3E+05	5.5E+05	1.2E+10	1.6E+05
				1.5E-02	I			V			1.36E+09	4.20E+05	1		Isopropalin	33820-53-0					1.8E+04			1.8E+04
				2.0E+00	P	2.0E-01	P	V		1.09E+05	1.36E+09	2.77E+04	1		Isopropanol	67-63-0					2.3E+06		2.4E+04	2.4E+04
				1.0E-01	I						1.36E+09		1	0.1	Isopropyl Methyl Phosphonic Acid	1832-54-8					1.2E+05	2.8E+05		8.2E+04
				5.0E-02	I						1.36E+09		1	0.1	Isoxaben	82558-50-7					5.8E+04	1.4E+05		4.1E+04
						3.0E-01	A	V			1.36E+09		1		JP-7	E1737665							1.8E+09	1.8E+09
				8.0E-03	O						1.36E+09		1	0.1	Lactofen	77501-63-4					9.3E+03	2.2E+04		6.6E+03
				2.0E-04	X						1.36E+09		1	0.1	Lactonitrile	78-97-7					2.3E+02	5.5E+02		1.6E+02
				5.0E-05	P						1.36E+09		1		Lanthanum	7439-91-0					5.8E+01			5.8E+01
				2.1E-05	P						1.36E+09		1	0.1	Lanthanum Acetate Hydrate	100587-90-4					2.4E+01	5.7E+01		1.7E+01
				1.9E-05	P						1.36E+09		1		Lanthanum Chloride Heptahydrate	10025-84-0					2.2E+01			2.2E+01
				2.8E-05	P						1.36E+09		1		Lanthanum Chloride, Anhydrous	10099-58-8					3.3E+01			3.3E+01
				1.6E-05	P						1.36E+09		1		Lanthanum Nitrate Hexahydrate	10277-43-7					1.9E+01			1.9E+01
															Lead Compounds									
8.5E-03	C	1.2E-05	C								1.36E+09		1		~Lead Phosphate	7446-27-7	3.8E+04		1.4E+08	3.8E+04				
2.1E-01	C	8.0E-05	C								1.36E+09		1	0.1	~Lead acetate	301-04-2	1.6E+03	3.7E+03	2.1E+07	1.1E+03				
											1.36E+09		1		~Lead and Compounds	7439-92-1								8.0E+02
3.8E-02	C	1.1E-05	C								1.36E+09		1	0.1	~Lead subacetate	1335-32-6	8.6E+03	2.0E+04	1.5E+08	6.0E+03				
				1.0E-07	I			V		2.43E+00	1.36E+09	1.91E+03	1		~Tetraethyl Lead	78-00-2					1.2E-01			1.2E-01
				5.0E-06	P			V		3.83E+02	1.36E+09	2.55E+04	1		Lewisite	541-25-3					5.8E+00			5.8E+00
				7.7E-03	O						1.36E+09		1	0.1	Linuron	330-55-2					9.0E+03	2.1E+04		6.3E+03
				2.0E-03	P						1.36E+09		1		Lithium	7439-93-2					2.3E+03			2.3E+03
				5.0E-04	I						1.36E+09		1	0.1	MCPA	94-74-6					5.8E+02	1.4E+03		4.1E+02
				4.4E-03	O						1.36E+09		1	0.1	MCPB	94-81-5					5.1E+03	1.2E+04		3.6E+03
				1.0E-03	I						1.36E+09		1	0.1	MCPP	93-65-2					1.2E+03	2.8E+03		8.2E+02
				2.0E-02	I						1.36E+09		1	0.1	Malathion	121-75-5					2.3E+04	5.5E+04		1.6E+04
				1.0E-01	I	7.0E-04	C				1.36E+09		1	0.1	Maleic Anhydride	108-31-6					1.2E+05	2.8E+05	4.2E+06	8.0E+04
				5.0E-01	I						1.36E+09		1	0.1	Maleic Hydrazide	123-33-1					5.8E+05	1.4E+06		4.1E+05
				1.0E-04	P						1.36E+09		1	0.1	Malononitrile	109-77-3					1.2E+02	2.8E+02		8.2E+01
				3.0E-02	H						1.36E+09		1	0.1	Mancozeb	8018-01-7					3.5E+04	8.3E+04		2.5E+04
				5.0E-03	I						1.36E+09		1	0.1	Maneb	12427-38-2					5.8E+03	1.4E+04		4.1E+03
				1.4E-01	I	5.0E-05	I						1		Manganese (Diet)	7439-96-5								
				2.4E-02	G	5.0E-05	I				1.36E+09		0.04		Manganese (Non-diet)	7439-96-5					2.8E+04		3.0E+05	2.6E+04
				9.0E-05	H						1.36E+09		1	0.1	Mephosfolan	950-10-7					1.1E+02	2.5E+02		7.4E+01
				3.0E-02	I						1.36E+09		1	0.1	Mepiquat Chloride	24307-26-4					3.5E+04	8.3E+04		2.5E+04
1.1E-02	P			4.0E-03	P						1.36E+09		1	0.1	Mercaptobenzothiazole, 2-	149-30-4	3.0E+04	7.0E+04		2.1E+04	4.7E+03	1.1E+04		3.3E+03
															Mercury Compounds									
				3.0E-04	I	3.0E-04	G				1.36E+09		0.07		~Mercuric Chloride (and other Mercury salts)	7487-94-7					3.5E+02		1.8E+06	3.5E+02
						3.0E-04	I	V		3.13E+00	1.36E+09	3.47E+04	1		~Mercury (elemental)	7439-97-6							4.6E+01	4.6E+01
				1.0E-04	I						1.36E+09		1		~Methyl Mercury	22967-92-6					1.2E+02			1.2E+02
				8.0E-05	I						1.36E+09		1	0.1	~Phenylmercuric Acetate	62-38-4					9.3E+01	2.2E+02		6.6E+01
				3.0E-05	I			V			1.36E+09	1.94E+06	1		Merphos	150-50-5					3.5E+01			3.5E+01
				6.0E-02	I						1.36E+09		1	0.1	Metalaxyl	57837-19-1					7.0E+04	1.7E+05		4.9E+04
				1.0E-04	I	3.0E-02	P	V		4.58E+03	1.36E+09	6.79E+03	1		Methacrylonitrile	126-98-7					1.2E+02		8.9E+02	1.0E+02
				5.0E-05	I						1.36E+09		1	0.1	Methamidophos	10265-92-6					5.8E+01	1.4E+02		4.1E+01
				2.0E+00	I	2.0E+01	I	V		1.06E+05	1.36E+09	2.90E+04	1		Methanol	67-56-1					2.3E+06		2.5E+06	1.2E+06
				1.5E-03	O						1.36E+09		1	0.1	Methidathion	950-37-8					1.8E+03	4.1E+03		1.2E+03
				2.5E-02	I						1.36E+09		1	0.1	Methomyl	16752-77-5					2.9E+04	6.9E+04		2.1E+04
4.9E-02	C										1.36E+09		1	0.1	Methoxy-5-nitroaniline, 2-	99-59-2	6.7E+03	1.6E+04		4.7E+03				
				5.0E-03	I						1.36E+09		1	0.1	Methoxychlor	72-43-5					5.8E+03	1.4E+04		4.1E+03
				8.0E-03	P	1.0E-03	P	V		1.15E+05	1.36E+09	1.24E+05	1		Methoxyethanol Acetate, 2-	110-49-6					9.3E+03		5.4E+02	5.1E+02
				5.0E-03	P	7.0E-03	P	V		1.06E+05	1.36E+09	1.01E+05	1		Methoxyethanol, 2-	109-86-4					5.8E+03		3.1E+03	2.0E+03
				1.0E+00	X			V		2.90E+04	1.36E+09	8.12E+03	1		Methyl Acetate	79-20-9					1.2E+06			1.2E+06
						2.0E-02	P	V		6.75E+03	1.36E+09	6.97E+03	1		Methyl Acrylate	96-33-3							6.1E+02	6.1E+02
				6.0E-01	I	5.0E+00	I	V		2.84E+04	1.36E+09	1.22E+04	1		Methyl Ethyl Ketone (2-Butanone)	78-93-3					7.0E+05		2.7E+05	1.9E+05
		1.0E-03	X	1.0E-03	P	2.0E-05	X	V		1.80E+05	1.36E+09	5.04E+04	1		Methyl Hydrazine	60-34-4			6.2E+01	6.2E+01	1.2E+03		4.4E+00	4.4E+00
						3.0E+00	I	V		3.36E+03	1.36E+09	1.06E+04	1		Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1							1.4E+05	1.4E+05
						1.0E-03	C	V		1.01E+04	1.36E+09	4.42E+03	1		Methyl Isocyanate	624-83-9							1.9E+01	1.9E+01
				1.4E+00	I	7.0E-01	I	V		2.36E+03	1.36E+09	6.33E+03	1		Methyl Methacrylate	80-62-6					1.6E+06		1.9E+04	1.9E+04
				2.5E-04	I						1.36E+09		1	0.1	Methyl Parathion	298-00-0					2.9E+02	6.9E+02		2.1E+02
				6.0E-02	X						1.36E+09		1	0.1	Methyl Phosphonic Acid	993-13-5					7.0E+04	1.7E+05		4.9E+04
				6.0E-03	H	4.0E-02	H	V		3.93E+02	1.36E+09	2.43E+04	1		Methyl Styrene (Mixed Isomers)	25013-15-4					7.0E+03		4.3E+03	2.6E+03
9.9E-02	C	2.8E-05	C								1.36E+09		1	0.1	Methyl methanesulfonate	66-27-3	3.3E+03	7.8E+03	6.0E+07	2.3E+03				
1.8E-03	C	2.6E-07	C			3.0E+00	I	V		8.87E+03	1.36E+09	4.90E+03	1		Methyl tert-Butyl Ether (MTBE)	1634-04-4	1.8E+05		2.3E+04	2.1E+04			6.4E+04	6.4E+04

RMLS
TR=1E-04
THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
				3.0E-04	X						1.36E+09		1	0.1	Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2					3.5E+02	8.3E+02		2.5E+02
						3.0E+00	X	V		2.45E+03	1.36E+09	1.72E+04	1		Methyl-2-Pentanol, 4-	108-11-2							2.3E+05	2.3E+05
9.0E-03	P			2.0E-02	X						1.36E+09		1	0.1	Methyl-5-Nitroaniline, 2-	99-55-8	3.6E+04	8.6E+04		2.6E+04	2.3E+04	5.5E+04		1.6E+04
8.3E+00	C	2.4E-03	C								1.36E+09		1	0.1	Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	3.9E+01	9.3E+01	6.9E+05	2.8E+01				
1.3E-01	C	3.7E-05	C								1.36E+09		1	0.1	Methylaniline Hydrochloride, 2-	636-21-5	2.5E+03	5.9E+03	4.5E+07	1.8E+03				
				1.0E-02	A						1.36E+09		1	0.1	Methylarsonic acid	124-58-3					1.2E+04	2.8E+04		8.2E+03
				2.0E-04	X						1.36E+09		1	0.1	Methylbenzene,1-4-diamine monohydrochloride, 2-	74612-12-7					2.3E+02	5.5E+02		1.6E+02
1.0E-01	X			3.0E-04	X						1.36E+09		1	0.1	Methylbenzene-1,4-diamine sulfate, 2-	615-50-9	3.3E+03	7.7E+03		2.3E+03	3.5E+02	8.3E+02		2.5E+02
2.2E+01	C	6.3E-03	C						M		1.36E+09		1	0.1	Methylcholanthrene, 3-	56-49-5	1.5E+01	3.5E+01	2.6E+05	1.0E+01				
2.0E-03	I	1.0E-08	I	6.0E-03	I	6.0E-01	I	V	M	3.32E+03	1.36E+09	2.19E+03	1		Methylene Chloride	75-09-2	1.6E+05		2.7E+05	1.0E+05	7.0E+03		5.8E+03	3.2E+03
1.0E-01	P	4.3E-04	C	2.0E-03	P				M		1.36E+09		1	0.1	Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	3.3E+03	7.7E+03	3.9E+06	2.3E+03	2.3E+03	5.5E+03		1.6E+03
4.6E-02	I	1.3E-05	C								1.36E+09		1	0.1	Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	7.1E+03	1.7E+04	1.3E+08	5.0E+03				
1.6E+00	C	4.6E-04	C			2.0E-02	C				1.36E+09		1	0.1	Methylenebisbenzenamine, 4,4'-	101-77-9	2.0E+02	4.8E+02	3.6E+06	1.4E+02			1.2E+08	1.2E+08
						6.0E-04	I				1.36E+09		1	0.1	Methylenediphenyl Diisocyanate	101-68-8							3.6E+06	3.6E+06
				7.0E-02	H			V		5.00E+02	1.36E+09	1.28E+04	1		Methylstyrene, Alpha-	98-83-9					8.2E+04			8.2E+04
				1.5E-01	I						1.36E+09		1	0.1	Metolachlor	51218-45-2					1.8E+05	4.1E+05		1.2E+05
				2.5E-02	I						1.36E+09		1	0.1	Metribuzin	21087-64-9					2.9E+04	6.9E+04		2.1E+04
				2.5E-01	I						1.36E+09		1	0.1	Metsulfuron-methyl	74223-64-6					2.9E+05	6.9E+05		2.1E+05
		4.5E-06	X	1.0E-02	X	1.0E-01	P	V		6.86E+00	1.36E+09	1.04E+03	1		Midrange Aliphatic Hydrocarbon Streams	E1790669			2.8E+02	2.8E+02	1.2E+04		4.6E+02	4.4E+02
				3.0E+00	P			V		3.42E-01	1.36E+09	1.38E+03	1		Mineral oils	8012-95-1					3.5E+06			3.5E+06
1.8E+01	C	5.1E-03	C	2.0E-04	I			V			1.36E+09	8.58E+05	1		Mirex	2385-85-5	1.8E+01		2.1E+02	1.7E+01	2.3E+02			2.3E+02
				2.0E-03	I						1.36E+09		1	0.1	Molinate	2212-67-1					2.3E+03	5.5E+03		1.6E+03
				5.0E-03	I	2.0E-03	A				1.36E+09		1		Molybdenum	7439-98-7					5.8E+03		1.2E+07	5.8E+03
				1.0E-01	I						1.36E+09		1		Monochloramine	10599-90-3					1.2E+05			1.2E+05
				2.0E-03	P						1.36E+09		1	0.1	Monomethylaniline	100-61-8					2.3E+03	5.5E+03		1.6E+03
				2.5E-02	I						1.36E+09		1	0.1	Myclobutanil	88671-89-0					2.9E+04	6.9E+04		2.1E+04
				3.0E-04	X						1.36E+09		1	0.1	N,N'-Diphenyl-1,4-benzenediamine	74-31-7					3.5E+02	8.3E+02		2.5E+02
				2.0E-03	I			V			1.36E+09	5.70E+04	1		Naled	300-76-5					2.3E+03			2.3E+03
				3.0E-02	X	1.0E-01	P	V			1.36E+09		1		Naphtha, High Flash Aromatic (HFAN)	64742-95-6					3.5E+04		6.0E+08	3.5E+04
1.8E+00	C	0.0E+00	C								1.36E+09		1	0.1	Naphthylamine, 2-	91-59-8	1.8E+02	4.3E+02		1.3E+02				
				1.2E-01	O						1.36E+09		1	0.1	Napropamide	15299-99-7					1.4E+05	3.3E+05		9.8E+04
		2.6E-04	C	1.1E-02	C	1.4E-05	C				1.36E+09		1	0.1	Nickel Acetate	373-02-4			6.4E+06	6.4E+06	1.3E+04	3.0E+04	8.3E+04	8.1E+03
		2.6E-04	C	1.1E-02	C	1.4E-05	C				1.36E+09		1	0.1	Nickel Carbonate	3333-67-3			6.4E+06	6.4E+06	1.3E+04	3.0E+04	8.3E+04	8.1E+03
		2.6E-04	C	1.1E-02	C	1.4E-05	C	V			1.36E+09		1		Nickel Carbonyl	13463-39-3			6.4E+06	6.4E+06	1.3E+04		8.3E+04	1.1E+04
		2.6E-04	C	1.1E-02	C	1.4E-05	C				1.36E+09			0.04	Nickel Hydroxide	12054-48-7			6.4E+06	6.4E+06	1.3E+04		8.3E+04	1.1E+04
		2.6E-04	C	1.1E-02	C	2.0E-05	C				1.36E+09			0.04	Nickel Oxide	1313-99-1			6.4E+06	6.4E+06	1.3E+04		1.2E+05	1.2E+04
		2.4E-04	I	1.1E-02	C	1.4E-05	C				1.36E+09			0.04	Nickel Refinery Dust	E715532			6.9E+06	6.9E+06	1.3E+04		8.3E+04	1.1E+04
		2.6E-04	C	2.0E-02	I	9.0E-05	A				1.36E+09			0.04	Nickel Soluble Salts	7440-02-0			6.4E+06	6.4E+06	2.3E+04		5.4E+05	2.2E+04
1.7E+00	C	4.8E-04	I	1.1E-02	C	1.4E-05	C				1.36E+09			0.04	Nickel Subsulfide	12035-72-2	1.9E+02		3.5E+06	1.9E+02	1.3E+04		8.3E+04	1.1E+04
9.1E-01	C	2.6E-04	C	1.1E-02	C	1.4E-05	C				1.36E+09		1	0.1	Nickelocene	1271-28-9	3.6E+02	8.5E+02	6.4E+06	2.5E+02	1.3E+04	3.0E+04	8.3E+04	8.1E+03
				1.6E+00	I						1.36E+09		1		Nitrate (measured as nitrogen)	14797-55-8					1.9E+06			1.9E+06
											1.36E+09		1		Nitrate + Nitrite (measured as nitrogen)	E701177								
				1.0E-01	I						1.36E+09		1		Nitrite (measured as nitrogen)	14797-65-0					1.2E+05			1.2E+05
				1.0E-02	X	5.0E-05	X				1.36E+09		1	0.1	Nitroaniline, 2-	88-74-4					1.2E+04	2.8E+04	3.0E+05	8.0E+03
2.0E-02	P			4.0E-03	P	6.0E-03	P				1.36E+09		1	0.1	Nitroaniline, 4-	100-01-6	1.6E+04	3.9E+04		1.1E+04	4.7E+03	1.1E+04	3.6E+07	3.3E+03
		4.0E-05	I	2.0E-03	I	9.0E-03	I	V		3.05E+03	1.36E+09	7.32E+04	1		Nitrobenzene	98-95-3			2.2E+03	2.2E+03	2.3E+03		2.9E+03	1.3E+03
				3.0E+03	P						1.36E+09		1	0.1	Nitrocellulose	9004-70-0					3.5E+09	8.3E+09		2.5E+09
				7.0E-02	H						1.36E+09		1	0.1	Nitrofurantoin	67-20-9					8.2E+04	1.9E+05		5.7E+04
1.3E+00	C	3.7E-04	C								1.36E+09		1	0.1	Nitrofurazone	59-87-0	2.5E+02	5.9E+02	4.5E+06	1.8E+02				
1.7E-02	P			1.0E-04	P						1.36E+09		1	0.1	Nitroglycerin	55-63-0	1.9E+04	4.5E+04		1.4E+04	1.2E+02	2.8E+02		8.2E+01
				1.0E-01	I						1.36E+09		1	0.1	Nitroguanidine	556-88-7					1.2E+05	2.8E+05		8.2E+04
		8.8E-06	P			5.0E-03	P	V		1.80E+04	1.36E+09	1.69E+04	1		Nitromethane	75-52-5			2.4E+03	2.4E+03			3.7E+02	3.7E+02
		5.8E-04	X			2.0E-02	I	V		4.86E+03	1.36E+09	1.31E+04	1		Nitropropane, 2-	79-46-9			2.8E+01	2.8E+01			1.2E+03	1.2E+03
2.7E+01	C	7.7E-03	C						M		1.36E+09		1	0.1	Nitroso-N-ethylurea, N-	759-73-9	1.2E+01	2.9E+01	2.2E+05	8.5E+00				
1.2E+02	C	3.4E-02	C						M		1.36E+09		1	0.1	Nitroso-N-methylurea, N-	684-93-5	2.7E+00	6.4E+00	4.9E+04	1.9E+00				
5.4E+00	I	1.6E-03	I					V			1.36E+09	2.43E+05	1		Nitroso-di-N-butylamine, N-	924-16-3	6.1E+01		1.9E+02	4.6E+01				
7.0E+00	I	2.0E-03	C								1.36E+09		1	0.1	Nitroso-di-N-propylamine, N-	621-64-7	4.7E+01	1.1E+02	8.3E+05	3.3E+01				
2.8E+00	I	8.0E-04	C								1.36E+09		1	0.1	Nitrosodiethanolamine, N-	1116-54-7	1.2E+02	2.8E+02	2.1E+06	8.2E+01				
1.5E+02	I	4.3E-02	I						M		1.36E+09		1	0.1	Nitrosodiethylamine, N-	55-18-5	2.2E+00	5.2E+00	3.9E+04	1.5E+00				
5.1E+01	I	1.4E-02	I	8.0E-06	P	4.0E-05	X	V	M	2.37E+05	1.36E+09	8.23E+04	1		Nitrosodimethylamine, N-	62-75-9	6.4E+00		7.2E+00	3.4E+00	9.3E+00		1.4E+01	5.7E+00
4.9E-03	I	2.6E-06	C								1.36E+09		1	0.1	Nitrosodiphenylamine, N-	86-30-6	6.7E+04	1.6E+05	6.4E+08	4.7E+04				
2.2E+01	I	6.3E-03	C					V		1.08E+05	1.36E+09	1.21E+05	1		Nitrosomethylethylamine, N-	10595-95-6	1.5E+01		2.4E+01	9.1E+00				
6.7E+00	C	1.9E-03	C								1.36E+09		1	0.1	Nitrosomorpholine [N-]	59-89-2	4.9E+01	1.2E+02	8.8E+05	3.4E+01				
9.4E+00	C	2.7E-03	C								1.36E+09		1	0.1	Nitrosopiperidine [N-]	100-75-4	3.5E+01	8.2E+01	6.2E+05	2.4E+01				
2.1E+00	I	6.1E-04	I								1.36E+09		1	0.1	Nitrosopyrrolidine, N-	930-55-2	1.6E+02	3.7E+02	2.7E+06	1.1E+02				
				1.0E-04	X						1.36E+09		1	0.1	Nitrotoluene, m-	99-08-1					1.2E+02	2.8E+02		8.2E+01
2.2E-01	P			9.0E-04	P			V		1.51E+03	1.36E+09	1.37E+05	1		Nitrotoluene, o-	88-72-2	1.5E+03			1.5E+03	1.1E+03			1.1E+03
1.6E-02	P			4.0E-03	P						1.36E+09		1	0.1	Nitrotoluene, p-	99-99-0	2.0E+04	4.8E+04		1.4E+04	4.7E+03	1.1E+04		3.3E+03
				3.0E-04	X	2.0E-02	P	V		6.86E+00	1.36E+09	1.04E+03	1		Nonane, n-	111-84-2					3.5E+02		9.1E+01	7.2E+01
				1.5E-03	O						1.36E+09		1	0.1	Norflurazon	27314-13-2					1.8E+03	4.1E+03		1.2E+03
				3.0E-03	I						1.36E+09		1	0.1	Octabromodiphenyl Ether	32536-52-0					3.5E+03	8.3E+03		2.5E+03
				5.0E-02	I						1.36E+09		1	0.006	Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0					5.8E+04	2.3E+06		5.7E+04
				2.0E-03	H						1.36E+09		1	0.1	Octamethylpyrophosphoramide	152-16-9					2.3E+03	5.5E+03		1.6E+03
7.8E-03	O			1.9E-01	O						1.36E+09		1	0.1	Oryzalin	19044-88-3	4.2E+04	9.9E+04		2.9E+04	2.2E+05	5.2E+05		1.6E+05
				5.0E-03	I						1.36E+09		1	0.1	Oxadiazon	19666-30-9					5.8E+03	1.4E+04		4.1E+03
				2.5E-02	I						1.36E+09		1	0.1	Oxamyl	23135-22-0					2.9E+04	6.9E+04		2.1E+04
7.3E-02	O			4.0E-02	O						1.36E+09		1	0.1	Oxyfluorfen	42874-03-3	4.5E+03	1.1E+04		3.1E+03	4.7E+04	1.1E+05		3.3E+04
				1.3E-02	I						1.36E+09		1	0.1	Paclobutrazol	76738-62-0					1.5E+04	3.6E+04		1.1E+04

RMLS
TR=1E-04
THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
				4.5E-03	I						1.36E+09		1	0.1	Paraquat Dichloride	1910-42-5					5.3E+03	1.2E+04		3.7E+03
				6.0E-03	H						1.36E+09		1	0.1	Parathion	56-38-2					7.0E+03	1.7E+04		4.9E+03
				5.0E-02	H			V			1.36E+09	4.49E+04	1		Pebulate	1114-71-2					5.8E+04			5.8E+04
				3.0E-01	O						1.36E+09		1	0.1	Pendimethalin	40487-42-1					3.5E+05	8.3E+05		2.5E+05
				2.0E-03	I			V		3.12E-01	1.36E+09	5.13E+05	1		Pentabromodiphenyl Ether	32534-81-9					2.3E+03			2.3E+03
				1.0E-04	I						1.36E+09		1	0.1	Pentabromodiphenyl ether, 2,2',4,4',5- (BDE-99)	60348-60-9					1.2E+02	2.8E+02		8.2E+01
				8.0E-04	I			V			1.36E+09	8.12E+04	1		Pentachlorobenzene	608-93-5					9.3E+02			9.3E+02
9.0E-02	P							V		4.57E+02	1.36E+09	9.65E+03	1		Pentachloroethane	76-01-7	3.6E+03			3.6E+03				
2.6E-01	H			3.0E-03	I			V			1.36E+09	4.32E+05	1		Pentachloronitrobenzene	82-68-8	1.3E+03			1.3E+03	3.5E+03			3.5E+03
4.0E-01	I	5.1E-06	C	5.0E-03	I						1.36E+09		1	0.25	Pentachlorophenol	87-86-5	8.2E+02	7.7E+02	3.3E+08	4.0E+02	5.8E+03	5.5E+03		2.8E+03
4.3E-03	X			9.0E-03	P						1.36E+09		1	0.1	Pentaerythritol tetranitrate (PETN)	78-11-5	7.6E+04	1.8E+05		5.3E+04	1.1E+04	2.5E+04		7.4E+03
				1.0E-04	X						1.36E+09		1	0.1	Pentamethylphosphoramide (PMPA)	10159-46-3					1.2E+02	2.8E+02		8.2E+01
						1.0E+00	P	V		3.88E+02	1.36E+09	7.79E+02	1		Pentane, n-	109-66-0							3.4E+03	3.4E+03
															Per- and Polyfluoroalkyl Substances (PFAS)									
				3.0E-06	D						1.36E+09		1	0.1	~Ammonium perfluoro-2-methyl-3-oxahexanoate	62037-80-3					3.5E+00	8.3E+00		2.5E+00
				3.0E-06	D			V			1.36E+09		1		~Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6					3.5E+00			3.5E+00
				3.0E-04	P						1.36E+09		1	0.1	~Perfluorobutanesulfonate	45187-15-3					3.5E+02	8.3E+02		2.5E+02
				3.0E-04	P						1.36E+09		1	0.1	~Perfluorobutanesulfonic acid (PFBS)	375-73-5					3.5E+02	8.3E+02		2.5E+02
				2.0E-05	A						1.36E+09		1	0.1	~Perfluorohexanesulfonate	108427-53-8					2.3E+01	5.5E+01		1.6E+01
				2.0E-05	A						1.36E+09		1	0.1	~Perfluorohexanesulfonic acid (PFHxS)	355-46-4					2.3E+01	5.5E+01		1.6E+01
				3.0E-06	A						1.36E+09		1	0.1	~Perfluorononanoate	72007-68-2					3.5E+00	8.3E+00		2.5E+00
				3.0E-06	A						1.36E+09		1	0.1	~Perfluorononanoic acid (PFNA)	375-95-1					3.5E+00	8.3E+00		2.5E+00
				2.0E-06	A						1.36E+09		1	0.1	~Perfluorooctanesulfonate	45298-90-6					2.3E+00	5.5E+00		1.6E+00
				2.0E-06	A						1.36E+09		1	0.1	~Perfluorooctanesulfonic acid (PFOS)	1763-23-1					2.3E+00	5.5E+00		1.6E+00
7.0E-02	D			3.0E-06	A						1.36E+09		1	0.1	~Perfluorooctanoate	45285-51-6	4.7E+03	1.1E+04		3.3E+03	3.5E+00	8.3E+00		2.5E+00
7.0E-02	D			3.0E-06	A						1.36E+09		1	0.1	~Perfluorooctanoic acid (PFOA)	335-67-1	4.7E+03	1.1E+04		3.3E+03	3.5E+00	8.3E+00		2.5E+00
				3.0E-04	P						1.36E+09		1	0.1	~Potassium perfluorobutanesulfonate	29420-49-3					3.5E+02	8.3E+02		2.5E+02
				2.0E-06	A						1.36E+09		1	0.1	~Potassium perfluorooctanesulfonate	2795-39-3					2.3E+00	5.5E+00		1.6E+00
															Perchlorates									
				7.0E-04	I						1.36E+09		1		~Ammonium Perchlorate	7790-98-9					8.2E+02			8.2E+02
				7.0E-04	I						1.36E+09		1		~Lithium Perchlorate	7791-03-9					8.2E+02			8.2E+02
				7.0E-04	I						1.36E+09		1		~Perchlorate and Perchlorate Salts	14797-73-0					8.2E+02			8.2E+02
				7.0E-04	I						1.36E+09		1		~Potassium Perchlorate	7778-74-7					8.2E+02			8.2E+02
				7.0E-04	I						1.36E+09		1		~Sodium Perchlorate	7601-89-0					8.2E+02			8.2E+02
				5.0E-02	I						1.36E+09		1	0.1	Permethrin	52645-53-1					5.8E+04	1.4E+05		4.1E+04
2.2E-03	C	6.3E-07	C								1.36E+09		1	0.1	Phenacetin	62-44-2	1.5E+05	3.5E+05	2.6E+09	1.0E+05				
				2.4E-01	O						1.36E+09		1	0.1	Phenmedipham	13684-63-4					2.8E+05	6.6E+05		2.0E+05
				3.0E-01	I	2.0E-01	C				1.36E+09		1	0.1	Phenol	108-95-2					3.5E+05	8.3E+05	1.2E+09	2.5E+05
				4.0E-03	I						1.36E+09		1	0.1	Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1					4.7E+03	1.1E+04		3.3E+03
				5.0E-04	X						1.36E+09		1	0.1	Phenothiazine	92-84-2					5.8E+02	1.4E+03		4.1E+02
				2.0E-04	X			V		1.29E+02	1.36E+09	7.06E+03	1		Phenyl Isothiocyanate	103-72-0					2.3E+02			2.3E+02
				6.0E-03	I						1.36E+09		1	0.1	Phenylenediamine, m-	108-45-2					7.0E+03	1.7E+04		4.9E+03
1.2E-01	P			4.0E-03	P						1.36E+09		1	0.1	Phenylenediamine, o-	95-54-5	2.7E+03	6.4E+03		1.9E+03	4.7E+03	1.1E+04		3.3E+03
				1.0E-03	X						1.36E+09		1	0.1	Phenylenediamine, p-	106-50-3					1.2E+03	2.8E+03		8.2E+02
1.9E-03	H										1.36E+09		1	0.1	Phenylphenol, 2-	90-43-7	1.7E+05	4.0E+05		1.2E+05				
				2.0E-04	H						1.36E+09		1	0.1	Phorate	298-02-2					2.3E+02	5.5E+02		1.6E+02
						3.0E-04	I	V		1.61E+03	1.36E+09	9.81E+02	1		Phosgene	75-44-5							1.3E+00	1.3E+00
				2.0E-02	I						1.36E+09		1	0.1	Phosmet	732-11-6					2.3E+04	5.5E+04		1.6E+04
				3.0E-04	I	3.0E-04	I	V			1.36E+09		1		Phosphine	7803-51-2					3.5E+02		1.8E+06	3.5E+02
						1.0E-02	I				1.36E+09		1		Phosphoric Acid	7664-38-2							6.0E+07	6.0E+07
				2.0E-05	I			V			1.36E+09	6.92E+03	1		Phosphorus, White	7723-14-0					2.3E+01			2.3E+01
															Phthalates									
1.4E-02	I	2.4E-06	C	2.0E-02	I						1.36E+09		1	0.1	~Bis(2-ethylhexyl)phthalate	117-81-7	2.3E+04	5.5E+04	6.9E+08	1.6E+04	2.3E+04	5.5E+04		1.6E+04
1.9E-03	P			2.0E-01	I						1.36E+09		1	0.1	~Butyl Benzyl Phthalate	85-68-7	1.7E+05	4.1E+05		1.2E+05	2.3E+05	5.5E+05		1.6E+05
				1.0E+00	I						1.36E+09		1	0.1	~Butylphthalyl Butylglycolate	85-70-1					1.2E+06	2.8E+06		8.2E+05
				1.0E-01	I						1.36E+09		1	0.1	~Dibutyl Phthalate	84-74-2					1.2E+05	2.8E+05		8.2E+04
				8.0E-01	I						1.36E+09		1	0.1	~Diethyl Phthalate	84-66-2					9.3E+05	2.2E+06		6.6E+05
				1.0E-01	I			V			1.36E+09	2.13E+04	1		~Dimethylterephthalate	120-61-6					1.2E+05			1.2E+05
				1.0E-02	P						1.36E+09		1	0.1	~Octyl Phthalate, di-N-	117-84-0					1.2E+04	2.8E+04		8.2E+03
				5.0E-01	X						1.36E+09		1	0.1	~Phthalic Acid, p-	100-21-0					5.8E+05	1.4E+06		4.1E+05
				2.0E+00	I	2.0E-02	C				1.36E+09		1	0.1	~Phthalic Anhydride	85-44-9					2.3E+06	5.5E+06	1.2E+08	1.6E+06
				7.0E-02	I						1.36E+09		1	0.1	Picloram	1918-02-1					8.2E+04	1.9E+05		5.7E+04
				1.0E-04	X						1.36E+09		1	0.1	Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3					1.2E+02	2.8E+02		8.2E+01
				2.0E-03	X						1.36E+09		1	0.1	Picric Acid (2,4,6-Trinitrophenol)	88-89-1					2.3E+03	5.5E+03		1.6E+03
				7.3E-04	O						1.36E+09		1	0.1	Pirimiphos, Methyl	29232-93-7					8.5E+02	2.0E+03		6.0E+02
3.0E+01	C	8.6E-03	C	7.0E-06	H						1.36E+09		1	0.1	Polybrominated Biphenyls	36355-01-8	1.1E+01	2.6E+01	1.9E+05	7.7E+00	8.2E+00	1.9E+01		5.7E+00
															Polychlorinated Biphenyls (PCBs)									
7.0E-02	G	2.0E-05	G	7.0E-05	I			V			1.36E+09	5.86E+05	1	0.14	~Aroclor 1016	12674-11-2	4.7E+03	7.9E+03	3.6E+04	2.7E+03	8.2E+01	1.4E+02		5.1E+01
2.0E+00	G	5.7E-04	G					V			1.36E+09	2.04E+05	1	0.14	~Aroclor 1221	11104-28-2	1.6E+02	2.8E+02	4.4E+02	8.3E+01				
2.0E+00	G	5.7E-04	G					V			1.36E+09	1.12E+05	1	0.14	~Aroclor 1232	11141-16-5	1.6E+02	2.8E+02	2.4E+02	7.2E+01				
2.0E+00	G	5.7E-04	G					V			1.36E+09	5.91E+05	1	0.14	~Aroclor 1242	53469-21-9	1.6E+02	2.8E+02	1.3E+03	9.5E+01				
2.0E+00	G	5.7E-04	G					V			1.36E+09	5.14E+05	1	0.14	~Aroclor 1248	12672-29-6	1.6E+02	2.8E+02	1.1E+03	9.4E+01				
2.0E+00	G	5.7E-04	G	2.0E-05	I			V			1.36E+09	8.43E+05	1	0.14	~Aroclor 1254	11097-69-1	1.6E+02	2.8E+02	1.8E+03	9.7E+01	2.3E+01	3.9E+01		1.5E+01
2.0E+00	G	5.7E-04	G					V			1.36E+09	1.31E+06	1	0.14	~Aroclor 1260	11096-82-5	1.6E+02	2.8E+02	2.8E+03	9.9E+01				
				6.0E-04	X			V			1.36E+09	1.15E+06	1	0.14	~Aroclor 5460	11126-42-4					7.0E+02	1.2E+03		4.4E+02
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	2.43E+06	1	0.14	~Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	8.4E+01	1.4E+02	2.6E+03	5.2E+01	2.7E+01	4.6E+01	1.4E+04	1.7E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.58E+06	1	0.14	~Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167)	52663-72-6	8.4E+01	1.4E+02	1.7E+03	5.1E+01	2.7E+01	4.6E+01	9.2E+03	1.7E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.04E+06	1	0.14	~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	8.4E+01	1.4E+02	1.1E+03	5.0E+01	2.7E+01	4.6E+01	6.1E+03	1.7E+01

RMLS
TR=1E-04
THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.11E+06	1	0.14	~Hexachlorobiphenyl, 2,3,3',4,4',5- (PCB 156)	38380-08-4	8.4E+01	1.4E+02	1.2E+03	5.0E+01	2.7E+01	4.6E+01	6.5E+03	1.7E+01
3.9E+03	W	1.1E+00	W	2.3E-08	W	1.3E-06	W	V			1.36E+09	1.58E+06	1	0.14	~Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	8.4E-02	1.4E-01	1.7E+00	5.1E-02	2.7E-02	4.6E-02	9.2E+00	1.7E-02
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	7.33E+05	1	0.14	~Pentachlorobiphenyl, 2',3,4,4',5- (PCB 123)	65510-44-3	8.4E+01	1.4E+02	7.9E+02	4.9E+01	2.7E+01	4.6E+01	4.3E+03	1.7E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	5.90E+05	1	0.14	~Pentachlorobiphenyl, 2,3',4,4',5- (PCB 118)	31508-00-6	8.4E+01	1.4E+02	6.3E+02	4.9E+01	2.7E+01	4.6E+01	3.4E+03	1.7E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	6.01E+05	1	0.14	~Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	8.4E+01	1.4E+02	6.5E+02	4.9E+01	2.7E+01	4.6E+01	3.5E+03	1.7E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.05E+06	1	0.14	~Pentachlorobiphenyl, 2,3,4,4',5- (PCB 114)	74472-37-0	8.4E+01	1.4E+02	1.1E+03	5.0E+01	2.7E+01	4.6E+01	6.1E+03	1.7E+01
1.3E+04	W	3.8E+00	W	7.0E-09	W	4.0E-07	W	V			1.36E+09	7.26E+05	1	0.14	~Pentachlorobiphenyl, 3,3',4,4',5- (PCB 126)	57465-28-8	2.5E-02	4.2E-02	2.3E-01	1.5E-02	8.2E-03	1.4E-02	1.3E+00	5.1E-03
2.0E+00	I	5.7E-04	I					V			1.36E+09	5.32E+05	1	0.14	~Polychlorinated Biphenyls (high risk)	1336-36-3	1.6E+02	2.8E+02	1.1E+03	9.4E+01				
4.0E-01	I	1.0E-04	I					V					1	0.14	~Polychlorinated Biphenyls (low risk)	1336-36-3								
7.0E-02	I	2.0E-05	I					V					1	0.14	~Polychlorinated Biphenyls (lowest risk)	1336-36-3								
1.3E+01	W	3.8E-03	W	7.0E-06	W	4.0E-04	W				1.36E+09		1	0.14	~Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	2.5E+01	4.2E+01	4.4E+05	1.6E+01	8.2E+00	1.4E+01	2.4E+06	5.1E+00
3.9E+01	W	1.1E-02	W	2.3E-06	W	1.3E-04	W	V			1.36E+09	5.09E+05	1	0.14	~Tetrachlorobiphenyl, 3,4,4',5- (PCB 81)	70362-50-4	8.4E+00	1.4E+01	5.5E+01	4.8E+00	2.7E+00	4.6E+00	3.0E+02	1.7E+00
						6.0E-04	I				1.36E+09		1	0.1	Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9							3.6E+06	3.6E+06
															Polynuclear Aromatic Hydrocarbons (PAHs)									
				6.0E-02	I			V			1.36E+09	1.41E+05	1	0.13	~Acenaphthene	83-32-9					7.0E+04	1.3E+05		4.5E+04
				3.0E-01	I			V			1.36E+09	5.23E+05	1	0.13	~Anthracene	120-12-7					3.5E+05	6.4E+05		2.3E+05
1.0E-01	E	6.0E-05	E					V	M		1.36E+09	4.41E+06	1	0.13	~Benz[a]anthracene	56-55-3	3.3E+03	5.9E+03	9.0E+04	2.1E+03				
				9.0E-05	X	2.0E-06	X				1.36E+09		1	0.1	~Benzo(e)pyrene	192-97-2					1.1E+02	2.5E+02	1.2E+04	7.3E+01
1.2E+00	C	1.1E-04	C								1.36E+09		1	0.13	~Benzo(i)fluoranthene	205-82-3	2.7E+02	5.0E+02	1.5E+07	1.8E+02				
1.0E+00	I	6.0E-04	I	3.0E-04	I	2.0E-06	I		M		1.36E+09		1	0.13	~Benzo[a]pyrene	50-32-8	3.3E+02	5.9E+02	2.8E+06	2.1E+02	3.5E+02	6.4E+02	1.2E+04	2.2E+02
1.0E-01	E	6.0E-05	E						M		1.36E+09		1	0.13	~Benzo[b]fluoranthene	205-99-2	3.3E+03	5.9E+03	2.8E+07	2.1E+03				
1.0E-02	E	6.0E-06	E						M		1.36E+09		1	0.13	~Benzo[k]fluoranthene	207-08-9	3.3E+04	5.9E+04	2.8E+08	2.1E+04				
				8.0E-02	I			V			1.36E+09	7.99E+04	1	0.13	~Chloronaphthalene, Beta-	91-58-7					9.3E+04	1.7E+05		6.0E+04
1.0E-03	E	6.0E-07	E						M		1.36E+09		1	0.13	~Chrysene	218-01-9	3.3E+05	5.9E+05	2.8E+09	2.1E+05				
1.0E+00	E	6.0E-04	E						M		1.36E+09		1	0.13	~Dibenz[a,h]anthracene	53-70-3	3.3E+02	5.9E+02	2.8E+06	2.1E+02				
1.2E+01	C	1.1E-03	C								1.36E+09		1	0.13	~Dibenzo(a,e)pyrene	192-65-4	2.7E+01	5.0E+01	1.5E+06	1.8E+01				
2.5E+02	C	7.1E-02	C						M		1.36E+09		1	0.13	~Dimethylbenz(a)anthracene, 7,12-	57-97-6	1.3E+00	2.4E+00	2.3E+04	8.4E-01				
				4.0E-02	I						1.36E+09		1	0.13	~Fluoranthene	206-44-0					4.7E+04	8.5E+04		3.0E+04
				4.0E-02	I			V			1.36E+09	2.81E+05	1	0.13	~Fluorene	86-73-7					4.7E+04	8.5E+04		3.0E+04
1.0E-01	E	6.0E-05	E						M		1.36E+09		1	0.13	~Indeno[1,2,3-cd]pyrene	193-39-5	3.3E+03	5.9E+03	2.8E+07	2.1E+03				
2.9E-02	P			7.0E-02	A			V		3.94E+02	1.36E+09	5.86E+04	1	0.13	~Methylnaphthalene, 1-	90-12-0	1.1E+04	2.0E+04		7.3E+03	8.2E+04	1.5E+05		5.3E+04
				4.0E-03	I			V			1.36E+09	5.80E+04	1	0.13	~Methylnaphthalene, 2-	91-57-6					4.7E+03	8.5E+03		3.0E+03
1.2E-01	C	3.4E-05	C	2.0E-02	I	3.0E-03	I	V			1.36E+09	4.63E+04	1	0.13	~Naphthalene	91-20-3	2.7E+03	5.0E+03	1.7E+03	8.6E+02	2.3E+04	4.2E+04	6.1E+02	5.9E+02
1.2E+00	C	1.1E-04	C								1.36E+09		1	0.13	~Nitropyrene, 4-	57835-92-4	2.7E+02	5.0E+02	1.5E+07	1.8E+02				
				3.0E-02	I			V			1.36E+09	2.38E+06	1	0.13	~Pyrene	129-00-0					3.5E+04	6.4E+04		2.3E+04
1.5E-01	I			9.0E-03	I						1.36E+09		1	0.1	Prochloraz	67747-09-5	2.2E+03	5.2E+03		1.5E+03	1.1E+04	2.5E+04		7.4E+03
				6.0E-03	H			V			1.36E+09	4.20E+05	1		Profluralin	26399-36-0					7.0E+03			7.0E+03
				1.5E-02	I						1.36E+09		1	0.1	Prometon	1610-18-0					1.8E+04	4.1E+04		1.2E+04
				4.0E-02	O						1.36E+09		1	0.1	Prometryn	7287-19-6					4.7E+04	1.1E+05		3.3E+04
				7.5E-02	I						1.36E+09		1	0.1	Pronamide	23950-58-5					8.8E+04	2.1E+05		6.2E+04
				1.3E-02	I						1.36E+09		1	0.1	Propachlor	1918-16-7					1.5E+04	3.6E+04		1.1E+04
				5.0E-03	I						1.36E+09		1	0.1	Propanil	709-98-8					5.8E+03	1.4E+04		4.1E+03
1.9E-01	O			4.0E-02	O						1.36E+09		1	0.1	Propargite	2312-35-8	1.7E+03	4.0E+03		1.2E+03	4.7E+04	1.1E+05		3.3E+04
				2.0E-03	I			V		1.11E+05	1.36E+09	6.27E+04	1		Propargyl Alcohol	107-19-7					2.3E+03			2.3E+03
				2.0E-02	I						1.36E+09		1	0.1	Propazine	139-40-2					2.3E+04	5.5E+04		1.6E+04
				2.0E-02	I						1.36E+09		1	0.1	Propham	122-42-9					2.3E+04	5.5E+04		1.6E+04
				1.0E-01	O						1.36E+09		1	0.1	Propiconazole	60207-90-1					1.2E+05	2.8E+05		8.2E+04
						8.0E-03	I	V		3.26E+04	1.36E+09	8.94E+03	1		Propionaldehyde	123-38-6							3.1E+02	3.1E+02
				1.0E-01	X	1.0E+00	X	V		2.64E+02	1.36E+09	6.99E+03	1		Propyl benzene	103-65-1					1.2E+05		3.1E+04	2.4E+04
						3.0E+00	C	V		3.49E+02	1.36E+09	7.04E+02	1		Propylene	115-07-1							9.3E+03	9.3E+03
				2.0E+01	P						1.36E+09		1	0.1	Propylene Glycol	57-55-6					2.3E+07	5.5E+07		1.6E+07
						2.7E-04	A				1.36E+09		1	0.1	Propylene Glycol Dinitrate	6423-43-4							1.6E+06	1.6E+06
				7.0E-01	H	2.0E+00	I	V		1.06E+05	1.36E+09	7.83E+04	1		Propylene Glycol Monomethyl Ether	107-98-2					8.2E+05		6.9E+05	3.7E+05
2.4E-01	I	3.7E-06	I			3.0E-02	I	V		7.77E+04	1.36E+09	1.03E+04	1		Propylene Oxide	75-56-9	1.4E+03		3.4E+03	9.7E+02			1.4E+03	1.4E+03
				1.0E-03	I			V		5.30E+05	1.36E+09	5.54E+04	1		Pyridine	110-86-1					1.2E+03			1.2E+03
				5.0E-04	I						1.36E+09		1	0.1	Quinalphos	13593-03-8					5.8E+02	1.4E+03		4.1E+02
3.0E+00	I										1.36E+09		1	0.1	Quinoline	91-22-5	1.1E+02	2.6E+02		7.7E+01				
				9.0E-03	I						1.36E+09		1	0.1	Quizalofop-ethyl	76578-14-8					1.1E+04	2.5E+04		7.4E+03
						3.0E+04	A						1		Refractory Ceramic Fibers (units in fibers)	E715557								
				3.0E-02	I						1.36E+09		1	0.1	Resmethrin	10453-86-8					3.5E+04	8.3E+04		2.5E+04
				5.0E-02	H			V			1.36E+09	4.65E+05	1		Ronnel	299-84-3					5.8E+04			5.8E+04
				4.0E-03	I						1.36E+09		1	0.1	Rotenone	83-79-4					4.7E+03	1.1E+04		3.3E+03
2.2E-01	C	6.3E-05	C						M		1.36E+09		1	0.1	Safrole	94-59-7	1.5E+03	3.5E+03	2.6E+07	1.0E+03				
				5.0E-03	I						1.36E+09		1		Selenious Acid	7783-00-8					5.8E+03			5.8E+03
				5.0E-03	I	2.0E-02	C				1.36E+09		1		Selenium	7782-49-2					5.8E+03		1.2E+08	5.8E+03
				5.0E-03	C	2.0E-02	C				1.36E+09		1		Selenium Sulfide	7446-34-6					5.8E+03		1.2E+08	5.8E+03
				1.4E-01	O						1.36E+09		1	0.1	Sethoxydim	74051-80-2					1.6E+05	3.9E+05		1.1E+05
						3.0E-03	C				1.36E+09		1		Silica (crystalline, respirable)	7631-86-9							1.8E+07	1.8E+07
				5.0E-03	I						1.36E+09		0.04		Silver	7440-22-4					5.8E+03			5.8E+03
1.2E-01	H			5.0E-03	I						1.36E+09		1	0.1	Simazine	122-34-9	2.7E+03	6.4E+03		1.9E+03	5.8E+03	1.4E+04		4.1E+03
				1.3E-02	I						1.36E+09		1	0.1	Sodium Acifluorfen	62476-59-9					1.5E+04	3.6E+04		1.1E+04
				4.0E-03	I						1.36E+09		1		Sodium Azide	26628-22-8					4.7E+03			4.7E+03
2.7E-01	H			3.0E-02	I						1.36E+09		1	0.1	Sodium Diethyldithiocarbamate	148-18-5	1.2E+03	2.9E+03		8.5E+02	3.5E+04	8.3E+04		2.5E+04
				5.0E-02	A	1.3E-02	C				1.36E+09		1		Sodium Fluoride	7681-49-4					5.8E+04		7.7E+07	5.8E+04
				2.0E-05	I						1.36E+09		1	0.1	Sodium Fluoroacetate	62-74-8					2.3E+01	5.5E+01		1.6E+01
				1.0E-03	H						1.36E+09		1		Sodium Metavanadate	13718-26-8					1.2E+03			1.2E+03
				8.0E-04	P						1.36E+09		1		Sodium Tungstate	13472-45-2					9.3E+02			9.3E+02
				8.0E-04	P						1.36E+09		1		Sodium Tungstate Dihydrate	10213-10-2					9.3E+02			9.3E+02

RMLS
TR=1E-04
THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
2.4E-02	H			3.0E-02	I						1.36E+09		1	0.1	Stirofos (Tetrachlorovinphos)	961-11-5	1.4E+04	3.2E+04		9.6E+03	3.5E+04	8.3E+04		2.5E+04
				6.0E-01	I						1.36E+09		1		Strontium, Stable	7440-24-6					7.0E+05			7.0E+05
				3.0E-04	I						1.36E+09		1	0.1	Strychnine	57-24-9					3.5E+02	8.3E+02		2.5E+02
				2.0E-01	I	1.0E+00	I	V		8.67E+02	1.36E+09	9.35E+03	1		Styrene	100-42-5					2.3E+05		4.1E+04	3.5E+04
				3.0E-03	P						1.36E+09		1	0.1	Styrene-Acrylonitrile (SAN) Trimer (THNA isomer)	57964-39-3					3.5E+03	8.3E+03		2.5E+03
				3.0E-03	P						1.36E+09		1	0.1	Styrene-Acrylonitrile (SAN) Trimer (THNP isomer)	57964-40-6					3.5E+03	8.3E+03		2.5E+03
				1.0E-03	P	2.0E-03	X				1.36E+09		1	0.1	Sulfolane	126-33-0					1.2E+03	2.8E+03	1.2E+07	8.2E+02
				8.0E-04	P						1.36E+09		1	0.1	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9					9.3E+02	2.2E+03		6.6E+02
						1.0E-03	C	V			1.36E+09		1		Sulfur Trioxide	7446-11-9							6.0E+06	6.0E+06
						1.0E-03	C				1.36E+09		1		Sulfuric Acid	7664-93-9							6.0E+06	6.0E+06
2.5E-02	I	7.1E-06	I	5.0E-02	H						1.36E+09		1	0.1	Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl	140-57-8	1.3E+04	3.1E+04	2.3E+08	9.2E+03	5.8E+04	1.4E+05		4.1E+04
				3.0E-02	H						1.36E+09		1	0.1	TCMTB	21564-17-0					3.5E+04	8.3E+04		2.5E+04
				7.0E-02	I						1.36E+09		1	0.1	Tebuthiuron	34014-18-1					8.2E+04	1.9E+05		5.7E+04
				2.0E-02	H						1.36E+09		1	0.1	Temephos	3383-96-8					2.3E+04	5.5E+04		1.6E+04
				1.3E-02	I						1.36E+09		1	0.1	Terbacil	5902-51-2					1.5E+04	3.6E+04		1.1E+04
				2.5E-05	H			V		3.09E+01	1.36E+09	2.64E+05	1		Terbufos	13071-79-9					2.9E+01			2.9E+01
				1.0E-03	I						1.36E+09		1	0.1	Terbutryn	886-50-0					1.2E+03	2.8E+03		8.2E+02
5.0E-03	C	1.3E-06	C					V			1.36E+09	3.99E+03	1		Tert-Butyl Acetate	540-88-5	6.5E+04		3.8E+03	3.6E+03				
				1.0E-04	I						1.36E+09		1	0.1	Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1					1.2E+02	2.8E+02		8.2E+01
				3.0E-05	P			V			1.36E+09	5.07E+04	1		Tetrachlorobenzene, 1,2,4,5-	95-94-3					3.5E+01			3.5E+01
2.6E-02	I	7.4E-06	I	3.0E-02	I			V		6.80E+02	1.36E+09	5.68E+03	1		Tetrachloroethane, 1,1,1,2-	630-20-6	1.3E+04		9.4E+02	8.8E+02	3.5E+04			3.5E+04
2.0E-01	I	5.8E-05	C	2.0E-02	I			V		1.90E+03	1.36E+09	1.51E+04	1		Tetrachloroethane, 1,1,2,2-	79-34-5	1.6E+03		3.2E+02	2.7E+02	2.3E+04			2.3E+04
2.1E-03	I	2.6E-07	I	6.0E-03	I	4.0E-02	I	V		1.66E+02	1.36E+09	2.35E+03	1		Tetrachloroethylene	127-18-4	1.6E+05		1.1E+04	1.0E+04	7.0E+03		4.1E+02	3.9E+02
				3.0E-02	I						1.36E+09		1	0.1	Tetrachlorophenol, 2,3,4,6-	58-90-2					3.5E+04	8.3E+04		2.5E+04
1.6E+01	X			6.0E-05	X			V			1.36E+09	1.05E+05	1		Tetrachlorotoluene, p- alpha, alpha, alpha-	5216-25-1	2.0E+01			2.0E+01	7.0E+01			7.0E+01
				5.0E-04	I						1.36E+09		1	0.1	Tetraethyl Dithiopyrophosphate	3689-24-5					5.8E+02	1.4E+03		4.1E+02
						8.0E+01	I	V		2.05E+03	1.36E+09	1.22E+03	1		Tetrafluoroethane, 1,1,1,2-	811-97-2							4.3E+05	4.3E+05
				1.0E-04	X						1.36E+09		1	0.1	Tetramethylphosphoramide, -N,N,N',N" (TMPA)	16853-36-4					1.2E+02	2.8E+02		8.2E+01
				2.0E-03	P						1.36E+09		1	0.00065	Tetryl (Trinitrophenylmethylnitramine)	479-45-8					2.3E+03	8.5E+05		2.3E+03
				2.0E-05	G						1.36E+09		1		Thallic Oxide	1314-32-5					2.3E+01			2.3E+01
				1.0E-05	X						1.36E+09		1		Thallium (I) Nitrate	10102-45-1					1.2E+01			1.2E+01
				1.0E-05	X						1.36E+09		1		Thallium (Soluble Salts)	7440-28-0					1.2E+01			1.2E+01
				1.0E-05	X			V			1.36E+09	1.40E+05	1		Thallium Acetate	563-68-8					1.2E+01			1.2E+01
				2.0E-05	X						1.36E+09		1	0.1	Thallium Carbonate	6533-73-9					2.3E+01	5.5E+01		1.6E+01
				1.0E-05	X						1.36E+09		1		Thallium Chloride	7791-12-0					1.2E+01			1.2E+01
				1.0E-05	G						1.36E+09		1		Thallium Selenite	12039-52-0					1.2E+01			1.2E+01
				2.0E-05	X						1.36E+09		1		Thallium Sulfate	7446-18-6					2.3E+01			2.3E+01
				4.3E-02	O						1.36E+09		1	0.1	Thifensulfuron-methyl	79277-27-3					5.0E+04	1.2E+05		3.5E+04
				1.0E-02	I						1.36E+09		1	0.1	Thiobencarb	28249-77-6					1.2E+04	2.8E+04		8.2E+03
				7.0E-02	X						1.36E+09		1	0.0075	Thiodiglycol	111-48-8					8.2E+04	2.6E+06		7.9E+04
				3.0E-04	H						1.36E+09		1	0.1	Thiofanox	39196-18-4					3.5E+02	8.3E+02		2.5E+02
1.2E-02	O			1.6E-01	O						1.36E+09		1	0.1	Thiophanate, Methyl	23564-05-8	2.8E+04	6.7E+04		2.0E+04	1.9E+05	4.4E+05		1.3E+05
				1.5E-02	O						1.36E+09		1	0.1	Thiram	137-26-8					1.8E+04	4.1E+04		1.2E+04
				6.0E-01	H						1.36E+09		1		Tin	7440-31-5					7.0E+05			7.0E+05
						1.0E-04	A	V			1.36E+09		1		Titanium Tetrachloride	7550-45-0							6.0E+05	6.0E+05
				8.0E-02	I	5.0E+00	I	V		8.18E+02	1.36E+09	4.29E+03	1		Toluene	108-88-3					9.3E+04		9.4E+04	4.7E+04
3.9E-02	C	1.1E-05	C			8.0E-06	C	V			1.36E+09	7.62E+05	1		Toluene-2,4-diisocyanate	584-84-9	8.4E+03		8.5E+04	7.6E+03			2.7E+01	2.7E+01
1.8E-01	X			2.0E-04	X						1.36E+09		1	0.1	Toluene-2,5-diamine	95-70-5	1.8E+03	4.3E+03		1.3E+03	2.3E+02	5.5E+02		1.6E+02
3.9E-02	C	1.1E-05	C			8.0E-06	C	V		1.71E+03	1.36E+09	6.32E+05	1		Toluene-2,6-diisocyanate	91-08-7	8.4E+03		7.0E+04	7.5E+03			2.2E+01	2.2E+01
				1.0E-04	X						1.36E+09		1	0.1	Toluenediamine, 2,3-	2687-25-4					1.2E+02	2.8E+02		8.2E+01
				1.0E-04	X						1.36E+09		1	0.1	Toluenediamine, 3,4-	496-72-0					1.2E+02	2.8E+02		8.2E+01
				5.0E-03	P						1.36E+09		1	0.1	Toluic Acid, p-	99-94-5					5.8E+03	1.4E+04		4.1E+03
1.6E-02	P	5.1E-05	C								1.36E+09		1	0.1	Toluidine, o- (Methylaniline, 2-)	95-53-4	2.0E+04	4.8E+04	3.3E+07	1.4E+04				
3.0E-02	P			4.0E-03	X						1.36E+09		1	0.1	Toluidine, p-	106-49-0	1.1E+04	2.6E+04		7.7E+03	4.7E+03	1.1E+04		3.3E+03
				3.0E+00	P			V		3.42E-01	1.36E+09	1.38E+03	1		Total Petroleum Hydrocarbons (Aliphatic High)	E1790670					3.5E+06			3.5E+06
						6.0E-01	P	V		1.41E+02	1.36E+09	8.29E+02	1		Total Petroleum Hydrocarbons (Aliphatic Low)	E1790666							2.2E+03	2.2E+03
				1.0E-02	X	1.0E-01	P	V		6.86E+00	1.36E+09	1.04E+03	1		Total Petroleum Hydrocarbons (Aliphatic Medium)	E1790668					1.2E+04		4.6E+02	4.4E+02
				4.0E-02	P						1.36E+09		1	0.13	Total Petroleum Hydrocarbons (Aromatic High)	E1790676					4.7E+04	8.5E+04		3.0E+04
				4.0E-03	P	3.0E-02	P	V		1.82E+03	1.36E+09	3.54E+03	1		Total Petroleum Hydrocarbons (Aromatic Low)	E1790672					4.7E+03		4.6E+02	4.2E+02
				4.0E-03	P	3.0E-03	P	V			1.36E+09	5.24E+04	1	0.13	Total Petroleum Hydrocarbons (Aromatic Medium)	E1790674					4.7E+03	8.5E+03	6.9E+02	5.6E+02
1.1E+00	I	3.2E-04	I	9.0E-05	P						1.36E+09		1	0.1	Toxaphene	8001-35-2	3.0E+02	7.0E+02	5.2E+06	2.1E+02	1.1E+02	2.5E+02		7.4E+01
				3.0E-05	X						1.36E+09		1	0.1	Toxaphene, Weathered	E1841606					3.5E+01	8.3E+01		2.5E+01
				7.5E-03	I						1.36E+09		1	0.1	Tralomethrin	66841-25-6					8.8E+03	2.1E+04		6.2E+03
				3.0E-04	A			V			1.36E+09	3.36E+03	1		Tri-n-butyltin	688-73-3					3.5E+02			3.5E+02
				8.0E+01	X						1.36E+09		1	0.1	Triacetin	102-76-1					9.3E+07	2.2E+08		6.6E+07
				3.4E-02	O						1.36E+09		1	0.1	Triadimefon	43121-43-3					4.0E+04	9.4E+04		2.8E+04
7.2E-02	O			2.5E-02	O			V			1.36E+09	3.62E+05	1		Triallate	2303-17-5	4.6E+03			4.6E+03	2.9E+04			2.9E+04
				1.0E-02	I						1.36E+09		1	0.1	Triasulfuron	82097-50-5					1.2E+04	2.8E+04		8.2E+03
				8.0E-03	I						1.36E+09		1	0.1	Tribenuron-methyl	101200-48-0					9.3E+03	2.2E+04		6.6E+03
				5.0E-03	I			V			1.36E+09	4.83E+04	1		Tribromobenzene, 1,2,4-	615-54-3					5.8E+03			5.8E+03
				9.0E-03	X						1.36E+09		1	0.1	Tribromophenol, 2,4,6-	118-79-6					1.1E+04	2.5E+04		7.4E+03
				2.0E-04	O						1.36E+09		1	0.1	Tribufos	78-48-8					2.3E+02	5.5E+02		1.6E+02
9.0E-03	P			1.0E-02	P						1.36E+09		1	0.1	Tributyl Phosphate	126-73-8	3.6E+04	8.6E+04		2.6E+04	1.2E+04	2.8E+04		8.2E+03
				3.0E-04	P						1.36E+09		1	0.1	Tributyltin Compounds	E1790679					3.5E+02	8.3E+02		2.5E+02
				3.0E-04	I						1.36E+09		1	0.1	Tributyltin Oxide	56-35-9					3.5E+02	8.3E+02		2.5E+02
											1.36E+09		1	0.1	Trichloramine	10025-85-1								
				3.0E+01	I	5.0E+00	P	V		9.10E+02	1.36E+09	1.29E+03	1		Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1					3.5E+07		2.8E+04	2.8E+04
7.0E-02	I			2.0E-02	I						1.36E+09		1	0.1	Trichloroacetic Acid	76-03-9	4.7E+03	1.1E+04		3.3E+03	2.3E+04	5.5E+04		1.6E+04

RMLS TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information																Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day)$^{-1}$	key	IUR (ug/m^3)$^{-1}$	key	RfD_o (mg/kg-day)	key	RfC_i (mg/m^3)	key	vol	mutagen	C_{sat} (mg/kg)	PEF (m^3/kg)	VF (m^3/kg)	GIABS	ABS_d	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
2.9E-02	H										1.36E+09		1	0.1	Trichloroaniline HCl, 2,4,6-	33663-50-2	1.1E+04	2.7E+04		7.9E+03				
7.0E-03	X			3.0E-05	X						1.36E+09		1	0.1	Trichloroaniline, 2,4,6-	634-93-5	4.7E+04	1.1E+05		3.3E+04	3.5E+01	8.3E+01		2.5E+01
				8.0E-04	X			V			1.36E+09	3.22E+04	1		Trichlorobenzene, 1,2,3-	87-61-6					9.3E+02			9.3E+02
2.9E-02	P			1.0E-02	I	2.0E-03	P	V		4.04E+02	1.36E+09	2.99E+04	1		Trichlorobenzene, 1,2,4-	120-82-1	1.1E+04			1.1E+04	1.2E+04		2.6E+02	2.6E+02
				2.0E+00	I	5.0E+00	I	V		6.40E+02	1.36E+09	1.65E+03	1		Trichloroethane, 1,1,1-	71-55-6					2.3E+06		3.6E+04	3.6E+04
5.7E-02	I	1.6E-05	I	4.0E-03	I	2.0E-04	X	V		2.16E+03	1.36E+09	7.22E+03	1		Trichloroethane, 1,1,2-	79-00-5	5.7E+03		5.5E+02	5.0E+02	4.7E+03		6.3E+00	6.3E+00
4.6E-02	I	4.1E-06	I	5.0E-04	I	2.0E-03	I	V	M	6.92E+02	1.36E+09	2.21E+03	1		Trichloroethylene	79-01-6	7.1E+03		6.6E+02	6.0E+02	5.8E+02		1.9E+01	1.9E+01
				3.0E-01	I			V		1.23E+03	1.36E+09	1.04E+03	1		Trichlorofluoromethane	75-69-4					3.5E+05			3.5E+05
				1.0E-01	I						1.36E+09		1	0.1	Trichlorophenol, 2,4,5-	95-95-4					1.2E+05	2.8E+05		8.2E+04
1.1E-02	I	3.1E-06	I	1.0E-03	P						1.36E+09		1	0.1	Trichlorophenol, 2,4,6-	88-06-2	3.0E+04	7.0E+04	5.4E+08	2.1E+04	1.2E+03	2.8E+03		8.2E+02
				1.0E-02	I						1.36E+09		1	0.1	Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5					1.2E+04	2.8E+04		8.2E+03
				8.0E-03	I						1.36E+09		1	0.1	Trichlorophenoxypropionic acid, -2,4,5	93-72-1					9.3E+03	2.2E+04		6.6E+03
				5.0E-03	I			V		1.28E+03	1.36E+09	1.50E+04	1		Trichloropropane, 1,1,2-	598-77-6					5.8E+03			5.8E+03
3.0E+01	I			4.0E-03	I	3.0E-04	I	V	M	1.40E+03	1.36E+09	1.57E+04	1		Trichloropropane, 1,2,3-	96-18-4	1.1E+01			1.1E+01	4.7E+03		2.1E+01	2.1E+01
				3.0E-03	X	3.0E-04	P	V		3.11E+02	1.36E+09	2.34E+03	1		Trichloropropene, 1,2,3-	96-19-5					3.5E+03		3.1E+00	3.1E+00
				2.0E-02	A						1.36E+09		1	0.1	Tricresyl Phosphate (TCP)	1330-78-5					2.3E+04	5.5E+04		1.6E+04
				3.0E-03	I						1.36E+09		1	0.1	Tridiphane	58138-08-2					3.5E+03	8.3E+03		2.5E+03
						7.0E-03	I	V		2.79E+04	1.36E+09	1.58E+04	1		Triethylamine	121-44-8							4.8E+02	4.8E+02
				2.0E+00	P						1.36E+09		1	0.1	Triethylene Glycol	112-27-6					2.3E+06	5.5E+06		1.6E+06
						2.0E+01	P	V		4.81E+03	1.36E+09	7.12E+02	1		Trifluoroethane, 1,1,1-	420-46-2							6.2E+04	6.2E+04
7.7E-03	I			7.5E-03	I			V			1.36E+09	5.13E+05	1		Trifluralin	1582-09-8	4.2E+04			4.2E+04	8.8E+03			8.8E+03
2.0E-02	P			1.0E-02	P						1.36E+09		1	0.1	Trimethyl Phosphate	512-56-1	1.6E+04	3.9E+04		1.1E+04	1.2E+04	2.8E+04		8.2E+03
				1.0E-02	I	6.0E-02	I	V		2.93E+02	1.36E+09	9.44E+03	1		Trimethylbenzene, 1,2,3-	526-73-8					1.2E+04		2.5E+03	2.0E+03
				1.0E-02	I	6.0E-02	I	V		2.19E+02	1.36E+09	7.91E+03	1		Trimethylbenzene, 1,2,4-	95-63-6					1.2E+04		2.1E+03	1.8E+03
				1.0E-02	I	6.0E-02	I	V		1.82E+02	1.36E+09	6.61E+03	1		Trimethylbenzene, 1,3,5-	108-67-8					1.2E+04		1.7E+03	1.5E+03
				1.0E-02	X			V		2.96E+01	1.36E+09	1.00E+03	1		Trimethylpentene, 2,4,4-	25167-70-8					1.2E+04			1.2E+04
				3.0E-02	I						1.36E+09		1	0.019	Trinitrobenzene, 1,3,5-	99-35-4					3.5E+04	4.4E+05		3.2E+04
3.0E-02	I			5.0E-04	I						1.36E+09		1	0.032	Trinitrotoluene, 2,4,6-	118-96-7	1.1E+04	8.0E+04		9.6E+03	5.8E+02	4.3E+03		5.1E+02
				2.0E-02	P						1.36E+09		1	0.1	Triphenylphosphine Oxide	791-28-6					2.3E+04	5.5E+04		1.6E+04
				2.0E-02	A						1.36E+09		1	0.1	Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8					2.3E+04	5.5E+04		1.6E+04
				1.0E-02	X						1.36E+09		1	0.1	Tris(1-chloro-2-propyl)phosphate	13674-84-5					1.2E+04	2.8E+04		8.2E+03
2.3E+00	C	6.6E-04	C					V		4.67E+02	1.36E+09	9.03E+05	1		Tris(2,3-dibromopropyl)phosphate	126-72-7	1.4E+02		1.7E+03	1.3E+02				
2.0E-02	P			7.0E-03	P						1.36E+09		1	0.1	Tris(2-chloroethyl)phosphate	115-96-8	1.6E+04	3.9E+04		1.1E+04	8.2E+03	1.9E+04		5.7E+03
3.2E-03	P			1.0E-01	P						1.36E+09		1	0.1	Tris(2-ethylhexyl)phosphate	78-42-2	1.0E+05	2.4E+05		7.2E+04	1.2E+05	2.8E+05		8.2E+04
				8.0E-04	P						1.36E+09		1		Tungsten	7440-33-7					9.3E+02			9.3E+02
				2.0E-04	A	4.0E-05	A				1.36E+09		1		Uranium	7440-61-1					2.3E+02		2.4E+05	2.3E+02
1.0E+00	C	2.9E-04	C						M		1.36E+09		1	0.1	Urethane	51-79-6	3.3E+02	7.7E+02	5.7E+06	2.3E+02				
		8.3E-03	P	9.0E-03	I	7.0E-06	P				1.36E+09		0.026		Vanadium Pentoxide	1314-62-1			2.0E+05	2.0E+05	1.1E+04		4.2E+04	8.4E+03
				5.0E-03	G	1.0E-04	A				1.36E+09		0.026		Vanadium and Compounds	7440-62-2					5.9E+03		6.0E+05	5.8E+03
				1.0E-03	I			V			1.36E+09	1.23E+05	1		Vernolate	1929-77-7					1.2E+03			1.2E+03
				1.2E-03	O						1.36E+09		1	0.1	Vinclozolin	50471-44-8					1.4E+03	3.3E+03		9.8E+02
				1.0E+00	H	2.0E-01	I	V		2.75E+03	1.36E+09	4.40E+03	1		Vinyl Acetate	108-05-4					1.2E+06		3.9E+03	3.8E+03
		1.5E-05	P			3.0E-03	I	V		2.47E+03	1.36E+09	1.37E+03	1		Vinyl Bromide	593-60-2			1.1E+02	1.1E+02			1.8E+01	1.8E+01
7.2E-01	I	4.4E-06	I	3.0E-03	I	8.0E-02	A	V	M	3.92E+03	1.36E+09	9.56E+02	1		Vinyl Chloride	75-01-4	4.5E+02		2.7E+02	1.7E+02	3.5E+03		3.3E+02	3.1E+02
				3.0E-04	I						1.36E+09		1	0.1	Warfarin	81-81-2					3.5E+02	8.3E+02		2.5E+02
				2.0E-01	G	1.0E-01	G	V		3.88E+02	1.36E+09	5.47E+03	1		Xylene, m-	108-38-3					2.3E+05		2.4E+03	2.4E+03
				2.0E-01	G	1.0E-01	G	V		4.34E+02	1.36E+09	6.45E+03	1		Xylene, o-	95-47-6					2.3E+05		2.8E+03	2.8E+03
				2.0E-01	G	1.0E-01	G	V		3.90E+02	1.36E+09	5.58E+03	1		Xylene, p-	106-42-3					2.3E+05		2.4E+03	2.4E+03
				2.0E-01	I	1.0E-01	I	V		2.60E+02	1.36E+09	5.74E+03	1		Xylenes	1330-20-7					2.3E+05		2.5E+03	2.5E+03
				3.0E-04	I						1.36E+09		1		Zinc Phosphide	1314-84-7					3.5E+02			3.5E+02
				3.0E-01	I						1.36E+09		1		Zinc and Compounds	7440-66-6					3.5E+05			3.5E+05
				5.0E-02	I						1.36E+09		1	0.1	Zineb	12122-67-7					5.8E+04	1.4E+05		4.1E+04
				8.0E-05	X						1.36E+09		1		Zirconium	7440-67-7					9.3E+01			9.3E+01

RMS
TR=1E-04
THQ=1

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
						Acephate	30560-19-1		
2.2E-06	I	9.0E-03	I	V		Acetaldehyde	75-07-0	5.6E+02	3.9E+01
						Acetochlor	34256-82-1		
				V		Acetone	67-64-1		
		2.0E-03	X			Acetone Cyanohydrin	75-86-5		8.8E+00
		6.0E-02	I	V		Acetonitrile	75-05-8		2.6E+02
				V		Acetophenone	98-86-2		
1.3E-03	C					Acetylaminofluorene, 2-	53-96-3	9.4E-01	
		2.0E-05	I	V		Acrolein	107-02-8		8.8E-02
1.0E-04	I	6.0E-03	I		M	Acrylamide	79-06-1	1.2E+01	2.6E+01
		2.0E-04	P	V		Acrylic Acid	79-10-7		8.8E-01
6.8E-05	I	2.0E-03	I	V		Acrylonitrile	107-13-1	1.8E+01	8.8E+00
		6.0E-03	P			Adiponitrile	111-69-3		2.6E+01
						Alachlor	15972-60-8		
						Aldicarb	116-06-3		
						Aldicarb Sulfone	1646-88-4		
						Aldicarb sulfoxide	1646-87-3		
4.9E-03	I			V		Aldrin	309-00-2	2.5E-01	
		1.0E-04	X	V		Allyl Alcohol	107-18-6		4.4E-01
6.0E-06	C	1.0E-03	I	V		Allyl Chloride	107-05-1	2.0E+02	4.4E+00
		5.0E-03	P			Aluminum	7429-90-5		2.2E+01
						Aluminum Phosphide	20859-73-8		
						Ametryn	834-12-8		
6.0E-03	C					Aminobiphenyl, 4-	92-67-1	2.0E-01	
						Aminophenol, m-	591-27-5		
						Aminophenol, o-	95-55-6		
						Aminophenol, p-	123-30-8		
						Amitraz	33089-61-1		
		5.0E-01	I	V		Ammonia	7664-41-7		2.2E+03
						Ammonium Picrate	131-74-8		
						Ammonium Sulfamate	7773-06-0		
		3.0E-03	X	V		Amyl Alcohol, tert-	75-85-4		1.3E+01
1.6E-06	C	1.0E-03	I			Aniline	62-53-3	7.7E+02	4.4E+00
						Anthraquinone, 9,10-	84-65-1		
		3.0E-04	A			Antimony (metallic)	7440-36-0		1.3E+00
						Antimony Pentoxide	1314-60-9		
						Antimony Tetroxide	1332-81-6		
		2.0E-04	I			Antimony Trioxide	1309-64-4		8.8E-01
4.3E-03	I	1.5E-05	C			Arsenic, Inorganic	7440-38-2	2.9E-01	6.6E-02
		5.0E-05	I			Arsine	7784-42-1		2.2E-01
						Asbestos (units in fibers)	1332-21-4		
						Asulam	3337-71-1		
						Atrazine	1912-24-9		
2.5E-04	C					Auramine	492-80-8	4.9E+00	
						Avermectin B1	65195-55-3		
		1.0E-02	A			Azinphos-methyl	86-50-0		4.4E+01
3.1E-05	I			V		Azobenzene	103-33-3	4.0E+01	
		7.0E-06	P			Azodicarbonamide	123-77-3		3.1E-02
		5.0E-04	H			Barium	7440-39-3		2.2E+00
				V		Benfluralin	1861-40-1		
						Benomyl	17804-35-2		
						Bensulfuron-methyl	83055-99-6		
						Bentazon	25057-89-0		
				V		Benzaldehyde	100-52-7		
7.8E-06	I	3.0E-02	I	V		Benzene	71-43-2	1.6E+02	1.3E+02
						Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1		
				V		Benzenethiol	108-98-5		
6.7E-02	I				M	Benzidine	92-87-5	1.8E-02	
						Benzoic Acid	65-85-0		
				V		Benzotrichloride	98-07-7		
						Benzyl Alcohol	100-51-6		
4.9E-05	C	1.0E-03	P	V		Benzyl Chloride	100-44-7	2.5E+01	4.4E+00
2.4E-03	I	2.0E-05	I			Beryllium and compounds	7440-41-7	5.1E-01	8.8E-02
						Bifenox	42576-02-3		
						Biphenthrin	82657-04-3		
		4.0E-04	X	V		Biphenyl, 1,1'-	92-52-4		1.8E+00
				V		Bis(2-chloro-1-methylethyl) ether	108-60-1		
						Bis(2-chloroethoxy)methane	111-91-1		
3.3E-04	I			V		Bis(2-chloroethyl)ether	111-44-4	3.7E+00	
6.2E-02	I			V		Bis(chloromethyl)ether	542-88-1	2.0E-02	
						Bisphenol A	80-05-7		
		2.0E-02	H			Boron And Borates Only	7440-42-8		8.8E+01
		2.0E-02	P	V		Boron Trichloride	10294-34-5		8.8E+01
		1.3E-02	C	V		Boron Trifluoride	7637-07-2		5.7E+01
						Bromate	15541-45-4		
		6.0E-05	X	V		Bromo-2-chloroethane, 1-	107-04-0		2.6E-01
				V		Bromo-3-fluorobenzene, 1-	1073-06-9		
				V		Bromo-4-fluorobenzene, 1-	460-00-4		
						Bromoacetic acid	79-08-3		
		6.0E-02	I	V		Bromobenzene	108-86-1		2.6E+02
		4.0E-02	X	V		Bromochloromethane	74-97-5		1.8E+02
3.7E-05	C			V		Bromodichloromethane	75-27-4	3.3E+01	
1.1E-06	I			V		Bromoform	75-25-2	1.1E+03	
		5.0E-03	I	V		Bromomethane	74-83-9		2.2E+01
				V		Bromophos	2104-96-3		
		1.0E-01	A	V		Bromopropane, 1-	106-94-5		4.4E+02
						Bromoxynil	1689-84-5		
				V		Bromoxynil Octanoate	1689-99-2		
3.0E-05	I	2.0E-03	I	V		Butadiene, 1,3-	106-99-0	4.1E+01	8.8E+00
				V		Butanol, N-	71-36-3		

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
		5.0E+00	I	V		Butyl Alcohol, t-	75-65-0		2.2E+04
		3.0E+01	P	V		Butyl alcohol, sec-	78-92-2		1.3E+05
				V		Butylate	2008-41-5		
5.7E-08	C					Butylated hydroxyanisole	25013-16-5	2.2E+04	
						Butylated hydroxytoluene	128-37-0		
				V		Butylbenzene, n-	104-51-8		
				V		Butylbenzene, sec-	135-98-8		
				V		Butylbenzene, tert-	98-06-6		
						Cacodylic Acid	75-60-5		
1.8E-03	I	1.0E-05	A			Cadmium (Diet)	7440-43-9	6.8E-01	4.4E-02
1.8E-03	I	1.0E-05	A			Cadmium (Water)	7440-43-9	6.8E-01	4.4E-02
		2.2E-03	C			Caprolactam	105-60-2		9.6E+00
4.3E-05	C					Captafol	2425-06-1	2.9E+01	
6.6E-07	C					Captan	133-06-2	1.9E+03	
						Carbaryl	63-25-2		
						Carbofuran	1563-66-2		
		7.0E-01	I	V		Carbon Disulfide	75-15-0		3.1E+03
6.0E-06	I	1.0E-01	I	V		Carbon Tetrachloride	56-23-5	2.0E+02	4.4E+02
		1.0E-01	P	V		Carbonyl Sulfide	463-58-1		4.4E+02
						Carbosulfan	55285-14-8		
						Carboxin	5234-68-4		
		9.0E-04	I			Ceric oxide	1306-38-3		3.9E+00
				V		Chloral Hydrate	302-17-0		
						Chloramben	133-90-4		
						Chloramines, Organic	E701235		
						Chloranil	118-75-2		
				V		Chlordane (alpha)	5103-71-9		
				V		Chlordane (gamma)	5103-74-2		
1.0E-04	I	7.0E-04	I	V		Chlordane (technical mixture)	12789-03-6	1.2E+01	3.1E+00
4.6E-03	C					Chlordecone (Kepone)	143-50-0	2.7E-01	
						Chlorfenvinphos	470-90-6		
						Chlorimuron, Ethyl-	90982-32-4		
		1.5E-04	A	V		Chlorine	7782-50-5		6.4E-01
		2.0E-04	I	V		Chlorine Dioxide	10049-04-4		8.8E-01
						Chlorite (Sodium Salt)	7758-19-2		
		5.0E+01	I	V		Chloro-1,1-difluoroethane, 1-	75-68-3		2.2E+05
3.0E-04	I	2.0E-02	I	V		Chloro-1,3-butadiene, 2-	126-99-8	4.1E+00	8.8E+01
						Chloro-2-methylaniline HCl, 4-	3165-93-3		
7.7E-05	C					Chloro-2-methylaniline, 4-	95-69-2	1.6E+01	
				V		Chloroacetaldehyde, 2-	107-20-0		
						Chloroacetic Acid	79-11-8		
		3.0E-05	I			Chloroacetophenone, 2-	532-27-4		1.3E-01
						Chloroaniline, p-	106-47-8		
		5.0E-02	P	V		Chlorobenzene	108-90-7		2.2E+02
						Chlorobenzene sulfonic acid, p-	98-66-8		
3.1E-05	C					Chlorobenzilate	510-15-6	4.0E+01	
						Chlorobenzoic Acid, p-	74-11-3		
8.6E-06	C	3.0E-01	P	V		Chlorobenzotrifluoride, 4-	98-56-6	1.4E+02	1.3E+03
				V		Chlorobutane, 1-	109-69-3		
		5.0E+01	I	V		Chlorodifluoromethane	75-45-6		2.2E+05
				V		Chloroethanol, 2-	107-07-3		
2.3E-05	I	9.8E-02	A	V		Chloroform	67-66-3	5.3E+01	4.3E+02
		9.0E-02	I	V		Chloromethane	74-87-3		3.9E+02
6.9E-04	C			V		Chloromethyl Methyl Ether	107-30-2	1.8E+00	
		1.0E-05	X			Chloronitrobenzene, o-	88-73-3		4.4E-02
		2.0E-03	P			Chloronitrobenzene, p-	100-00-5		8.8E+00
				V		Chlorophenol, 2-	95-57-8		
		4.0E-04	C	V		Chloropicrin	76-06-2		1.8E+00
						Chlorothalonil	1897-45-6		
				V		Chlorotoluene, o-	95-49-8		
				V		Chlorotoluene, p-	106-43-4		
6.9E-02	C					Chlorozotocin	54749-90-5	1.8E-02	
						Chlorpropham	101-21-3		
						Chlorpyrifos	2921-88-2		
						Chlorpyrifos Methyl	5598-13-0		
						Chlorsulfuron	64902-72-3		
						Chlorthal-dimethyl	1861-32-1		
						Chlorthiophos	60238-56-4		
						Chromium(III), Insoluble Salts	16065-83-1		
8.4E-02	G	1.0E-04	I		M	Chromium(VI)	18540-29-9	1.5E-02	4.4E-01
						Chromium, Total	7440-47-3		
						Clofentezine	74115-24-5		
9.0E-03	P	6.0E-06	P			Cobalt	7440-48-4	1.4E-01	2.6E-02
6.2E-04	I			V	M	Coke Oven Emissions	E649830	2.0E+00	
						Copper	7440-50-8		
		6.0E-01	C			Cresol, m-	108-39-4		2.6E+03
		6.0E-01	C			Cresol, o-	95-48-7		2.6E+03
		6.0E-01	C			Cresol, p-	106-44-5		2.6E+03
						Cresol, p-chloro-m-	59-50-7		
		6.0E-01	C			Cresols	1319-77-3		2.6E+03
				V		Crotonaldehyde, trans-	123-73-9		
		4.0E-01	I	V		Cumene	98-82-8		1.8E+03
6.3E-05	C					Cupferron	135-20-6	1.9E+01	
						Cyanazine	21725-46-2		
						Cyanides			
						~Calcium Cyanide	592-01-8		
						~Copper Cyanide	544-92-3		
		8.0E-04	G	V		~Cyanide (CN-)	57-12-5		3.5E+00
				V		~Cyanogen	460-19-5		
				V		~Cyanogen Bromide	506-68-3		

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
				V		~Cyanogen Chloride	506-77-4		
		8.0E-04	I	V		~Hydrogen Cyanide	74-90-8		3.5E+00
						~Potassium Cyanide	151-50-8		
						~Potassium Silver Cyanide	506-61-6		
						~Silver Cyanide	506-64-9		
						~Sodium Cyanide	143-33-9		
						~Thiocyanates	E1790665		
				V		~Thiocyanic Acid	463-56-9		
						~Zinc Cyanide	557-21-1		
		6.0E+00	I	V		Cyclohexane	110-82-7		2.6E+04
						Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3		
		7.0E-01	P	V		Cyclohexanone	108-94-1		3.1E+03
		1.0E+00	X	V		Cyclohexene	110-83-8		4.4E+03
				V		Cyclohexylamine	108-91-8		
						Cyfluthrin	68359-37-5		
						Cypermethrin	52315-07-8		
						Cyromazine	66215-27-8		
6.9E-05	C					DDD, p,p`- (DDD)	72-54-8	1.8E+01	
9.7E-05	C			V		DDE, p,p'-	72-55-9	1.3E+01	
9.7E-05	I					DDT	50-29-3	1.3E+01	
						Dalapon	75-99-0		
5.1E-06	C					Daminozide	1596-84-5	2.4E+02	
						Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5		
						Demeton	8065-48-3		
						Di(2-ethylhexyl)adipate	103-23-1		
						Diallate	2303-16-4		
						Diazinon	333-41-5		
				V		Dibenzothiophene	132-65-0		
6.0E-03	P	2.0E-04	I	V	M	Dibromo-3-chloropropane, 1,2-	96-12-8	2.0E-01	8.8E-01
						Dibromoacetic acid	631-64-1		
				V		Dibromobenzene, 1,3-	108-36-1		
				V		Dibromobenzene, 1,4-	106-37-6		
				V		Dibromochloromethane	124-48-1		
6.0E-04	I	9.0E-03	I	V		Dibromoethane, 1,2-	106-93-4	2.0E+00	3.9E+01
		4.0E-03	X	V		Dibromomethane (Methylene Bromide)	74-95-3		1.8E+01
						Dibutyltin Compounds	E1790661		
						Dicamba	1918-00-9		
						Dichloramine	3400-09-7		
4.2E-03	P			V		Dichloro-2-butene, 1,4-	764-41-0	2.9E-01	
4.2E-03	P			V		Dichloro-2-butene, cis-1,4-	1476-11-5	2.9E-01	
4.2E-03	P			V		Dichloro-2-butene, trans-1,4-	110-57-6	2.9E-01	
						Dichloroacetic Acid	79-43-6		
		2.0E-01	H	V		Dichlorobenzene, 1,2-	95-50-1		8.8E+02
1.1E-05	C	8.0E-01	I	V		Dichlorobenzene, 1,4-	106-46-7	1.1E+02	3.5E+03
3.4E-04	C					Dichlorobenzidine, 3,3'-	91-94-1	3.6E+00	
						Dichlorobenzophenone, 4,4'-	90-98-2		
		1.0E-01	X	V		Dichlorodifluoromethane	75-71-8		4.4E+02
1.6E-06	C			V		Dichloroethane, 1,1-	75-34-3	7.7E+02	
2.6E-05	I	7.0E-03	P	V		Dichloroethane, 1,2-	107-06-2	4.7E+01	3.1E+01
		2.0E-01	I	V		Dichloroethylene, 1,1-	75-35-4		8.8E+02
				V		Dichloroethylene, cis-1,2-	156-59-2		
		4.0E-02	X	V		Dichloroethylene, trans-1,2-	156-60-5		1.8E+02
						Dichlorophenol, 2,4-	120-83-2		
						Dichlorophenoxy Acetic Acid, 2,4-	94-75-7		
3.7E-06	P	4.0E-03	I	V		Dichloropropane, 1,2-	78-87-5	3.3E+02	1.8E+01
				V		Dichloropropane, 1,3-	142-28-9		
						Dichloropropanol, 2,3-	616-23-9		
4.0E-06	I	2.0E-02	I	V		Dichloropropene, 1,3-	542-75-6	3.1E+02	8.8E+01
8.3E-05	C	5.0E-04	I			Dichlorvos	62-73-7	1.5E+01	2.2E+00
						Dicrotophos	141-66-2		
		3.0E-04	X	V		Dicyclopentadiene	77-73-6		1.3E+00
4.6E-03	I					Dieldrin	60-57-1	2.7E-01	
3.0E-04	C	5.0E-03	I			Diesel Engine Exhaust	E17136615	4.1E+00	2.2E+01
		2.0E-04	P			Diethanolamine	111-42-2		8.8E-01
		1.0E-04	P			Diethylene Glycol Monobutyl Ether	112-34-5		4.4E-01
		3.0E-04	P			Diethylene Glycol Monoethyl Ether	111-90-0		1.3E+00
				V		Diethylformamide	617-84-5		
1.0E-01	C					Diethylstilbestrol	56-53-1	1.2E-02	
						Difenzoquat	43222-48-6		
						Diflubenzuron	35367-38-5		
		4.0E+01	I	V		Difluoroethane, 1,1-	75-37-6		1.8E+05
		3.0E+01	X	V		Difluoropropane, 2,2-	420-45-1		1.3E+05
1.3E-05	C			V		Dihydrosafrole	94-58-6	9.4E+01	
		7.0E-01	P	V		Diisopropyl Ether	108-20-3		3.1E+03
				V		Diisopropyl Methylphosphonate	1445-75-6		
						Dimethipin	55290-64-7		
						Dimethoate	60-51-5		
						Dimethoxybenzidine, 3,3'-	119-90-4		
						Dimethyl methylphosphonate	756-79-6		
1.3E-03	C					Dimethylamino azobenzene [p-]	60-11-7	9.4E-01	
						Dimethylaniline HCl, 2,4-	21436-96-4		
						Dimethylaniline, 2,4-	95-68-1		
				V		Dimethylaniline, N,N-	121-69-7		
						Dimethylbenzidine, 3,3'-	119-93-7		
		3.0E-02	I	V		Dimethylformamide	68-12-2		1.3E+02
		2.0E-06	X	V		Dimethylhydrazine, 1,1-	57-14-7		8.8E-03
1.6E-01	C			V		Dimethylhydrazine, 1,2-	540-73-8	7.7E-03	
						Dimethylphenol, 2,4-	105-67-9		
						Dimethylphenol, 2,6-	576-26-1		
						Dimethylphenol, 3,4-	95-65-8		

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 $(ug\ or\ fibers/m^3)$
1.3E-05	C			V		Dimethylvinylchloride	513-37-1	9.4E+01	
						Dinitro-o-cresol, 4,6-	534-52-1		
						Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5		
		2.0E-03	X			Dinitroaniline, 3,5-	618-87-1		8.8E+00
						Dinitrobenzene, 1,2-	528-29-0		
						Dinitrobenzene, 1,3-	99-65-0		
						Dinitrobenzene, 1,4-	100-25-4		
						Dinitrophenol, 2,4-	51-28-5		
						Dinitrotoluene Mixture, 2,4/2,6-	E1615210		
8.9E-05	C					Dinitrotoluene, 2,4-	121-14-2	1.4E+01	
						Dinitrotoluene, 2,6-	606-20-2		
						Dinitrotoluene, 2-Amino-4,6-	35572-78-2		
						Dinitrotoluene, 4-Amino-2,6-	19406-51-0		
						Dinitrotoluene, Technical grade	25321-14-6		
						Dinoseb	88-85-7		
5.0E-06	I	3.0E-02	I	V		Dioxane, 1,4-	123-91-1	2.5E+02	1.3E+02
						Dioxins			
1.3E+00	I					~Hexachlorodibenzo-p-dioxin, Mixture	34465-46-8	9.4E-04	
3.8E+01	C	4.0E-08	C	V		~TCDD, 2,3,7,8-	1746-01-6	3.2E-05	1.8E-04
						Diphenamid	957-51-7		
		4.0E-04	X	V		Diphenyl Ether	101-84-8		1.8E+00
						Diphenyl Sulfone	127-63-9		
						Diphenylamine	122-39-4		
2.2E-04	I					Diphenylhydrazine, 1,2-	122-66-7	5.6E+00	
						Diquat	2764-72-9		
2.1E-03	C					Direct Black 38	1937-37-7	5.8E-01	
2.1E-03	C					Direct Blue 6	2602-46-2	5.8E-01	
1.9E-03	C					Direct Brown 95	16071-86-6	6.5E-01	
						Disulfoton	298-04-4		
				V		Dithiane, 1,4-	505-29-3		
						Diuron	330-54-1		
						Dodine	2439-10-3		
				V		EPTC	759-94-4		
				V		Endosulfan	115-29-7		
						Endosulfan Sulfate	1031-07-8		
						Endothall	145-73-3		
						Endrin	72-20-8		
1.2E-06	I	1.0E-03	I	V		Epichlorohydrin	106-89-8	1.0E+03	4.4E+00
		2.0E-02	I	V		Epoxybutane, 1,2-	106-88-7		8.8E+01
						Ethanol, 2-(2-methoxyethoxy)-	111-77-3		
						Ethephon	16672-87-0		
						Ethion	563-12-2		
		6.0E-02	P	V		Ethoxyethanol Acetate, 2-	111-15-9		2.6E+02
		4.0E-02	P	V		Ethoxyethanol, 2-	110-80-5		1.8E+02
		7.0E-02	P	V		Ethyl Acetate	141-78-6		3.1E+02
		8.0E-03	P	V		Ethyl Acrylate	140-88-5		3.5E+01
		4.0E+00	P	V		Ethyl Chloride (Chloroethane)	75-00-3		1.8E+04
				V		Ethyl Ether	60-29-7		
		3.0E-01	P	V		Ethyl Methacrylate	97-63-2		1.3E+03
8.0E-08	I	4.0E+01	I	V		Ethyl Tertiary Butyl Ether (ETBE)	637-92-3	1.5E+04	1.8E+05
						Ethyl-p-nitrophenyl Phosphonate	2104-64-5		
2.5E-06	C	1.0E+00	I	V		Ethylbenzene	100-41-4	4.9E+02	4.4E+03
						Ethylene Cyanohydrin	109-78-4		
				V		Ethylene Diamine	107-15-3		
		4.0E-01	C			Ethylene Glycol	107-21-1		1.8E+03
		1.6E+00	I			Ethylene Glycol Monobutyl Ether	111-76-2		7.0E+03
3.0E-03	I	3.0E-02	C	V	M	Ethylene Oxide	75-21-8	4.1E-01	1.3E+02
1.3E-05	C					Ethylene Thiourea	96-45-7	9.4E+01	
1.9E-02	C			V		Ethyleneimine	151-56-4	6.5E-02	
						Ethylphthalyl Ethyl Glycolate	84-72-0		
						Fenamiphos	22224-92-6		
						Fenpropathrin	39515-41-8		
						Fenvalerate	51630-58-1		
						Fluometuron	2164-17-2		
		1.3E-02	C			Fluoride	16984-48-8		5.7E+01
		1.3E-02	C			Fluorine (Soluble Fluoride)	7782-41-4		5.7E+01
						Fluridone	59756-60-4		
						Flurprimidol	56425-91-3		
						Flusilazole	85509-19-9		
						Flutolanil	66332-96-5		
						Fluvalinate	69409-94-5		
						Folpet	133-07-3		
						Fomesafen	72178-02-0		
						Fonofos	944-22-9		
1.3E-05	I	9.8E-03	A	V		Formaldehyde	50-00-0	9.4E+01	4.3E+01
		3.0E-04	X	V		Formic Acid	64-18-6		1.3E+00
						Fosetyl-AL	39148-24-8		
						Furans			
				V		~Dibenzofuran	132-64-9		
				V		~Furan	110-00-9		
		2.0E+00	I	V		~Tetrahydrofuran	109-99-9		8.8E+03
						Furazolidone	67-45-8		
		5.0E-02	H	V		Furfural	98-01-1		2.2E+02
4.3E-04	C					Furium	531-82-8	2.9E+00	
8.6E-06	C					Furmecyclox	60568-05-0	1.4E+02	
						Glufosinate, Ammonium	77182-82-2		
		8.0E-05	C			Glutaraldehyde	111-30-8		3.5E-01
		1.0E-03	X	V		Glycidaldehyde	765-34-4		4.4E+00
						Glyphosate	1071-83-6		
				V		Guanidine	113-00-8		

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
						Guanidine Chloride	50-01-1		
						Guanidine Nitrate	506-93-4		
						Haloxyfop, Methyl	69806-40-2		
1.3E-03	I			V		Heptachlor	76-44-8	9.4E-01	
2.6E-03	I			V		Heptachlor Epoxide	1024-57-3	4.7E-01	
		3.0E-03	X	V		Heptanal, n-	111-71-7		1.3E+01
		4.0E-01	P	V		Heptane, N-	142-82-5		1.8E+03
				V		Hexabromobenzene	87-82-1		
						Hexabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-153)	68631-49-2		
4.6E-04	I			V		Hexachlorobenzene	118-74-1	2.7E+00	
2.2E-05	I			V		Hexachlorobutadiene	87-68-3	5.6E+01	
1.8E-03	I					Hexachlorocyclohexane, Alpha-	319-84-6	6.8E-01	
5.3E-04	I					Hexachlorocyclohexane, Beta-	319-85-7	2.3E+00	
3.1E-04	C					Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	4.0E+00	
5.1E-04	I					Hexachlorocyclohexane, Technical	608-73-1	2.4E+00	
		2.0E-04	I	V		Hexachlorocyclopentadiene	77-47-4		8.8E-01
1.1E-05	C	3.0E-02	I	V		Hexachloroethane	67-72-1	1.1E+02	1.3E+02
						Hexachlorophene	70-30-4		
						Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4		
		1.0E-05	I	V		Hexamethylene Diisocyanate, 1,6-	822-06-0		4.4E-02
		4.0E-04	C			Hexamethylene diisocyanate biuret	4035-89-6		1.8E+00
		4.0E-04	C			Hexamethylene diisocyanate isocyanurate	3779-63-3		1.8E+00
						Hexamethylphosphoramide	680-31-9		
2.0E-07	X	6.0E-01	P	V		Hexane, Commercial	E5241997	6.1E+03	2.6E+03
		7.0E-01	I	V		Hexane, N-	110-54-3		3.1E+03
						Hexanedioic Acid	124-04-9		
		4.0E-04	P	V		Hexanol, 1-,2-ethyl- (2-Ethyl-1-hexanol)	104-76-7		1.8E+00
		3.0E-02	I	V		Hexanone, 2-	591-78-6		1.3E+02
						Hexazinone	51235-04-2		
						Hexythiazox	78587-05-0		
						Hydramethylnon	67485-29-4		
4.9E-03	I	3.0E-05	P	V		Hydrazine	302-01-2	2.5E-01	1.3E-01
4.9E-03	I					Hydrazine Sulfate	10034-93-2	2.5E-01	
		2.0E-02	I	V		Hydrogen Chloride	7647-01-0		8.8E+01
		1.4E-02	C	V		Hydrogen Fluoride	7664-39-3		6.1E+01
		2.0E-03	I	V		Hydrogen Sulfide	7783-06-4		8.8E+00
						Hydroquinone	123-31-9		
						Imazalil	35554-44-0		
						Imazaquin	81335-37-7		
						Imazethapyr	81335-77-5		
						Iodine	7553-56-2		
						Iprodione	36734-19-7		
						Iron	7439-89-6		
				V		Isobutyl Alcohol	78-83-1		
		2.0E+00	C			Isophorone	78-59-1		8.8E+03
				V		Isopropalin	33820-53-0		
		2.0E-01	P	V		Isopropanol	67-63-0		8.8E+02
						Isopropyl Methyl Phosphonic Acid	1832-54-8		
						Isoxaben	82558-50-7		
		3.0E-01	A	V		JP-7	E1737665		1.3E+03
						Lactofen	77501-63-4		
						Lactonitrile	78-97-7		
						Lanthanum	7439-91-0		
						Lanthanum Acetate Hydrate	100587-90-4		
						Lanthanum Chloride Heptahydrate	10025-84-0		
						Lanthanum Chloride, Anhydrous	10099-58-8		
						Lanthanum Nitrate Hexahydrate	10277-43-7		
						Lead Compounds			
1.2E-05	C					~Lead Phosphate	7446-27-7	1.0E+02	
8.0E-05	C					~Lead acetate	301-04-2	1.5E+01	
						~Lead and Compounds	7439-92-1		
1.1E-05	C					~Lead subacetate	1335-32-6	1.1E+02	
				V		~Tetraethyl Lead	78-00-2		
				V		Lewisite	541-25-3		
						Linuron	330-55-2		
						Lithium	7439-93-2		
						MCPA	94-74-6		
						MCPB	94-81-5		
						MCPP	93-65-2		
						Malathion	121-75-5		
		7.0E-04	C			Maleic Anhydride	108-31-6		3.1E+00
						Maleic Hydrazide	123-33-1		
						Malononitrile	109-77-3		
						Mancozeb	8018-01-7		
						Maneb	12427-38-2		
		5.0E-05	I			Manganese (Diet)	7439-96-5		2.2E-01
		5.0E-05	I			Manganese (Non-diet)	7439-96-5		2.2E-01
						Mephosfolan	950-10-7		
						Mepiquat Chloride	24307-26-4		
						Mercaptobenzothiazole, 2-	149-30-4		
						Mercury Compounds			
		3.0E-04	G			~Mercuric Chloride (and other Mercury salts)	7487-94-7		1.3E+00
		3.0E-04	I	V		~Mercury (elemental)	7439-97-6		1.3E+00
						~Methyl Mercury	22967-92-6		
						~Phenylmercuric Acetate	62-38-4		
				V		Merphos	150-50-5		
						Metalaxyl	57837-19-1		
		3.0E-02	P	V		Methacrylonitrile	126-98-7		1.3E+02
						Methamidophos	10265-92-6		
		2.0E+01	I	V		Methanol	67-56-1		8.8E+04

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
						Methidathion	950-37-8		
						Methomyl	16752-77-5		
						Methoxy-5-nitroaniline, 2-	99-59-2		
						Methoxychlor	72-43-5		
		1.0E-03	P	V		Methoxyethanol Acetate, 2-	110-49-6		4.4E+00
		7.0E-03	P	V		Methoxyethanol, 2-	109-86-4		3.1E+01
				V		Methyl Acetate	79-20-9		
		2.0E-02	P	V		Methyl Acrylate	96-33-3		8.8E+01
		5.0E+00	I	V		Methyl Ethyl Ketone (2-Butanone)	78-93-3		2.2E+04
1.0E-03	X	2.0E-05	X	V		Methyl Hydrazine	60-34-4	1.2E+00	8.8E-02
		3.0E+00	I	V		Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1		1.3E+04
		1.0E-03	C	V		Methyl Isocyanate	624-83-9		4.4E+00
		7.0E-01	I	V		Methyl Methacrylate	80-62-6		3.1E+03
						Methyl Parathion	298-00-0		
						Methyl Phosphonic Acid	993-13-5		
		4.0E-02	H	V		Methyl Styrene (Mixed Isomers)	25013-15-4		1.8E+02
2.8E-05	C					Methyl methanesulfonate	66-27-3	4.4E+01	
2.6E-07	C	3.0E+00	I	V		Methyl tert-Butyl Ether (MTBE)	1634-04-4	4.7E+03	1.3E+04
						Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2		
		3.0E+00	X	V		Methyl-2-Pentanol, 4-	108-11-2		1.3E+04
						Methyl-5-Nitroaniline, 2-	99-55-8		
2.4E-03	C					Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	5.1E-01	
3.7E-05	C					Methylaniline Hydrochloride, 2-	636-21-5	3.3E+01	
						Methylarsonic acid	124-58-3		
						Methylbenzene,1-4-diamine monohydrochloride, 2-	74612-12-7		
						Methylbenzene-1,4-diamine sulfate, 2-	615-50-9		
6.3E-03	C				M	Methylcholanthrene, 3-	56-49-5	1.9E-01	
1.0E-08	I	6.0E-01	I	V	M	Methylene Chloride	75-09-2	1.2E+05	2.6E+03
4.3E-04	C				M	Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	2.9E+00	
1.3E-05	C					Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	9.4E+01	
4.6E-04	C	2.0E-02	C			Methylenebisbenzenamine, 4,4'-	101-77-9	2.7E+00	8.8E+01
		6.0E-04	I			Methylenediphenyl Diisocyanate	101-68-8		2.6E+00
				V		Methylstyrene, Alpha-	98-83-9		
						Metolachlor	51218-45-2		
						Metribuzin	21087-64-9		
						Metsulfuron-methyl	74223-64-6		
4.5E-06	X	1.0E-01	P	V		Midrange Aliphatic Hydrocarbon Streams	E1790669	2.7E+02	4.4E+02
				V		Mineral oils	8012-95-1		
5.1E-03	C			V		Mirex	2385-85-5	2.4E-01	
						Molinate	2212-67-1		
		2.0E-03	A			Molybdenum	7439-98-7		8.8E+00
						Monochloramine	10599-90-3		
						Monomethylaniline	100-61-8		
						Myclobutanil	88671-89-0		
						N,N'-Diphenyl-1,4-benzenediamine	74-31-7		
				V		Naled	300-76-5		
		1.0E-01	P	V		Naphtha, High Flash Aromatic (HFAN)	64742-95-6		4.4E+02
0.0E+00	C					Naphthylamine, 2-	91-59-8		
						Napropamide	15299-99-7		
2.6E-04	C	1.4E-05	C			Nickel Acetate	373-02-4	4.7E+00	6.1E-02
2.6E-04	C	1.4E-05	C			Nickel Carbonate	3333-67-3	4.7E+00	6.1E-02
2.6E-04	C	1.4E-05	C	V		Nickel Carbonyl	13463-39-3	4.7E+00	6.1E-02
2.6E-04	C	1.4E-05	C			Nickel Hydroxide	12054-48-7	4.7E+00	6.1E-02
2.6E-04	C	2.0E-05	C			Nickel Oxide	1313-99-1	4.7E+00	8.8E-02
2.4E-04	I	1.4E-05	C			Nickel Refinery Dust	E715532	5.1E+00	6.1E-02
2.6E-04	C	9.0E-05	A			Nickel Soluble Salts	7440-02-0	4.7E+00	3.9E-01
4.8E-04	I	1.4E-05	C			Nickel Subsulfide	12035-72-2	2.6E+00	6.1E-02
2.6E-04	C	1.4E-05	C			Nickelocene	1271-28-9	4.7E+00	6.1E-02
						Nitrate (measured as nitrogen)	14797-55-8		
						Nitrate + Nitrite (measured as nitrogen)	E701177		
						Nitrite (measured as nitrogen)	14797-65-0		
		5.0E-05	X			Nitroaniline, 2-	88-74-4		2.2E-01
		6.0E-03	P			Nitroaniline, 4-	100-01-6		2.6E+01
4.0E-05	I	9.0E-03	I	V		Nitrobenzene	98-95-3	3.1E+01	3.9E+01
						Nitrocellulose	9004-70-0		
						Nitrofurantoin	67-20-9		
3.7E-04	C					Nitrofurazone	59-87-0	3.3E+00	
						Nitroglycerin	55-63-0		
						Nitroguanidine	556-88-7		
8.8E-06	P	5.0E-03	P	V		Nitromethane	75-52-5	1.4E+02	2.2E+01
5.8E-04	X	2.0E-02	I	V		Nitropropane, 2-	79-46-9	2.1E+00	8.8E+01
7.7E-03	C				M	Nitroso-N-ethylurea, N-	759-73-9	1.6E-01	
3.4E-02	C				M	Nitroso-N-methylurea, N-	684-93-5	3.6E-02	
1.6E-03	I			V		Nitroso-di-N-butylamine, N-	924-16-3	7.7E-01	
2.0E-03	C					Nitroso-di-N-propylamine, N-	621-64-7	6.1E-01	
8.0E-04	C					Nitrosodiethanolamine, N-	1116-54-7	1.5E+00	
4.3E-02	I				M	Nitrosodiethylamine, N-	55-18-5	2.9E-02	
1.4E-02	I	4.0E-05	X	V	M	Nitrosodimethylamine, N-	62-75-9	8.8E-02	1.8E-01
2.6E-06	C					Nitrosodiphenylamine, N-	86-30-6	4.7E+02	
6.3E-03	C			V		Nitrosomethylethylamine, N-	10595-95-6	1.9E-01	
1.9E-03	C					Nitrosomorpholine [N-]	59-89-2	6.5E-01	
2.7E-03	C					Nitrosopiperidine [N-]	100-75-4	4.5E-01	
6.1E-04	I					Nitrosopyrrolidine, N-	930-55-2	2.0E+00	
						Nitrotoluene, m-	99-08-1		
				V		Nitrotoluene, o-	88-72-2		
						Nitrotoluene, p-	99-99-0		
		2.0E-02	P	V		Nonane, n-	111-84-2		8.8E+01
						Norflurazon	27314-13-2		
						Octabromodiphenyl Ether	32536-52-0		
						Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0		

RMLS TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
						Octamethylpyrophosphoramide	152-16-9		
						Oryzalin	19044-88-3		
						Oxadiazon	19666-30-9		
						Oxamyl	23135-22-0		
						Oxyfluorfen	42874-03-3		
						Paclobutrazol	76738-62-0		
						Paraquat Dichloride	1910-42-5		
						Parathion	56-38-2		
					V	Pebulate	1114-71-2		
						Pendimethalin	40487-42-1		
					V	Pentabromodiphenyl Ether	32534-81-9		
						Pentabromodiphenyl ether, 2,2',4,4',5- (BDE-99)	60348-60-9		
					V	Pentachlorobenzene	608-93-5		
					V	Pentachloroethane	76-01-7		
					V	Pentachloronitrobenzene	82-68-8		
5.1E-06	C					Pentachlorophenol	87-86-5	2.4E+02	
						Pentaerythritol tetranitrate (PETN)	78-11-5		
						Pentamethylphosphoramide (PMPA)	10159-46-3		
		1.0E+00	P	V		Pentane, n-	109-66-0		4.4E+03
						Per- and Polyfluoroalkyl Substances (PFAS)			
						~Ammonium perfluoro-2-methyl-3-oxahexanoate	62037-80-3		
					V	~Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6		
						~Perfluorobutanesulfonate	45187-15-3		
						~Perfluorobutanesulfonic acid (PFBS)	375-73-5		
						~Perfluorohexanesulfonate	108427-53-8		
						~Perfluorohexanesulfonic acid (PFHxS)	355-46-4		
						~Perfluorononanoate	72007-68-2		
						~Perfluorononanoic acid (PFNA)	375-95-1		
						~Perfluorooctanesulfonate	45298-90-6		
						~Perfluorooctanesulfonic acid (PFOS)	1763-23-1		
						~Perfluorooctanoate	45285-51-6		
						~Perfluorooctanoic acid (PFOA)	335-67-1		
						~Potassium perfluorobutanesulfonate	29420-49-3		
						~Potassium perfluorooctanesulfonate	2795-39-3		
						Perchlorates			
						~Ammonium Perchlorate	7790-98-9		
						~Lithium Perchlorate	7791-03-9		
						~Perchlorate and Perchlorate Salts	14797-73-0		
						~Potassium Perchlorate	7778-74-7		
						~Sodium Perchlorate	7601-89-0		
						Permethrin	52645-53-1		
6.3E-07	C					Phenacetin	62-44-2	1.9E+03	
						Phenmedipham	13684-63-4		
		2.0E-01	C			Phenol	108-95-2		8.8E+02
						Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1		
						Phenothiazine	92-84-2		
					V	Phenyl Isothiocyanate	103-72-0		
						Phenylenediamine, m-	108-45-2		
						Phenylenediamine, o-	95-54-5		
						Phenylenediamine, p-	106-50-3		
						Phenylphenol, 2-	90-43-7		
						Phorate	298-02-2		
		3.0E-04	I	V		Phosgene	75-44-5		1.3E+00
						Phosmet	732-11-6		
		3.0E-04	I	V		Phosphine	7803-51-2		1.3E+00
		1.0E-02	I			Phosphoric Acid	7664-38-2		4.4E+01
					V	Phosphorus, White	7723-14-0		
						Phthalates			
2.4E-06	C					~Bis(2-ethylhexyl)phthalate	117-81-7	5.1E+02	
						~Butyl Benzyl Phthalate	85-68-7		
						~Butylphthalyl Butylglycolate	85-70-1		
						~Dibutyl Phthalate	84-74-2		
						~Diethyl Phthalate	84-66-2		
					V	~Dimethylterephthalate	120-61-6		
						~Octyl Phthalate, di-N-	117-84-0		
						~Phthalic Acid, p-	100-21-0		
		2.0E-02	C			~Phthalic Anhydride	85-44-9		8.8E+01
						Picloram	1918-02-1		
						Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3		
						Picric Acid (2,4,6-Trinitrophenol)	88-89-1		
						Pirimiphos, Methyl	29232-93-7		
8.6E-03	C					Polybrominated Biphenyls	36355-01-8	1.4E-01	
						Polychlorinated Biphenyls (PCBs)			
2.0E-05	G				V	~Aroclor 1016	12674-11-2	6.1E+01	
5.7E-04	G				V	~Aroclor 1221	11104-28-2	2.1E+00	
5.7E-04	G				V	~Aroclor 1232	11141-16-5	2.1E+00	
5.7E-04	G				V	~Aroclor 1242	53469-21-9	2.1E+00	
5.7E-04	G				V	~Aroclor 1248	12672-29-6	2.1E+00	
5.7E-04	G				V	~Aroclor 1254	11097-69-1	2.1E+00	
5.7E-04	G				V	~Aroclor 1260	11096-82-5	2.1E+00	
					V	~Aroclor 5460	11126-42-4		
1.1E-03	W	1.3E-03	W	V		~Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	1.1E+00	5.8E+00
1.1E-03	W	1.3E-03	W	V		~Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167)	52663-72-6	1.1E+00	5.8E+00
1.1E-03	W	1.3E-03	W	V		~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	1.1E+00	5.8E+00
1.1E-03	W	1.3E-03	W	V		~Hexachlorobiphenyl, 2,3,3',4,4',5- (PCB 156)	38380-08-4	1.1E+00	5.8E+00
1.1E+00	W	1.3E-06	W	V		~Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	1.1E-03	5.8E-03
1.1E-03	W	1.3E-03	W	V		~Pentachlorobiphenyl, 2',3,4,4',5- (PCB 123)	65510-44-3	1.1E+00	5.8E+00
1.1E-03	W	1.3E-03	W	V		~Pentachlorobiphenyl, 2,3',4,4',5- (PCB 118)	31508-00-6	1.1E+00	5.8E+00
1.1E-03	W	1.3E-03	W	V		~Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	1.1E+00	5.8E+00
1.1E-03	W	1.3E-03	W	V		~Pentachlorobiphenyl, 2,3,4,4',5- (PCB 114)	74472-37-0	1.1E+00	5.8E+00

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
3.8E+00	W	4.0E-07	W	V		~Pentachlorobiphenyl, 3,3',4,4',5- (PCB 126)	57465-28-8	3.2E-04	1.8E-03
5.7E-04	I			V		~Polychlorinated Biphenyls (high risk)	1336-36-3	2.1E+00	
1.0E-04	I			V		~Polychlorinated Biphenyls (low risk)	1336-36-3	1.2E+01	
2.0E-05	I			V		~Polychlorinated Biphenyls (lowest risk)	1336-36-3	6.1E+01	
3.8E-03	W	4.0E-04	W			~Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	3.2E-01	1.8E+00
1.1E-02	W	1.3E-04	W	V		~Tetrachlorobiphenyl, 3,4,4',5- (PCB 81)	70362-50-4	1.1E-01	5.8E-01
		6.0E-04	I			Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9		2.6E+00
						Polynuclear Aromatic Hydrocarbons (PAHs)			
				V		~Acenaphthene	83-32-9		
				V		~Anthracene	120-12-7		
6.0E-05	E			V	M	~Benz[a]anthracene	56-55-3	2.0E+01	
		2.0E-06	X			~Benzo(e)pyrene	192-97-2		8.8E-03
1.1E-04	C					~Benzo(j)fluoranthene	205-82-3	1.1E+01	
6.0E-04	I	2.0E-06	I		M	~Benzo[a]pyrene	50-32-8	2.0E+00	8.8E-03
6.0E-05	E				M	~Benzo[b]fluoranthene	205-99-2	2.0E+01	
6.0E-06	E				M	~Benzo[k]fluoranthene	207-08-9	2.0E+02	
				V		~Chloronaphthalene, Beta-	91-58-7		
6.0E-07	E				M	~Chrysene	218-01-9	2.0E+03	
6.0E-04	E				M	~Dibenz[a,h]anthracene	53-70-3	2.0E+00	
1.1E-03	C					~Dibenzo(a,e)pyrene	192-65-4	1.1E+00	
7.1E-02	C				M	~Dimethylbenz(a)anthracene, 7,12-	57-97-6	1.7E-02	
						~Fluoranthene	206-44-0		
				V		~Fluorene	86-73-7		
6.0E-05	E				M	~Indeno[1,2,3-cd]pyrene	193-39-5	2.0E+01	
				V		~Methylnaphthalene, 1-	90-12-0		
				V		~Methylnaphthalene, 2-	91-57-6		
3.4E-05	C	3.0E-03	I	V		~Naphthalene	91-20-3	3.6E+01	1.3E+01
1.1E-04	C					~Nitropyrene, 4-	57835-92-4	1.1E+01	
				V		~Pyrene	129-00-0		
						Prochloraz	67747-09-5		
				V		Profluralin	26399-36-0		
						Prometon	1610-18-0		
						Prometryn	7287-19-6		
						Pronamide	23950-58-5		
						Propachlor	1918-16-7		
						Propanil	709-98-8		
						Propargite	2312-35-8		
				V		Propargyl Alcohol	107-19-7		
						Propazine	139-40-2		
						Propham	122-42-9		
						Propiconazole	60207-90-1		
		8.0E-03	I	V		Propionaldehyde	123-38-6		3.5E+01
		1.0E+00	X	V		Propyl benzene	103-65-1		4.4E+03
		3.0E+00	C	V		Propylene	115-07-1		1.3E+04
						Propylene Glycol	57-55-6		
		2.7E-04	A			Propylene Glycol Dinitrate	6423-43-4		1.2E+00
		2.0E+00	I	V		Propylene Glycol Monomethyl Ether	107-98-2		8.8E+03
3.7E-06	I	3.0E-02	I	V		Propylene Oxide	75-56-9	3.3E+02	1.3E+02
				V		Pyridine	110-86-1		
						Quinalphos	13593-03-8		
						Quinoline	91-22-5		
						Quizalofop-ethyl	76578-14-8		
		3.0E+04	A			Refractory Ceramic Fibers (units in fibers)	E715557		1.3E+05
						Resmethrin	10453-86-8		
				V		Ronnel	299-84-3		
						Rotenone	83-79-4		
6.3E-05	C				M	Safrole	94-59-7	1.9E+01	
						Selenious Acid	7783-00-8		
		2.0E-02	C			Selenium	7782-49-2		8.8E+01
		2.0E-02	C			Selenium Sulfide	7446-34-6		8.8E+01
						Sethoxydim	74051-80-2		
		3.0E-03	C			Silica (crystalline, respirable)	7631-86-9		1.3E+01
						Silver	7440-22-4		
						Simazine	122-34-9		
						Sodium Acifluorfen	62476-59-9		
						Sodium Azide	26628-22-8		
						Sodium Diethyldithiocarbamate	148-18-5		
		1.3E-02	C			Sodium Fluoride	7681-49-4		5.7E+01
						Sodium Fluoroacetate	62-74-8		
						Sodium Metavanadate	13718-26-8		
						Sodium Tungstate	13472-45-2		
						Sodium Tungstate Dihydrate	10213-10-2		
						Stirofos (Tetrachlorovinphos)	961-11-5		
						Strontium, Stable	7440-24-6		
						Strychnine	57-24-9		
		1.0E+00	I	V		Styrene	100-42-5		4.4E+03
						Styrene-Acrylonitrile (SAN) Trimer (THNA isomer)	57964-39-3		
						Styrene-Acrylonitrile (SAN) Trimer (THNP isomer)	57964-40-6		
		2.0E-03	X			Sulfolane	126-33-0		8.8E+00
						Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9		
		1.0E-03	C	V		Sulfur Trioxide	7446-11-9		4.4E+00
		1.0E-03	C			Sulfuric Acid	7664-93-9		4.4E+00
7.1E-06	I					Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl ester	140-57-8	1.7E+02	
						TCMTB	21564-17-0		
						Tebuthiuron	34014-18-1		
						Temephos	3383-96-8		
						Terbacil	5902-51-2		
				V		Terbufos	13071-79-9		
						Terbutryn	886-50-0		
1.3E-06	C			V		Tert-Butyl Acetate	540-88-5	9.4E+02	

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
						Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1		
				V		Tetrachlorobenzene, 1,2,4,5-	95-94-3		
7.4E-06	I			V		Tetrachloroethane, 1,1,1,2-	630-20-6	1.7E+02	
5.8E-05	C			V		Tetrachloroethane, 1,1,2,2-	79-34-5	2.1E+01	
2.6E-07	I	4.0E-02	I	V		Tetrachloroethylene	127-18-4	4.7E+03	1.8E+02
						Tetrachlorophenol, 2,3,4,6-	58-90-2		
				V		Tetrachlorotoluene, p- alpha, alpha, alpha-	5216-25-1		
						Tetraethyl Dithiopyrophosphate	3689-24-5		
		8.0E+01	I	V		Tetrafluoroethane, 1,1,1,2-	811-97-2		3.5E+05
						Tetramethylphosphoramide, -N,N,N',N" (TMPA)	16853-36-4		
						Tetryl (Trinitrophenylmethylnitramine)	479-45-8		
						Thallic Oxide	1314-32-5		
						Thallium (I) Nitrate	10102-45-1		
						Thallium (Soluble Salts)	7440-28-0		
				V		Thallium Acetate	563-68-8		
						Thallium Carbonate	6533-73-9		
						Thallium Chloride	7791-12-0		
						Thallium Selenite	12039-52-0		
						Thallium Sulfate	7446-18-6		
						Thifensulfuron-methyl	79277-27-3		
						Thiobencarb	28249-77-6		
						Thiodiglycol	111-48-8		
						Thiofanox	39196-18-4		
						Thiophanate, Methyl	23564-05-8		
						Thiram	137-26-8		
						Tin	7440-31-5		
		1.0E-04	A	V		Titanium Tetrachloride	7550-45-0		4.4E-01
		5.0E+00	I	V		Toluene	108-88-3		2.2E+04
1.1E-05	C	8.0E-06	C	V		Toluene-2,4-diisocyanate	584-84-9	1.1E+02	3.5E-02
						Toluene-2,5-diamine	95-70-5		
1.1E-05	C	8.0E-06	C	V		Toluene-2,6-diisocyanate	91-08-7	1.1E+02	3.5E-02
						Toluenediamine, 2,3-	2687-25-4		
						Toluenediamine, 3,4-	496-72-0		
						Toluic Acid, p-	99-94-5		
5.1E-05	C					Toluidine, o- (Methylaniline, 2-)	95-53-4	2.4E+01	
						Toluidine, p-	106-49-0		
				V		Total Petroleum Hydrocarbons (Aliphatic High)	E1790670		
		6.0E-01	P	V		Total Petroleum Hydrocarbons (Aliphatic Low)	E1790666		2.6E+03
		1.0E-01	P	V		Total Petroleum Hydrocarbons (Aliphatic Medium)	E1790668		4.4E+02
						Total Petroleum Hydrocarbons (Aromatic High)	E1790676		
		3.0E-02	P	V		Total Petroleum Hydrocarbons (Aromatic Low)	E1790672		1.3E+02
		3.0E-03	P	V		Total Petroleum Hydrocarbons (Aromatic Medium)	E1790674		1.3E+01
3.2E-04	I					Toxaphene	8001-35-2	3.8E+00	
						Toxaphene, Weathered	E1841606		
						Tralomethrin	66841-25-6		
				V		Tri-n-butyltin	688-73-3		
						Triacetin	102-76-1		
						Triadimefon	43121-43-3		
				V		Triallate	2303-17-5		
						Triasulfuron	82097-50-5		
						Tribenuron-methyl	101200-48-0		
				V		Tribromobenzene, 1,2,4-	615-54-3		
						Tribromophenol, 2,4,6-	118-79-6		
						Tribufos	78-48-8		
						Tributyl Phosphate	126-73-8		
						Tributyltin Compounds	E1790679		
						Tributyltin Oxide	56-35-9		
						Trichloramine	10025-85-1		
		5.0E+00	P	V		Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1		2.2E+04
						Trichloroacetic Acid	76-03-9		
						Trichloroaniline HCl, 2,4,6-	33663-50-2		
						Trichloroaniline, 2,4,6-	634-93-5		
				V		Trichlorobenzene, 1,2,3-	87-61-6		
		2.0E-03	P	V		Trichlorobenzene, 1,2,4-	120-82-1		8.8E+00
		5.0E+00	I	V		Trichloroethane, 1,1,1-	71-55-6		2.2E+04
1.6E-05	I	2.0E-04	X	V		Trichloroethane, 1,1,2-	79-00-5	7.7E+01	8.8E-01
4.1E-06	I	2.0E-03	I	V	M	Trichloroethylene	79-01-6	3.0E+02	8.8E+00
				V		Trichlorofluoromethane	75-69-4		
						Trichlorophenol, 2,4,5-	95-95-4		
3.1E-06	I					Trichlorophenol, 2,4,6-	88-06-2	4.0E+02	
						Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5		
						Trichlorophenoxypropionic acid, -2,4,5	93-72-1		
				V		Trichloropropane, 1,1,2-	598-77-6		
		3.0E-04	I	V	M	Trichloropropane, 1,2,3-	96-18-4		1.3E+00
		3.0E-04	P	V		Trichloropropene, 1,2,3-	96-19-5		1.3E+00
						Tricresyl Phosphate (TCP)	1330-78-5		
						Tridiphane	58138-08-2		
		7.0E-03	I	V		Triethylamine	121-44-8		3.1E+01
						Triethylene Glycol	112-27-6		
		2.0E+01	P	V		Trifluoroethane, 1,1,1-	420-46-2		8.8E+04
				V		Trifluralin	1582-09-8		
						Trimethyl Phosphate	512-56-1		
		6.0E-02	I	V		Trimethylbenzene, 1,2,3-	526-73-8		2.6E+02
		6.0E-02	I	V		Trimethylbenzene, 1,2,4-	95-63-6		2.6E+02
		6.0E-02	I	V		Trimethylbenzene, 1,3,5-	108-67-8		2.6E+02
				V		Trimethylpentene, 2,4,4-	25167-70-8		
						Trinitrobenzene, 1,3,5-	99-35-4		
						Trinitrotoluene, 2,4,6-	118-96-7		
						Triphenylphosphine Oxide	791-28-6		
						Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8		

RMLs TR=1E-04 THQ=1.0

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; ca = cancer; nc = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; max = ceiling limit exceeded; sat = Csat exceeded.

Toxicity and Chemical-specific Information						Contaminant		Carcinogenic Target Risk (TR) = 1E-04	Noncancer Hazard Index (HI) = 1
IUR $(ug/m^3)^{-1}$	key	RfC_i (mg/m^3)	key	vol	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1E-04 (ug/m^3)	Noncarcinogenic SL THI=1 (ug or fibers/m^3)
						Tris(1-chloro-2-propyl)phosphate	13674-84-5		
6.6E-04	C			V		Tris(2,3-dibromopropyl)phosphate	126-72-7	1.9E+00	
						Tris(2-chloroethyl)phosphate	115-96-8		
						Tris(2-ethylhexyl)phosphate	78-42-2		
						Tungsten	7440-33-7		
		4.0E-05	A			Uranium	7440-61-1		1.8E-01
2.9E-04	C				M	Urethane	51-79-6	4.2E+00	
8.3E-03	P	7.0E-06	P			Vanadium Pentoxide	1314-62-1	1.5E-01	3.1E-02
		1.0E-04	A			Vanadium and Compounds	7440-62-2		4.4E-01
				V		Vernolate	1929-77-7		
						Vinclozolin	50471-44-8		
		2.0E-01	I	V		Vinyl Acetate	108-05-4		8.8E+02
1.5E-05	P	3.0E-03	I	V		Vinyl Bromide	593-60-2	8.2E+01	1.3E+01
4.4E-06	I	8.0E-02	A	V	M	Vinyl Chloride	75-01-4	2.8E+02	3.5E+02
						Warfarin	81-81-2		
		1.0E-01	G	V		Xylene, m-	108-38-3		4.4E+02
		1.0E-01	G	V		Xylene, o-	95-47-6		4.4E+02
		1.0E-01	G	V		Xylene, p-	106-42-3		4.4E+02
		1.0E-01	I	V		Xylenes	1330-20-7		4.4E+02
						Zinc Phosphide	1314-84-7		
						Zinc and Compounds	7440-66-6		
						Zineb	12122-67-7		
						Zirconium	7440-67-7		

RMLs
TR=1E-04
THQ=1.0

Contaminant		Molecular Weight		Volatility Parameters					Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients					Water Solubility		Tapwater Dermal Parameters				
Analyte	CAS No.	MW	MW Ref	H^+ (unitless)	HLC (atm-m^3/mole)	H^+ and HLC Ref	VP (mmHg)	VP Ref	MP (C)	MP Ref	Density (g/cm^3)	Density Ref	D_{oa} (cm^2/s)	D_{iw} (cm^2/s)	D_{oa} and D_{iw} Ref	K_d (L/kg)	K_d Ref	K_{oc} (L/kg)	K_{oc} Ref	log K_{ow} (unitless)	log K_{ow} Ref	S (mg/L)	S Ref	B (unitless)	t_{event} (hr/event)	t^* (hr)	K_p (cm/hr)	K_p Ref
Acephate	30560-19-1	1.8E+02	PHYSPROP	2.0E-11	5.0E-13	EPI	1.7E-06	PHYSPROP	8.8E+01	PHYSPROP	1.4E+00	CRC	3.7E-02	8.0E-06	WATER9 (U.S. EPA, 2001)			1.0E+01	EPI	-8.5E-01	PHYSPROP	8.2E+05	PHYSPROP	2.1E-04	1.1E+00	2.7E+00	4.0E-05	EPI
Acetaldehyde	75-07-0	4.4E+01	PHYSPROP	2.7E-03	6.7E-05	PHYSPROP	9.0E+02	PHYSPROP	-1.2E+02	PHYSPROP	7.8E-01	CRC	1.3E-01	1.4E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-3.4E-01	PHYSPROP	1.0E+06	PHYSPROP	1.3E-03	1.9E-01	4.5E-01	5.3E-04	EPI
Acetochlor	34256-82-1	2.7E+02	PHYSPROP	9.1E-07	2.2E-08	PHYSPROP	2.8E-05	PHYSPROP	0.0E+00	EPI	1.1E+00	PubChem	2.2E-02	5.6E-06	WATER9 (U.S. EPA, 2001)			3.0E+02	EPI	3.0E+00	PHYSPROP	2.2E+02	PHYSPROP	3.1E-02	3.4E+00	8.2E+00	5.0E-03	EPI
Acetone	67-64-1	5.8E+01	PHYSPROP	1.4E-03	3.5E-05	PHYSPROP	2.3E+02	PHYSPROP	-9.5E+01	PHYSPROP	7.8E-01	CRC	1.1E-01	1.1E-05	WATER9 (U.S. EPA, 2001)			2.4E+00	EPI	-2.4E-01	PHYSPROP	1.0E+06	PHYSPROP	1.5E-03	2.2E-01	5.3E-01	5.1E-04	EPI
Acetone Cyanohydrin	75-86-5	8.5E+01	PHYSPROP	8.1E-08	2.0E-09	PHYSPROP	3.4E-01	PHYSPROP	-1.9E+01	PHYSPROP	9.3E-01	CRC	8.6E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-3.0E-02	PHYSPROP	1.0E+06	PHYSPROP	1.8E-03	3.2E-01	7.6E-01	5.0E-04	EPI
Acetonitrile	75-05-8	4.1E+01	PHYSPROP	1.4E-03	3.5E-05	PHYSPROP	8.9E+01	PHYSPROP	-4.4E+01	PHYSPROP	7.9E-01	CRC	1.3E-01	1.4E-05	WATER9 (U.S. EPA, 2001)			4.7E+00	EPI	-3.4E-01	PHYSPROP	1.0E+06	PHYSPROP	1.4E-03	1.8E-01	4.3E-01	5.5E-04	EPI
Acetophenone	98-86-2	1.2E+02	PHYSPROP	4.3E-04	1.0E-05	PHYSPROP	4.0E-01	PHYSPROP	2.0E+01	PHYSPROP	1.0E+00	CRC	6.5E-02	8.7E-06	WATER9 (U.S. EPA, 2001)			1.6E+02	EPI	1.6E+00	PHYSPROP	6.1E+03	PHYSPROP	1.6E-02	5.0E-01	1.2E+00	3.7E-03	EPI
Acetylaminofluorene, 2-	53-96-3	1.2E+02	PHYSPROP	7.8E-09	1.9E-10	PHYSPROP	9.4E-08	PHYSPROP	1.9E+02	PHYSPROP			5.2E-02	8.0E-06	WATER9 (U.S. EPA, 2001)			2.2E+03	EPI	3.1E+00	PHYSPROP	5.5E+00	PHYSPROP	6.0E-02	1.9E+00	4.5E+00	1.0E-02	EPI
Acrolein	107-02-8	5.6E+01	PHYSPROP	5.0E-03	1.2E-04	PHYSPROP	2.7E+02	PHYSPROP	-8.8E+01	PHYSPROP	8.4E-01	CRC	1.1E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-1.0E-02	PHYSPROP	2.1E+05	PHYSPROP	2.2E-03	2.2E-01	5.2E-01	7.5E-04	EPI
Acrylamide	79-06-1	7.1E+01	PHYSPROP	7.0E-08	1.7E-09	EPI	7.0E-03	PHYSPROP	8.5E+01	PHYSPROP	1.2E+00	LANGE	1.1E-01	1.3E-05	WATER9 (U.S. EPA, 2001)			5.7E+00	EPI	-6.7E-01	PHYSPROP	3.9E+05	PHYSPROP	7.3E-04	2.6E-01	6.3E-01	2.2E-04	EPI
Acrylic Acid	79-10-7	7.2E+01	PHYSPROP	1.5E-05	3.7E-07	EPI	4.0E+00	PHYSPROP	1.3E+01	PHYSPROP	1.1E+00	CRC	1.0E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			1.4E+00	EPI	3.5E-01	PHYSPROP	1.0E+06	PHYSPROP	3.4E-03	2.7E-01	6.4E-01	1.1E-03	EPI
Acrylonitrile	107-13-1	5.3E+01	PHYSPROP	5.6E-03	1.4E-04	PHYSPROP	1.1E+02	PHYSPROP	-8.4E+01	PHYSPROP	8.0E-01	CRC	1.1E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			8.5E+00	EPI	2.5E-01	PHYSPROP	7.5E+04	PHYSPROP	3.3E-03	2.1E-01	5.0E-01	1.2E-03	EPI
Adiponitrile	111-69-3	1.1E+02	PHYSPROP	4.9E-08	1.2E-09	EPI	6.8E-04	PHYSPROP	1.0E+00	PHYSPROP	9.7E-01	CRC	7.1E-02	9.0E-06	WATER9 (U.S. EPA, 2001)			2.0E+01	EPI	-3.2E-01	PHYSPROP	8.0E+04	PHYSPROP	9.5E-04	4.2E-01	1.0E+00	2.4E-04	EPI
Alachlor	15972-60-8	2.7E+02	PHYSPROP	3.4E-07	8.3E-09	PHYSPROP	2.2E-05	PHYSPROP	4.0E+01	PHYSPROP	1.1E+00	CRC	2.3E-02	5.7E-06	WATER9 (U.S. EPA, 2001)			3.1E+02	EPI	3.5E+00	PHYSPROP	2.4E+02	PHYSPROP	6.6E-02	3.4E+00	8.2E+00	1.1E-02	EPI
Aldicarb	116-06-3	1.9E+02	PHYSPROP	5.9E-08	1.4E-09	EPI	3.5E-05	PHYSPROP	9.9E+01	PHYSPROP	1.2E+00	CRC	3.2E-02	7.2E-06	WATER9 (U.S. EPA, 2001)			2.5E+01	EPI	1.1E+00	PHYSPROP	6.0E+03	PHYSPROP	4.0E-03	1.2E+00	2.9E+00	7.6E-04	EPI
Aldicarb Sulfone	1646-88-4	2.2E+02	PHYSPROP	1.4E-07	3.4E-09	EPI	9.0E-05	PHYSPROP	1.4E+02	PHYSPROP			5.2E-02	6.1E-06	WATER9 (U.S. EPA, 2001)			1.0E+01	EPI	-5.7E-01	PHYSPROP	1.0E+04	PHYSPROP	2.1E-04	1.8E+00	4.4E+00	3.7E-05	EPI
Aldicarb sulfoxide	1646-87-3	2.1E+02	PHYSPROP	4.0E-08	9.7E-10	EPI	1.0E-04	PHYSPROP	7.8E+01	EPI			5.4E-02	6.4E-06	WATER9 (U.S. EPA, 2001)			1.0E+01	EPI	-7.8E-01	PHYSPROP	2.8E+04	PHYSPROP	1.8E-04	1.5E+00	3.6E+00	3.3E-05	EPI
Aldrin	309-00-2	3.6E+02	PHYSPROP	1.8E-03	4.4E-05	PHYSPROP	1.2E-04	PHYSPROP	1.0E+02	PHYSPROP	1.6E+00	PubChem	2.3E-02	5.8E-06	WATER9 (U.S. EPA, 2001)			8.2E+04	EPI	6.5E+00	PHYSPROP	1.7E-02	PHYSPROP	2.2E+00	1.2E+01	4.8E+01	2.9E-01	EPI
Allyl Alcohol	107-18-6	5.8E+01	PHYSPROP	2.0E-04	5.0E-06	PHYSPROP	2.6E+01	PHYSPROP	-1.3E+02	PHYSPROP	8.5E-01	CRC	1.1E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			1.9E+00	EPI	1.7E-01	PHYSPROP	1.0E+06	PHYSPROP	2.8E-03	2.2E-01	5.3E-01	9.6E-04	EPI
Allyl Chloride	107-05-1	7.7E+01	PHYSPROP	4.5E-01	1.1E-02	EPI	3.7E+02	PHYSPROP	-1.3E+02	PHYSPROP	9.4E-01	CRC	9.4E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			4.0E+01	EPI	1.9E+00	PHYSPROP	3.4E+03	PHYSPROP	3.8E-02	2.8E-01	6.8E-01	1.1E-02	EPI
Aluminum	7429-90-5	2.7E+01	CRC				0.0E+00	NIOSH	6.6E+02	CRC	2.7E+00	CRC				1.5E+03	BAES							2.0E-03	1.5E-01	3.6E-01	1.0E-03	RAGSE
Aluminum Phosphide	20859-73-8	5.8E+01	PHYSPROP						2.6E+03	CRC	2.4E+00	CRC												2.9E-03	2.2E-01	5.3E-01	1.0E-03	RAGSE
Ametryn	834-12-8	2.3E+02	PHYSPROP	9.9E-08	2.4E-09	EPI	2.7E-06	PHYSPROP	8.8E+01	PHYSPROP			5.1E-02	6.0E-06	WATER9 (U.S. EPA, 2001)			4.3E+02	EPI	3.0E+00	PHYSPROP	2.1E+02	PHYSPROP	4.6E-02	2.0E+00	4.7E+00	7.9E-03	EPI
Aminobiphenyl, 4-	92-67-1	1.7E+02	PHYSPROP	6.0E-06	1.5E-07	PHYSPROP	1.2E-04	PHYSPROP	5.4E+01	PHYSPROP			6.2E-02	7.3E-06	WATER9 (U.S. EPA, 2001)			2.5E+03	EPI	2.9E+00	PHYSPROP	2.2E+02	PHYSPROP	7.0E-02	9.3E-01	2.2E+00	1.4E-02	EPI
Aminophenol, m-	591-27-5	1.1E+02	PHYSPROP	8.1E-09	2.0E-10	PHYSPROP	9.6E-03	PHYSPROP	1.2E+02	PHYSPROP			8.3E-02	9.7E-06	WATER9 (U.S. EPA, 2001)			9.0E+01	EPI	2.1E-01	PHYSPROP	2.7E+04	PHYSPROP	2.1E-03	4.3E-01	1.0E+00	5.3E-04	EPI
Aminophenol, o-	95-55-6	1.1E+02	PHYSPROP	8.1E-09	2.0E-10	PHYSPROP	9.6E-03	PHYSPROP	1.7E+02	PHYSPROP	1.3E+00	CRC	8.0E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			9.2E+01	EPI	6.2E-01	PHYSPROP	2.0E+04	PHYSPROP	4.0E-03	4.3E-01	1.0E+00	9.9E-04	EPI
Aminophenol, p-	123-30-8	1.1E+02	PHYSPROP	1.5E-08	3.6E-10	EPI	4.0E-05	EPI	1.9E+02	PHYSPROP			8.3E-02	9.7E-06	WATER9 (U.S. EPA, 2001)			9.0E+01	EPI	4.0E-02	PHYSPROP	1.6E+04	PHYSPROP	1.6E-03	4.3E-01	1.0E+00	4.1E-04	EPI
Amitraz	33089-61-1	2.9E+02	PHYSPROP	4.0E-04	9.9E-06	PHYSPROP	2.0E-06	PHYSPROP	8.6E+01	PHYSPROP	1.1E+00	CRC	2.2E-02	5.4E-06	WATER9 (U.S. EPA, 2001)			2.6E+05	EPI	5.5E+00	PHYSPROP	1.0E+00	PHYSPROP	1.1E+00	4.6E+00	1.8E+01	1.6E-01	EPI
Ammonia	7664-41-7	1.7E+01	PHYSPROP	6.6E-04	1.6E-05	PHYSPROP	7.5E+03	PHYSPROP	-7.8E+01	PHYSPROP	7.0E-01	CRC	2.3E-01	2.2E-05	WATER9 (U.S. EPA, 2001)					2.3E-01	PHYSPROP	4.8E+05	PHYSPROP	1.6E-03	1.3E-01	3.1E-01	1.0E-03	RAGSE
Ammonium Picrate	131-74-8	2.3E+02	PHYSPROP	7.0E-10	1.7E-11	EPI	7.5E-07	PHYSPROP	1.2E+02	PHYSPROP	1.8E+00	PERRY	3.0E-02	8.2E-06	WATER9 (U.S. EPA, 2001)			2.3E+03	EPI	1.4E+00	PHYSPROP	1.3E+04	PHYSPROP	3.6E-03	2.0E+00	4.8E+00	6.2E-04	EPI
Ammonium Sulfamate	7773-06-0	1.1E+02	CRC				0.0E+00	NIOSH	1.3E+02	CRC	1.8E+00	PubChem										1.3E+06	PERRY	4.1E-03	4.6E-01	1.1E+00	1.0E-03	RAGSE
Amyl Alcohol, tert-	75-85-4	8.8E+01	PHYSPROP	5.6E-04	1.4E-05	PHYSPROP	1.7E+01	PHYSPROP	-9.1E+00	PHYSPROP	8.1E-01	CRC	7.9E-02	9.1E-06	WATER9 (U.S. EPA, 2001)			4.1E+00	EPI	8.9E-01	PHYSPROP	1.1E+05	PHYSPROP	7.1E-03	3.3E-01	7.9E-01	2.0E-03	EPI
Aniline	62-53-3	9.3E+01	PHYSPROP	8.3E-05	2.0E-06	PHYSPROP	6.7E-01	PHYSPROP	-6.0E+00	PHYSPROP	1.0E+00	CRC	8.3E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			7.0E+01	EPI	9.0E-01	PHYSPROP	3.6E+04	PHYSPROP	6.9E-03	3.5E-01	8.4E-01	1.9E-03	EPI
Anthraquinone, 9,10-	84-65-1	2.1E+02	PHYSPROP	9.6E-07	2.4E-08	EPI	1.2E-07	PHYSPROP	2.9E+02	PHYSPROP			5.4E-02	6.3E-06	WATER9 (U.S. EPA, 2001)			5.0E+03	EPI	3.4E+00	PHYSPROP	1.4E+00	PHYSPROP	1.1E-01	1.5E+00	3.7E+00	1.9E-02	EPI
Antimony (metallic)	7440-36-0	1.2E+02	CRC				0.0E+00	NIOSH	6.3E+02	PHYSPROP	6.7E+00	CRC				4.5E+01	SSL							4.2E-03	5.1E-01	1.2E+00	1.0E-03	RAGSE
Antimony Pentoxide	1314-60-9	3.2E+02	CRC								3.8E+00	CRC										3.0E+03	CRC	6.9E-03	6.8E+00	1.6E+01	1.0E-03	RAGSE
Antimony Tetroxide	1332-81-6	3.1E+02	EPI								6.6E+00	CRC												6.7E-03	5.5E+00	1.3E+01	1.0E-03	RAGSE
Antimony Trioxide	1309-64-4	2.9E+02	EPI						5.7E+02	CRC	5.6E+00	CRC												6.6E-03	4.5E+00	1.1E+01	1.0E-03	RAGSE
Arsenic, Inorganic	7440-38-2	7.5E+01	CRC						2.7E+02	CRC	4.9E+00	CRC				2.9E+01	SSL							3.3E-03	2.8E-01	6.6E-01	1.0E-03	RAGSE
Arsine	7784-42-1	7.8E+01	PHYSPROP						-1.2E+02	PHYSPROP	3.2E+00	CRC										2.0E+05	PERRY	3.4E-03	2.9E-01	6.9E-01	1.0E-03	RAGSE
Asbestos (units in fibers)	1332-21-4						0.0E+00	NIOSH																			1.0E-03	RAGSE
Asulam	3337-71-1	2.3E+02	PHYSPROP	7.0E-11	1.7E-12	PHYSPROP	1.4E-06	PHYSPROP	1.4E+02	PHYSPROP			5.1E-02	5.9E-06	WATER9 (U.S. EPA, 2001)			2.8E+01	EPI	-2.7E-01	PHYSPROP	5.0E+03	PHYSPROP	3.1E-04	2.0E+00	4.9E+00	5.3E-05	EPI
Atrazine	1912-24-9	2.2E+02	PHYSPROP	9.6E-08	2.4E-09	EPI	2.9E-07	PHYSPROP	1.7E+02	PHYSPROP	1.2E+00	PubChem	2.6E-02	6.8E-06	WATER9 (U.S. EPA, 2001)			2.2E+02	EPI	2.6E+00	PHYSPROP	3.5E+01	PHYSPROP	3.0E-02	1.7E+00	4.1E+00	5.2E-03	EPI
Auramine	492-80-8	2.7E+02	PHYSPROP	1.5E-07	3.6E-09	PHYSPROP	1.3E-06	PHYSPROP	1.4E+02	PHYSPROP			4.6E-02	5.3E-06	WATER9 (U.S. EPA, 2001)			4.5E+03	EPI	3.0E+00	PHYSPROP	5.4E+01	PHYSPROP	3.0E-02	3.3E+00	7.9E+00	4.8E-03	EPI
Avermectin B1	65195-55-3	8.8E+02	PHYSPROP	5.4E-26	1.3E-27	PHYSPROP	1.5E-30	PHYSPROP	3.5E+02	EPI			2.1E-02	2.4E-06	WATER9 (U.S. EPA, 2001)			8.8E+05	EPI	4.5E+00	PHYSPROP	3.5E-04	PHYSPROP	2.1E-04	8.4E+03	2.0E+04	1.8E-05	EPI
Azinphos-methyl	86-50-0	3.2E+02	PHYSPROP	9.8E-07	2.4E-08	EPI	1.6E-06	PHYSPROP	7.3E+01	PHYSPROP	1.4E+00	CRC	2.3E-02	5.6E-06	WATER9 (U.S. EPA, 2001)			5.2E+01	EPI	2.9E+00	PHYSPROP	2.1E+01	PHYSPROP	1.2E-02	6.3E+00	1.5E+01	1.8E-03	EPI
Azobenzene	103-33-3	1.8E+02	PHYSPROP	5.5E-04	1.4E-05	EPI	3.6E-04	PHYSPROP	6.8E+01	PHYSPROP	1.2E+00	PERRY	3.6E-02	7.5E-06	WATER9 (U.S. EPA, 2001)			3.8E+03	EPI	3.8E+00	PHYSPROP	6.4E+00	PHYSPROP	2.7E-01	1.1E+00	2.6E+00	5.1E-02	EPI
Azodicarbonamide	123-77-3	1.2E+02	PHYSPROP	3.4E-11	8.2E-13	EPI	1.9E-10	PHYSPROP	2.1E+02	EPI	1.7E+00	GuideChem	8.3E-02	1.2E-05	WATER9 (U.S. EPA, 2001)			7.0E+01	EPI	-1.7E+00	PHYSPROP	3.5E+01	PHYSPROP	1.1E-04	4.7E-01	1.1E+00	2.6E-05	EPI
Barium	7440-39-3	1.4E+02	EPI						7.1E+02	PHYSPROP	3.6E+00	CRC				4.1E+01	SSL							4.5E-03	6.2E-01	1.5E+00	1.0E-03	RAGSE
Benfluralin	1861-40-1	3.4E+02	PHYSPROP	1.2E-02	2.9E-04	EPI	6.5E-05	PHYSPROP	6.6E+01	PHYSPROP	1.3E+00	ChemNet	2.2E-02	5.5E-06	WATER9 (U.S. EPA, 2001)			1.6E+04	EPI	5.3E+00	PHYSPROP	1.0E-01	PHYSPROP	4.8E-01	7.9E+00	1.9E+01	6.8E-02	EPI
Benomyl	17804-35-2	2.9E+02	PHYSPROP	2.0E-10	4.9E-12	PHYSPROP	3.7E-09	PHYSPROP	1.4E+02	EPI			4.3E-02	5.1E-06	WATER9 (U.S. EPA, 2001)			3.4E+02	EPI	2.1E+00	PHYSPROP	3.8E+00	PHYSPROP	6.2E-03	4.4E+00	1.1E+01	9.4E-04	EPI
Bensulfuron-methyl	83055-99-6	4.1E+02	PHYSPROP	1.5E-13	3.8E-15	EPI	2.1E-14	PHYSPROP	1.9E+02	PHYSPROP			3.4E-02	4.0E-06	WATER9 (U.S. EPA, 2001)			2.8E+01	EPI	2.2E+00	PHYSPROP	1.2E+02	PHYSPROP	1.7E-03	2.1E+01	5.0E+01	2.2E-04	EPI
Bentazon	25057-89-0	2.4E+02	PHYSPROP	8.9E-08	2.2E-09	EPI	1.6E-06	PHYSPROP	1.4E+02	PHYSPROP			4.9E-02	5.1E-06	WATER9 (U.S. EPA, 2001)			1.0E+01	EPI	2.3E+00	PHYSPROP	5.0E+02	PHYSPROP	1.5E-02	2.3E+00	5.6E+00	2.5E-03	EPI
Benzaldehyde	100-52-7	1.1E+02	PHYSPROP	1.1E-03	2.7E-05	PHYSPROP	1.3E+00	PHYSPROP	-2.6E+01	PHYSPROP	1.0E+00	CRC	7.4E-02	9.5E-06	WATER9 (U.S. EPA, 2001)			1.1E+01	EPI	1.5E+00	PHYSPROP	7.0E+03	PHYSPROP	1.5E-02	4.1E-01	9.9E-01	3.8E-03	EPI
Benzene	71-43-2	7.8E+01	PHYSPROP	2.3E-01	5.6E-03	PHYSPROP	9.5E+01	PHYSPROP	5.5E+00	PHYSPROP	8.8E-01	CRC	9.0E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.5E+02	EPI	2.1E+00	PHYSPROP	1.8E+03	PHYSPROP	5.1E-02	2.9E-01	6.9E-01	1.5E-02	EPI
Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1	2.2E+02	EPI	8.9E-22	2.2E-23	EPI	2.9E-14	EPI	2.4E+02	EPI			5.2E-02	6.1E-06	WATER9 (U.S. EPA, 2001)			3.8E+01	EPI	-3.7E+00	EPI	1.0E+06	EPI	1.7E-06	1.8E+00	4.3E+00	3.0E-07	EPI
Benzenethiol	108-98-5	1.1E+02	PHYSPROP	1.4E-02	3.4E-04	EPI	1.9E+00	PHYSPROP	-1.5E+01	PHYSPROP	1.1E+00	CRC	7.3E-02	9.5E-06	WATER9 (U.S. EPA, 2001)			2.3E+02	EPI	2.5E+00	PHYSPROP	8.4E+02	PHYSPROP	7.2E-02	4.4E-01	1.0E+00	1.8E-02	EPI
Benzidine	92-87-5	1.8E+02	PHYSPROP	2.1E-09	5.2E-11	PHYSPROP	9.0E-07	PHYSPROP	1.2E+02	PHYSPROP	1.2E+00	Yaws 2008	3.5E-02	7.5E-06	WATER9 (U.S. EPA, 2001)			1.2E+03	EPI	1.3E+00	PHYSPROP	3.2E+02	PHYSPROP	5.9E-03	1.1E+00	2.7E+00	1.1E-03	EPI
Benzoic Acid	65-85-0	1.2E+02	PHYSPROP	1.6E-06	3.8E-08	EPI	7.0E-04	PHYSPROP	1.2E+02	PHYSPROP	1.3E+00	CRC	7.0E-02	9.8E-06	WATER9 (U.S. EPA, 2001)			6.0E-01	SSL	1.9E+00	PHYSPROP	3.4E+03	PHYSPROP	2.4E-02	5.1E-01	1.2E+00	5.7E-03	EPI
Benzotrichloride	98-07-7	2.0E+02	PHYSPROP	1.1E-02	2.6E-04	PHYSPROP	4.1E-01	EPI	-5.0E+00	PHYSPROP	1.4E+00	CRC	3.1E-02	7.7E-06	WATER9 (U.S. EPA, 2001)			1.0E+03	EPI	3.9E+00	PHYSPROP	5.3E+01	PHYSPROP	2.6E-01	1.3E+00	3.1E+00	4.9E-02	EPI
Benzyl Alcohol	100-51-6	1.1E+02	PHYSPROP	1.4E-05	3.4E-07	PHYSPROP	9.4E-02	PHYSPROP	-1.5E+01	PHYSPROP	1.0E+00	CRC	7.3E-02	9.4E-06	WATER9 (U.S. EPA, 2001)			2.1E+01	EPI	1.1E+00	PHYSPROP	4.3E+04	PHYSPROP	8.4E-03	4.2E-01	1.0E+00	2.1E-03	EPI
Benzyl Chloride	100-44-7	1.3E+02	PHYSPROP	1.7E-02	4.1E-04	EPI	1.2E+00	PHYSPROP	-4.5E+01	PHYSPROP	1.1E+00	CRC	6.3E-02	8.8E-06	WATER9 (U.S. EPA, 2001)			4.5E+02	EPI	2.3E+00	PHYSPROP	5.3E+02	PHYSPROP	4.5E-02	5.4E-01	1.3E+00	1.0E-02	EPI
Beryllium and compounds	7440-41-7	9.0E+00	EPI				0.0E+00	NIOSH	9.9E+02	PHYSPROP	1.9E+00	CRC				7.9E+02	SSL							1.2E-03	1.2E-01	2.8E-01	1.0E-03	RAGSE
Bifenox	42576-02-3	3.4E+02	PHYSPROP	4.4E-06	1.1E-07	EPI	1.0E-07	PHYSPROP	8.5E+01	PHYSPROP	1.2E+00	PubChem	2.0E-02	5.0E-06	WATER9 (U.S. EPA, 2001)			3.7E+03	EPI	4.5E+00	PHYSPROP	4.0E-01	PHYSPROP	1.3E-01	8.7E+00	2.1E+01	1.8E-02	EPI
Biphenthrin	82657-04-3	4.2E+02	PHYSPROP	4.1E-05	1.0E-06	EPI	1.8E-07	PHYSPROP	6.9E+01	PHYSPROP	1.2E+00	CRC	1.8E-02	4.5E-06	WATER9 (U.S. EPA, 2001)			2.3E+06	EPI	6.0E+00	PHYSPROP	1.0E-03	PHYSPROP	1.4E+01	2.5E+01	1.1E+02	1.7E+00	EPI
Biphenyl, 1,1'-	92-52-4	1.5E+02	PHYSPROP	1.3E-02	3.1E-04	PHYSPROP	8.9E-03	PHYSPROP	6.9E+01	PHYSPROP	1.0E+00	CRC	4.7E-02	6.1E-06	WATER9 (U.S. EPA, 2001)			5.1E+03	EPI	4.0E+00	PHYSPROP	7.5E+00	PHYSPROP	4.5E-01	7.7E-01	1.8E+00	9.4E-02	EPI
Bis(2-chloro-1-methylethyl) ether	108-60-1	1.7E+02	PHYSPROP	3.0E-03	7.4E-05	EPI	5.6E-01	PHYSPROP	-9.7E+01	PHYSPROP	1.1E+00	CRC	4.0E-02	7.4E-06	WATER9 (U.S. EPA, 2001)			8.3E+01	EPI	2.5E+00	PHYSPROP	1.7E+03	PHYSPROP	3.8E-02	9.5E-01	2.3E+00	7.6E-03	EPI
Bis(2-chloroethoxy)methane	111-91-1	1.7E+02	PHYSPROP	1.6E-04	3.9E-06	EPI	1.3E-01	EPI	-3.2E+01	PHYSPROP			6.1E-02	7.1E-06	WATER9 (U.S. EPA, 2001)			1.4E+01	EPI	1.3E+00	PHYSPROP	7.8E+03	PHYSPROP	6.2E-03	9.8E-01	2.4E+00	1.2E-03	EPI
Bis(2-chloroethyl)ether	111-44-4	1.4E+02	PHYSPROP	7.0E-04	1.7E-05	EPI	1.6E+00	PHYSPROP	-5.2E+01	PHYSPROP	1.2E+00	CRC	5.7E-02	8.7E-06	WATER9 (U.S. EPA, 2001)			3.2E+01	EPI	1.3E+00	PHYSPROP	1.7E+04	PHYSPROP	8.2E-03	6.6E-01	1.6E+00	1.8E-03	EPI
Bis(chloromethyl)ether	542-88-1	1.1E+02	PHYSPROP	1.8E-01	4.4E-03	EPI	2.9E+01	PHYSPROP	-4.2E+01	PHYSPROP	1.3E+00	CRC	7.6E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			9.7E+00	EPI	5.7E-01	PHYSPROP	2.2E+04	PHYSPROP	3.5E-03	4.6E-01	1.1E+00	8.6E-04	EPI
Bisphenol A	80-05-7	2.3E+02	PHYSPROP	4.1E-10	1.0E-11	PHYSPROP	3.9E-07	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	PubChem	2.5E-02	6.5E-06	WATER9 (U.S. EPA, 2001)			3.8E+04	EPI	3.3E+00	PHYSPROP	1.2E+02	PHYSPROP	7.7E-02	2.0E+00	4.8E+00	1.3E-02	EPI
Boron Acid Borates Only	7440-42-8	1.4E+01	CRC						2.1E+03	CRC	2.3E+00	CRC				3.0E+00	BAES							1.4E-03	1.3E-01	3.0E-01	1.0E-03	RAGSE
Boron Trichloride	10294-34-5	1.2E+02	PHYSPROP	7.5E-01	1.8E-02	PHYSPROP	1.0E+00	PHYSPROP	-1.1E+02	PHYSPROP	4.8E-03	CRC	1.2E-01	2.2E-05	WATER9 (U.S. EPA, 2001)					1.2E+00	PHYSPROP			4.2E-03	4.8E-01	1.1E+00	1.0E-03	RAGSE
Boron Trifluoride	7637-07-2

Contaminant		Molecular Weight		Volatility Parameters					Melting Point		Density		Diffusivity in Air and Water			Partition Coefficients						Water Solubility		Tapwater Dermal Parameters				
Analyte	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m^3/mole)	H' and HLC Ref	VP (mmHg)	VP Ref	MP (C)	MP Ref	Density (g/cm^3)	Density Ref	D_{oa} (cm^2/s)	D_{iw} (cm^2/s)	D_{ia} and D_{iw} Ref	K_d (L/kg)	K_d Ref	K_{oc} (L/kg)	K_{oc} Ref	log K_{ow} (unitless)	log K_{ow} Ref	S (mg/L)	S Ref	B (unitless)	t_{event} (hr/event)	t^* (hr)	K_p (cm/hr)	K_p Ref
Carbosulfan	55285-14-8	3.8E+02	PHYSPROP	2.1E-05	5.1E-07	EPI	3.1E-07	PHYSPROP	2.5E+01	PHYSPROP	1.1E+00	CRC	1.8E-02	4.4E-06	WATER9 (U.S. EPA, 2001)			1.2E+04	EPI	5.6E+00	PHYSPROP	3.0E-01	PHYSPROP	4.3E-01	1.4E+01	3.4E+01	5.8E-02	EPI
Carboxin	5234-68-4	2.4E+02	PHYSPROP	1.3E-08	3.2E-10	EPI	1.5E-07	PHYSPROP	9.2E+01	PHYSPROP			5.0E-02	5.8E-06	WATER9 (U.S. EPA, 2001)			1.7E+02	EPI	2.1E+00	PHYSPROP	1.5E+02	PHYSPROP	1.2E-02	2.2E+00	5.2E+00	2.0E-03	EPI
Ceric oxide	1306-38-3	1.7E+02	CRC						2.5E+03	CRC	7.2E+00	CRC												5.0E-03	9.7E-01	2.3E+00	1.0E-03	RAGSE
Chloral Hydrate	302-17-0	1.7E+02	PHYSPROP	2.3E-07	5.7E-09	PHYSPROP	1.5E+01	PHYSPROP	5.7E+01	PHYSPROP	1.9E+00	CRC	5.4E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	9.9E-01	PHYSPROP	7.9E+05	PHYSPROP	4.2E-03	8.9E-01	2.1E+00	8.4E-04	EPI
Chloramben	133-90-4	2.1E+02	PHYSPROP	1.6E-09	3.9E-11	EPI	1.0E-07	PHYSPROP	2.0E+02	PHYSPROP			5.4E-02	6.4E-06	WATER9 (U.S. EPA, 2001)			2.1E+01	EPI	1.9E+00	PHYSPROP	7.0E+02	PHYSPROP	1.1E-02	1.5E+00	3.6E+00	2.0E-03	EPI
Chloramines, Organic	E701235																											
Chloranil	118-75-2	2.5E+02	PHYSPROP	1.3E-08	3.3E-10	PHYSPROP	2.3E-06	PHYSPROP	2.9E+02	PHYSPROP			4.8E-02	5.7E-06	WATER9 (U.S. EPA, 2001)			3.1E+02	EPI	2.2E+00	PHYSPROP	2.5E+02	PHYSPROP	1.2E-02	2.5E+00	6.0E+00	1.9E-03	EPI
Chlordane (alpha)	5103-71-9	4.1E+02	PHYSPROP	2.0E-03	4.9E-05	EPI	3.6E-05	PHYSPROP	1.1E+02	EPI	1.8E+00	LookChem	2.3E-02	5.8E-06	WATER9 (U.S. EPA, 2001)			6.8E+04	EPI	6.1E+00	PHYSPROP	5.6E-02	PHYSPROP	2.6E-01	2.1E+01	5.0E+01	3.4E-02	RAGSE
Chlordane (gamma)	5103-74-2	4.1E+02	PHYSPROP	2.0E-03	4.9E-05	EPI	5.0E-05	PHYSPROP	1.1E+02	EPI	1.8E+00	ChemNet	2.3E-02	5.8E-06	WATER9 (U.S. EPA, 2001)			6.8E+04	EPI	6.2E+00	PHYSPROP	5.6E-02	PHYSPROP	2.6E-01	2.1E+01	5.0E+01	3.4E-02	RAGSE
Chlordane (technical mixture)	12789-03-6	4.1E+02	PHYSPROP	2.0E-03	4.9E-05	EPI	1.0E-05	PHYSPROP	1.1E+02	EPI	1.6E+00	CRC	2.1E-02	5.4E-06	WATER9 (U.S. EPA, 2001)			6.8E+04	EPI	6.2E+00	EPI	5.6E-02	EPI	8.3E-01	2.1E+01	8.0E+01	1.1E-01	EPI
Chlordecone (Kepone)	143-50-0	4.9E+02	PHYSPROP	2.2E-06	5.4E-08	EPI	2.3E-07	PHYSPROP	3.5E+02	EPI	1.6E+00	CRC	2.0E-02	4.9E-06	WATER9 (U.S. EPA, 2001)			1.8E+04	EPI	5.4E+00	PHYSPROP	2.7E+00	PHYSPROP	9.3E-02	5.9E+01	1.4E+02	1.1E-02	EPI
Chlorfenvinphos	470-90-6	3.6E+02	PHYSPROP	1.2E-06	2.9E-08	EPI	7.5E-06	PHYSPROP	-2.0E+01	PHYSPROP			3.8E-02	4.4E-06	WATER9 (U.S. EPA, 2001)			1.3E+03	EPI	3.8E+00	PHYSPROP	1.2E+02	PHYSPROP	3.7E-02	1.1E+01	2.6E+01	5.1E-03	EPI
Chlorimuron, Ethyl-	90982-32-4	4.1E+02	PHYSPROP	7.4E-14	1.8E-15	EPI	4.0E-12	PHYSPROP	1.8E+02	PHYSPROP			3.4E-02	4.0E-06	WATER9 (U.S. EPA, 2001)			7.2E+01	EPI	2.5E+00	PHYSPROP	1.2E+03	PHYSPROP	2.6E-03	2.2E+01	5.3E+01	3.4E-04	EPI
Chlorine	7782-50-5	7.1E+01	PHYSPROP	4.8E-01	1.2E-02	PHYSPROP	5.9E+03	PHYSPROP	-1.0E+02	PHYSPROP	2.9E+00	CRC	1.5E-01	2.2E-05	WATER9 (U.S. EPA, 2001)	2.5E-01	BAES			8.5E-01	PHYSPROP	6.3E+03	PHYSPROP	3.2E-03	2.6E-01	6.3E-01	1.0E-03	RAGSE
Chlorine Dioxide	10049-04-4	6.7E+01	EPI	1.6E+00	4.0E-02	Toxnet HSDB	7.6E+02	Toxnet HSDB	-5.9E+01	CRC	2.8E+00	CRC	1.6E-01	2.2E-05	WATER9 (U.S. EPA, 2001)							8.0E+03	PubChem	3.2E-03	2.5E-01	6.0E-01	1.0E-03	RAGSE
Chlorite (Sodium Salt)	7758-19-2	9.0E+01	EPI						1.8E+02	CRC												6.4E+05	CRC	3.7E-03	3.4E-01	8.1E-01	1.0E-03	RAGSE
Chloro-1,1-difluoroethane, 1-	75-68-3	1.0E+02	PHYSPROP	2.4E+00	5.9E-02	PHYSPROP	2.5E+03	PHYSPROP	-1.3E+02	PHYSPROP	1.1E+00	CRC	8.0E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			4.4E+01	EPI	2.1E+00	PHYSPROP	1.4E+03	PHYSPROP	3.8E-02	3.8E-01	9.2E-01	9.9E-03	EPI
Chloro-1,3-butadiene, 2-	126-99-8	8.9E+01	PHYSPROP	2.3E+00	5.6E-02	PHYSPROP	2.2E+02	PHYSPROP	-1.3E+02	PHYSPROP	9.6E-01	CRC	8.4E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			6.1E+01	EPI	2.5E+00	PHYSPROP	8.7E+02	PHYSPROP	8.6E-02	3.3E-01	7.9E-01	2.4E-02	EPI
Chloro-2-methylaniline HCl, 4-	3165-93-3	1.8E+02	PHYSPROP	6.4E-05	1.6E-06	PHYSPROP	4.1E-02	PHYSPROP	1.6E+02	EPI			6.0E-02	7.0E-06	WATER9 (U.S. EPA, 2001)			3.5E+02	EPI	2.3E+00	PHYSPROP	9.5E+02	PHYSPROP	9.2E-05	1.0E+00	2.5E+00	1.8E-05	EPI
Chloro-2-methylaniline, 4-	95-69-2	1.4E+02	PHYSPROP	8.1E-05	2.0E-06	PHYSPROP	4.1E-02	PHYSPROP	3.0E+01	PHYSPROP			7.0E-02	8.2E-06	WATER9 (U.S. EPA, 2001)			1.8E+02	EPI	2.3E+00	PHYSPROP	9.5E+02	PHYSPROP	3.7E-02	6.5E-01	1.6E+00	8.1E-03	EPI
Chloroacetaldehyde, 2-	107-20-0	7.8E+01	PHYSPROP	9.8E-04	2.4E-05	PHYSPROP	6.4E+01	PHYSPROP	-1.6E+01	PHYSPROP	1.2E+00	CRC	1.0E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	9.0E-02	PHYSPROP	1.1E+05	PHYSPROP	2.2E-03	2.9E-01	6.9E-01	6.5E-04	EPI
Chloroacetic Acid	79-11-8	9.4E+01	PHYSPROP	3.8E-07	9.3E-09	PHYSPROP	6.5E-02	PHYSPROP	6.3E+01	PHYSPROP	1.4E+00	CRC	9.4E-02	1.2E-05	WATER9 (U.S. EPA, 2001)			1.4E+00	EPI	2.2E-01	PHYSPROP	8.6E+05	PHYSPROP	2.4E-03	3.6E-01	8.5E-01	6.5E-04	EPI
Chloroacetophenone, 2-	532-27-4	1.5E+02	PHYSPROP	1.4E-04	3.5E-06	PHYSPROP	5.4E-03	PHYSPROP	5.7E+01	PHYSPROP	1.3E+00	CRC	5.2E-02	8.7E-06	WATER9 (U.S. EPA, 2001)			9.9E+01	EPI	1.9E+00	PHYSPROP	1.1E+03	PERRY	1.9E-02	7.7E-01	1.9E+00	4.1E-03	EPI
Chloroaniline, p-	106-47-8	1.3E+02	PHYSPROP	4.7E-05	1.2E-06	EPI	2.7E-02	PHYSPROP	7.3E+01	PHYSPROP	1.4E+00	CRC	7.0E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.1E+02	EPI	1.8E+00	PHYSPROP	3.9E+03	PHYSPROP	2.2E-02	5.4E-01	1.3E+00	5.0E-03	EPI
Chlorobenzene	108-90-7	1.1E+02	PHYSPROP	1.3E-01	3.1E-03	PHYSPROP	1.2E+01	PHYSPROP	-4.5E+01	PHYSPROP	1.1E+00	CRC	7.2E-02	9.5E-06	WATER9 (U.S. EPA, 2001)			2.3E+02	EPI	2.8E+00	PHYSPROP	5.0E+02	PHYSPROP	1.2E-01	4.5E-01	1.1E+00	2.8E-02	EPI
Chlorobenzene sulfonic acid, p-	98-66-8	1.9E+02	PHYSPROP	7.6E-08	1.9E-09	PHYSPROP	4.3E-06	PHYSPROP	6.7E+01	PHYSPROP			5.7E-02	6.7E-06	WATER9 (U.S. EPA, 2001)			1.6E+01	EPI	-5.2E-01	PHYSPROP	3.1E+05	PHYSPROP	3.1E-04	1.3E+00	3.0E+00	5.9E-05	EPI
Chlorobenzilate	510-15-6	3.3E+02	PHYSPROP	3.0E-06	7.2E-08	EPI	2.2E-06	PHYSPROP	3.7E+01	PHYSPROP	1.3E+00	CRC	3.2E-02	5.5E-06	WATER9 (U.S. EPA, 2001)			1.5E+03	EPI	4.7E+00	PHYSPROP	1.3E+01	PHYSPROP	2.3E-01	7.0E+00	1.7E+01	3.3E-02	EPI
Chlorobenzoic Acid, p-	74-11-3	1.6E+02	PHYSPROP	3.3E-06	8.0E-08	PHYSPROP	2.3E-03	PHYSPROP	2.4E+02	PHYSPROP	1.5E+00	PERRY	5.5E-02	9.5E-06	WATER9 (U.S. EPA, 2001)			2.7E+01	EPI	2.7E+00	PHYSPROP	7.2E+01	PHYSPROP	5.8E-02	7.9E-01	1.9E+00	1.2E-02	EPI
Chlorobenzotrifluoride, 4-	98-56-6	1.8E+02	PHYSPROP	1.4E+00	3.5E-02	PHYSPROP	7.6E+00	PHYSPROP	-3.3E+01	PHYSPROP	1.3E+00	CRC	3.8E-02	8.0E-06	WATER9 (U.S. EPA, 2001)			1.6E+03	EPI	3.6E+00	PHYSPROP	2.9E+01	PHYSPROP	1.9E-01	1.1E+00	2.6E+00	3.8E-02	EPI
Chlorobutane, 1-	109-69-3	9.3E+01	PHYSPROP	6.8E-01	1.7E-02	PHYSPROP	1.0E+02	PHYSPROP	-1.2E+02	PHYSPROP	8.9E-01	CRC	7.8E-02	9.3E-06	WATER9 (U.S. EPA, 2001)			7.2E+01	EPI	2.6E+00	PHYSPROP	1.1E+03	PHYSPROP	1.0E-01	3.5E-01	8.3E-01	2.7E-02	EPI
Chlorodifluoromethane	75-45-6	8.6E+01	PHYSPROP	1.7E+00	4.1E-02	PHYSPROP	7.3E+03	PHYSPROP	-1.6E+02	PHYSPROP	1.5E+00	CRC	1.0E-01	1.3E-05	WATER9 (U.S. EPA, 2001)			3.2E+01	EPI	1.1E+00	PHYSPROP	2.8E+03	PHYSPROP	9.6E-03	3.2E-01	7.7E-01	2.7E-03	EPI
Chloroethanol, 2-	107-07-3	8.1E+01	PHYSPROP	3.1E-05	7.6E-07	EPI	7.2E+00	PHYSPROP	-6.8E+01	PHYSPROP	1.2E+00	CRC	1.0E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			1.9E+00	EPI	3.0E-02	PHYSPROP	1.0E+06	PHYSPROP	2.0E-03	3.0E-01	7.1E-01	5.8E-04	EPI
Chloroform	67-66-3	1.2E+02	PHYSPROP	1.5E-01	3.7E-03	PHYSPROP	2.0E+02	PHYSPROP	-5.4E+01	PHYSPROP	1.5E+00	CRC	7.0E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			3.2E+01	EPI	2.0E+00	PHYSPROP	8.0E+03	PHYSPROP	2.9E-02	4.9E-01	1.2E+00	6.8E-03	EPI
Chloromethane	74-87-3	5.0E+01	PHYSPROP	3.6E-01	8.8E-03	PHYSPROP	4.3E+03	PHYSPROP	-9.8E+01	PHYSPROP	9.1E-01	CRC	1.2E-01	1.4E-05	WATER9 (U.S. EPA, 2001)			1.3E+01	EPI	9.1E-01	PHYSPROP	5.3E+03	PHYSPROP	9.0E-03	2.0E-01	4.8E-01	3.3E-03	EPI
Chloromethyl Methyl Ether	107-30-2	8.1E+01	PHYSPROP	1.2E-02	3.0E-04	PHYSPROP	3.0E+01	PHYSPROP	-1.0E+02	PHYSPROP	1.1E+00	CRC	9.5E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			5.3E+00	EPI	3.2E-01	PHYSPROP	6.9E+04	PHYSPROP	3.1E-03	3.0E-01	7.1E-01	9.1E-04	EPI
Chloronitrobenzene, o-	88-73-3	1.6E+02	PHYSPROP	3.8E-04	9.3E-06	PHYSPROP	1.8E-02	EPI	3.3E+01	PHYSPROP	1.4E+00	CRC	5.1E-02	8.8E-06	WATER9 (U.S. EPA, 2001)			3.7E+02	EPI	2.2E+00	PHYSPROP	4.4E+02	PHYSPROP	3.0E-02	8.0E-01	1.9E+00	6.3E-03	EPI
Chloronitrobenzene, p-	100-00-5	1.6E+02	PHYSPROP	2.0E-04	4.9E-06	PHYSPROP	2.2E-02	EPI	8.4E+01	PHYSPROP	1.3E+00	CRC	5.0E-02	8.5E-06	WATER9 (U.S. EPA, 2001)			3.6E+02	EPI	2.4E+00	PHYSPROP	2.3E+02	PHYSPROP	3.8E-02	8.0E-01	1.9E+00	7.9E-03	EPI
Chlorophenol, 2-	95-57-8	1.3E+02	PHYSPROP	4.6E-04	1.1E-05	PHYSPROP	2.5E+00	PHYSPROP	9.8E+00	PHYSPROP	1.3E+00	CRC	6.6E-02	9.5E-06	WATER9 (U.S. EPA, 2001)			3.9E+02	SSL	2.2E+00	PHYSPROP	1.1E+04	PHYSPROP	3.5E-02	5.5E-01	1.3E+00	8.0E-03	EPI
Chloropicrin	76-06-2	1.6E+02	PHYSPROP	8.4E-02	2.1E-03	PHYSPROP	2.4E+01	PHYSPROP	-6.4E+01	PHYSPROP	1.7E+00	CRC	5.2E-02	9.6E-06	WATER9 (U.S. EPA, 2001)			4.4E+01	EPI	2.1E+00	PHYSPROP	1.6E+03	PHYSPROP	2.3E-02	8.8E-01	2.1E+00	4.6E-03	EPI
Chlorotrilonil	1897-45-6	2.7E+02	PHYSPROP	8.2E-05	2.0E-06	PHYSPROP	5.7E-07	PHYSPROP	2.5E+02	PHYSPROP	1.7E+00	CRC	3.8E-02	7.3E-06	WATER9 (U.S. EPA, 2001)			1.0E+03	EPI	3.1E+00	PHYSPROP	8.1E-01	PHYSPROP	3.4E-02	3.2E+01	7.8E+00	5.4E-03	EPI
Chlorotoluene, o-	95-49-8	1.3E+02	PHYSPROP	1.5E-01	3.6E-03	PHYSPROP	3.4E+00	PHYSPROP	-3.6E+01	PHYSPROP	1.1E+00	CRC	6.3E-02	8.7E-06	WATER9 (U.S. EPA, 2001)			3.8E+02	EPI	3.4E+00	PHYSPROP	3.7E+02	PHYSPROP	2.5E-01	5.4E-01	1.3E+00	5.7E-02	EPI
Chlorotoluene, p-	106-43-4	1.3E+02	PHYSPROP	1.8E-01	4.4E-03	EPI	2.7E+00	PHYSPROP	7.5E+00	PHYSPROP	1.1E+00	CRC	6.3E-02	8.7E-06	WATER9 (U.S. EPA, 2001)			3.8E+02	EPI	3.3E+00	PHYSPROP	1.1E+02	PHYSPROP	2.2E-01	5.4E-01	1.3E+00	5.0E-02	EPI
Chlorozotocin	54749-90-5	2.7E+02	PHYSPROP	1.5E-20	3.7E-22	PHYSPROP	4.0E-14	PHYSPROP	1.5E+02	EPI			4.6E-02	5.4E-06	WATER9 (U.S. EPA, 2001)			1.0E+01	EPI	-1.0E+00	PHYSPROP	1.8E+03	PHYSPROP	6.2E-05	3.2E+00	7.8E+00	9.9E-06	EPI
Chloropropham	101-21-3	2.1E+02	PHYSPROP	2.3E-05	5.7E-07	EPI	1.8E-04	PHYSPROP	4.1E+01	PHYSPROP	1.2E+00	CRC	2.6E-02	6.7E-06	WATER9 (U.S. EPA, 2001)			3.5E+02	EPI	3.5E+00	PHYSPROP	8.9E+01	PHYSPROP	1.2E-01	1.7E+00	4.0E+00	2.1E-02	EPI
Chlorpyrifos	2921-88-2	3.5E+02	PHYSPROP	1.2E-04	2.9E-06	PHYSPROP	2.0E-05	PHYSPROP	4.2E+01	PHYSPROP			3.8E-02	4.5E-06	WATER9 (U.S. EPA, 2001)			7.3E+03	EPI	5.0E+00	PHYSPROP	1.1E+00	PHYSPROP	2.4E-01	9.7E+00	2.3E+01	3.3E-02	EPI
Chlorpyrifos Methyl	5598-13-0	3.2E+02	PHYSPROP	1.5E-04	3.8E-06	EPI	4.2E-05	PHYSPROP	4.3E+01	PHYSPROP			4.0E-02	4.7E-06	WATER9 (U.S. EPA, 2001)			2.0E+03	EPI	4.3E+00	PHYSPROP	4.9E+00	PHYSPROP	1.2E-01	6.7E+00	1.6E+01	1.8E-02	EPI
Chlorsulfuron	64902-72-3	3.6E+02	PHYSPROP	1.4E-14	3.4E-16	EPI	2.3E-11	PHYSPROP	1.8E+02	PHYSPROP			3.8E-02	4.4E-06	WATER9 (U.S. EPA, 2001)			3.2E+02	EPI	2.0E+00	PHYSPROP	3.1E+04	PHYSPROP	2.4E-03	1.1E+01	2.5E+01	3.3E-04	EPI
Chlorthal-dimethyl	1861-32-1	3.3E+02	PHYSPROP	8.9E-05	2.2E-06	EPI	2.5E-06	PHYSPROP	1.6E+02	PHYSPROP			4.0E-02	4.6E-06	WATER9 (U.S. EPA, 2001)			5.1E+02	EPI	4.3E+00	PHYSPROP	5.0E-01	PHYSPROP	1.1E-01	7.6E+00	1.8E+01	1.5E-02	EPI
Chlorthiophos	60238-56-4	3.6E+02	PHYSPROP	4.9E-05	1.2E-06	PHYSPROP	4.0E-01	PHYSPROP	8.6E+01	EPI			3.7E-02	4.4E-06	WATER9 (U.S. EPA, 2001)			1.3E+04	EPI	5.8E+00	PHYSPROP	3.0E-01	PHYSPROP	7.7E-01	1.1E+01	4.3E+01	1.1E-01	EPI
Chromium(III), Insoluble Salts	16065-83-1	5.2E+01	EPI								5.2E+00	CRC				1.8E+06	SSL							2.8E-03	2.1E-01	4.9E-01	1.0E-03	RAGSE
Chromium(VI)	18540-29-9	5.2E+01	EPI													1.9E+01	SSL					1.7E+06	CRC	5.5E-03	2.1E-01	4.9E-01	2.0E-03	RAGSE
Chromium, Total	7440-47-3	5.2E+01	PHYSPROP						1.9E+03	PHYSPROP	7.2E+00	CRC				1.8E+06	SSL							2.8E-03	2.1E-01	4.9E-01	1.0E-03	RAGSE
Clofentezine	74115-24-5	3.0E+02	PHYSPROP	1.6E-08	3.9E-10	EPI	9.8E-10	PHYSPROP	1.8E+02	PHYSPROP			4.2E-02	4.9E-06	WATER9 (U.S. EPA, 2001)			3.0E+04	EPI	3.1E+00	PHYSPROP	1.0E+00	PHYSPROP	2.4E-02	5.2E+00	1.3E+01	3.6E-03	EPI
Cobalt	7440-48-4	5.9E+01	EPI				0.0E+00	NIOSH	1.5E+03	CRC	8.9E+00	CRC				4.5E+01	BAES							1.2E-03	2.2E-01	5.4E-01	4.0E-04	RAGSE
Coke Oven Emissions	E649830			4.5E-01	1.1E-02	Toxnet HSDB	9.5E+01	Toxnet HSDB					1.0E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			1.6E+04										
Copper	7440-50-8	6.4E+01	PHYSPROP				0.0E+00	NIOSH	1.1E+03	PHYSPROP	9.0E+00	CRC				3.5E+01	BAES							3.1E-03	2.4E-01	5.7E-01	1.0E-03	RAGSE
Cresol, m-	108-39-4	1.1E+02	PHYSPROP	3.5E-05	8.6E-07	PHYSPROP	1.1E-01	PHYSPROP	1.2E+01	PHYSPROP	1.0E+00	CRC	7.3E-02	9.3E-06	WATER9 (U.S. EPA, 2001)			3.0E+02	EPI	2.0E+00	PHYSPROP	2.3E+04	PHYSPROP	3.1E-02	4.2E-01	1.0E+00	7.8E-03	EPI
Cresol, o-	95-48-7	1.1E+02	PHYSPROP	4.9E-05	1.2E-06	PHYSPROP	3.0E-01	EPI	3.0E+01	PHYSPROP	1.0E+00	CRC	7.3E-02	9.3E-06	WATER9 (U.S. EPA, 2001)			3.1E+02	EPI	2.0E+00	PHYSPROP	2.6E+04	PHYSPROP	3.1E-02	4.2E-01	1.0E+00	7.7E-03	EPI
Cresol, p-	106-44-5	1.1E+02	PHYSPROP	4.1E-05	1.0E-06	PHYSPROP	1.1E-01	PHYSPROP	3.6E+01	PHYSPROP	1.0E+00	CRC	7.2E-02	9.2E-06	WATER9 (U.S. EPA, 2001)			3.0E+02	EPI	1.9E+00	PHYSPROP	2.2E+04	PHYSPROP	3.0E-02	4.2E-01	1.0E+00	7.5E-03	EPI
Cresol, p-chloro-m-	59-50-7	1.4E+02	PHYSPROP	1.0E-04	2.5E-06	EPI	5.0E-02	PHYSPROP	6.7E+01	PHYSPROP			7.0E-02	8.1E-06	WATER9 (U.S. EPA, 2001)			4.9E+02	EPI	3.1E+00	PHYSPROP	3.8E+03	PHYSPROP	1.3E-01	6.6E-01	1.6E+00	2.9E-02	EPI
Cresols	1319-77-3	3.2E+02	PHYSPROP	2.5E-05	6.2E-07	PHYSPROP	1.7E-01	PHYSPROP	3.0E+01	EPI			4.0E-02	4.7E-06	WATER9 (U.S. EPA, 2001)			3.1E+02	EPI	2.0E+00	PHYSPROP	9.1E+03	PHYSPROP	5.3E-02	6.9E+00	1.7E+01	7.7E-03	EPI
Crotonaldehyde, trans-	123-73-9	7.0E+01	PHYSPROP	7.9E-04	1.9E-05	PHYSPROP	3.0E+01	PHYSPROP	-7.6E+01	PHYSPROP	8.5E-01	CRC	9.6E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			1.8E+00	EPI	6.0E-01	PHYSPROP	1.5E+05	PHYSPROP	5.1E-03	2.6E-01	6.2E-01	1.6E-03	EPI
Cumene	98-82-8	1.2E+02	PHYSPROP	4.7E-01	1.2E-02	PHYSPROP	4.5E+00	PHYSPROP	-9.6E+01	PHYSPROP	8.6E-01	CRC	6.0E-02	7.9E-06	WATER9 (U.S. EPA, 2001)			7.0E+02	EPI	3.7E+00	PHYSPROP	6.1E+01	PHYSPROP	3.8E-01	5.0E-01	1.2E+00	9.0E-02	EPI
Cuoferron	135-20-6	1.6E+02	PHYSPROP	1.5E-07	3.6E-09	PHYSPROP	6.3E-05	PHYSPROP	1.6E+02	PHYSPROP			6.6E-02	7.7E-06	WATER9 (U.S. EPA, 2001)			7.6E+02	EPI	-1.7E+00	PHYSPROP	6.1E+05	PHYSPROP	8.0E-06	7.8E-01	1.9E+00	1.7E-06	EPI
Cyanazine	21725-46-2	2.4E+02	PHYSPROP	1.1E-10	2.6E-12	EPI	1.4E-07	PHYSPROP	1.7E+02	PHYSPROP			4.9E-02	5.7E-06	WATER9 (U.S. EPA, 2001)			1.3E+02	EPI	2.2E+00	PHYSPROP	1.7E+02	PHYSPROP	1.2E-02	2.3E+00	5.6E+00	2.1E-03	EPI
Cyanides																												
~Calcium Cyanide	592-01-8	9.2E+01	PHYSPROP																					3.7E-03	3.4E-01	8.3E-01	1.0E-03	RAGSE
~Copper Cyanide	544-92-3	9.0E+01	PHYSPROP								4.7E+02	PHYSPROP	2.9E+00	CRC										3.6E-03	3.3E-01	8.0E-01	1.0E-03	RAGSE
~Cyanide (CN-)	57-12-5	2.6E+01	PHYSPROP	4.2E-03	1.0E-04	Ma et al 2010	3.1E+02	PHYSPROP			7.0E-01	CHEM GUIDE	2.1E-01	2.5E-05	WATER9 (U.S. EPA, 2001)	9.9E+00	SSL					9.5E+04	PHYSPROP	2.0E-03	1.5E-01	3.5E-01	1.0E-03	RAGSE
~Cyanogen	460-19-5	5.2E+01	PHYSPROP	2.2E-01	5.4E-03	EPI	4.3E+03	PHYSPROP	-2.8E+01	PHYSPROP	9.5E-01	CRC	1.2E-01	1.4E-05	WATER9 (U.S. EPA, 2001)					7.0E-02	PHYSPROP	8.0E+03	CRC	2.5E-03	2.1E-01	4.9E-01	8.9E-04	RAGSE
~Cyanogen Bromide	506-68-3	1.1E+02	PHYSPROP	1.0E+00	2.5E-02	EPI	1.2E+02	PHYSPROP	5.2E+01	PHYSPROP	2.0E+00	CRC	9.8E-02	1.4E-05	WATER9 (U.S. EPA, 2001)									1.0E-03	4.1E-01	9.9E-01	2.6E-04	RAGSE
~Cyanogen Chloride	506-77-4	6.1E+01	PHYSPROP	7.9E-02	1.9E-03	YAWS	1.2E+03	PHYSPROP	-6.6E+00	PHYSPROP	1.2E+00	CRC	1.2E-01	1.4E-05	WATER9 (U.S. EPA, 2001)							6.0E+04	PHYSPROP	1.2E-03	2.3E-01	5.6E-01	3.9E-04	RAGSE
~Hydrogen Cyanide	74-90-8	2.7E+01	PHYSPROP	5.4E-03	1.3E-04	PHYSPROP	7.4E+02	PHYSPROP	-1.3E+01	PHYSPROP	6.9E-01	CRC	1.7E-01	1.7E-05	WATER9 (U.S. EPA, 2001)	9.9E+00	SSL			-2.5E-01	PHYSPROP	1.0E+06	PHYSPROP	2.0E-03	1.5E-01	3.6E-01	1.0E-03	RAGSE
~Potassium Cyanide	151-50-8	6.5E+01	PHYSPROP				0.0E+00	NIOSH	6.3E+02	PHYSPROP	1.6E+00	CRC									

Contaminant		Molecular Weight		Volatility Parameters					Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients						Water Solubility		Tapwater Dermal Parameters			
Analyte	CAS No.	MW	MW Ref	H^+ (unitless)	HLC (atm-m^3/mole)	H^+ and HLC Ref	VP (mmHg)	VP Ref	MP (C)	MP Ref	Density (g/cm^3)	Density Ref	D_{ia} (cm^2/s)	D_{iw} (cm^2/s)	D_{ia} and D_{iw} Ref	K_d (L/kg)	K_d Ref	K_{oc} (L/kg)	K_{oc} Ref	log K_{ow} (unitless)	log K_{ow} Ref	S (mg/L)	S Ref	B (unitless)	t_{event} (hr/event)	t^* (hr)	K_p (cm/hr)	K_p Ref
Dichloro-2-butene, 1,4-	764-41-0	1.3E+02	PHYSPROP	3.5E-01	8.5E-03	PHYSPROP	3.0E+00	EPI	3.5E+00	PHYSPROP	1.2E+00	LANGE	6.7E-02	9.3E-06	WATER9 (U.S. EPA, 2001)					1.3E+02	EPI	2.6E+00	PHYSPROP	5.8E+02	PHYSPROP	7.1E-02	5.3E-01	1.3E+00	1.7E-02	EPI
Dichloro-2-butene, cis-1,4-	1476-11-5	1.3E+02	PHYSPROP	2.7E-02	6.6E-04	EPI	4.1E+00	PHYSPROP	-4.8E+01	PHYSPROP	1.2E+00	CRC	6.7E-02	9.3E-06	WATER9 (U.S. EPA, 2001)					1.3E+02	EPI	2.6E+00	PHYSPROP	5.8E+02	PHYSPROP	7.1E-02	5.3E-01	1.3E+00	1.7E-02	EPI
Dichloro-2-butene, trans-1,4-	110-57-6	1.3E+02	PHYSPROP	2.7E-02	6.6E-04	EPI	3.4E+00	PHYSPROP	2.0E+00	PHYSPROP	1.2E+00	CRC	6.6E-02	9.3E-06	WATER9 (U.S. EPA, 2001)					1.3E+02	EPI	2.6E+00	PHYSPROP	8.5E+02	PHYSPROP	7.1E-02	5.3E-01	1.3E+00	1.7E-02	EPI
Dichloroacetic Acid	79-43-6	1.3E+02	PHYSPROP	3.4E-07	8.4E-09	PHYSPROP	1.8E-01	PHYSPROP	1.4E+01	PHYSPROP	1.6E+00	CRC	7.2E-02	1.1E-05	WATER9 (U.S. EPA, 2001)					2.3E+00	EPI	9.2E-01	PHYSPROP	1.0E+06	PHYSPROP	5.3E-03	5.5E-01	1.3E+00	1.2E-03	EPI
Dichlorobenzene. 1,2-	95-50-1	1.5E+02	PHYSPROP	7.8E-02	1.9E-03	PHYSPROP	1.4E+00	PHYSPROP	-1.7E+01	PHYSPROP	1.3E+00	CRC	5.6E-02	8.9E-06	WATER9 (U.S. EPA, 2001)					3.8E+02	EPI	3.4E+00	PHYSPROP	1.6E+02	PHYSPROP	2.1E-01	7.0E-01	1.7E+00	4.5E-02	EPI
Dichlorobenzene, 1,4-	106-46-7	1.5E+02	PHYSPROP	9.9E-02	2.4E-03	PHYSPROP	1.7E+00	PHYSPROP	5.2E+01	PHYSPROP	1.2E+00	CRC	5.5E-02	8.7E-06	WATER9 (U.S. EPA, 2001)					3.8E+02	EPI	3.4E+00	PHYSPROP	8.1E+01	PHYSPROP	2.1E-01	7.0E-01	1.7E+00	4.5E-02	EPI
Dichlorobenzidine, 3,3'-	91-94-1	2.5E+02	PHYSPROP	1.2E-09	2.8E-11	PHYSPROP	2.6E-07	PHYSPROP	1.3E+02	PHYSPROP			4.7E-02	5.5E-06	WATER9 (U.S. EPA, 2001)					3.2E+03	EPI	3.5E+00	PHYSPROP	7.8E-02	2.8E+00	6.6E+00	1.3E-02	EPI
Dichlorobenzophenone, 4,4'-	90-98-2	2.5E+02	PHYSPROP	4.4E-05	1.1E-06	PHYSPROP	6.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.5E+00	CRC	2.6E-02	6.3E-06	WATER9 (U.S. EPA, 2001)					2.9E+03	EPI	4.4E+00	PHYSPROP	8.3E-01	PHYSPROP	3.3E-01	2.7E+00	6.4E+00	5.4E-02	EPI
Dichlorodifluoromethane	75-71-8	1.2E+02	PHYSPROP	1.4E+01	3.4E-01	PHYSPROP	4.8E+03	PHYSPROP	-1.6E+02	PHYSPROP	1.5E+00	PERRY	7.6E-02	1.1E-05	WATER9 (U.S. EPA, 2001)					4.4E+01	EPI	2.2E+00	PHYSPROP	2.8E+02	PHYSPROP	3.8E-02	5.0E-01	1.2E+00	9.0E-03	EPI
Dichloroethane, 1,1-	75-34-3	9.9E+01	PHYSPROP	2.3E-01	5.6E-03	PHYSPROP	2.3E+02	PHYSPROP	-9.7E+01	PHYSPROP	1.2E+00	CRC	8.4E-02	1.1E-05	WATER9 (U.S. EPA, 2001)					3.2E+01	EPI	1.8E+00	PHYSPROP	5.0E+03	PHYSPROP	2.6E-02	3.8E-01	9.0E-01	6.8E-03	EPI
Dichloroethane. 1,2-	107-06-2	9.9E+01	PHYSPROP	4.8E-02	1.2E-03	PHYSPROP	7.9E+01	PHYSPROP	-3.6E+01	PHYSPROP	1.2E+00	CRC	8.6E-02	1.1E-05	WATER9 (U.S. EPA, 2001)					4.0E+01	EPI	1.5E+00	PHYSPROP	8.6E+03	PHYSPROP	1.6E-02	3.8E-01	9.0E-01	4.2E-03	EPI
Dichloroethylene. 1,1-	75-35-4	9.7E+01	PHYSPROP	1.1E+00	2.6E-02	PHYSPROP	6.0E+02	PHYSPROP	-1.2E+02	PHYSPROP	1.2E+00	CRC	8.6E-02	1.1E-05	WATER9 (U.S. EPA, 2001)					3.2E+01	EPI	2.1E+00	PHYSPROP	2.4E+03	PHYSPROP	4.4E-02	3.7E-01	8.8E-01	1.2E-02	EPI
Dichloroethylene, cis-1,2-	156-59-2	9.7E+01	PHYSPROP	1.7E-01	4.1E-03	PHYSPROP	2.0E+02	PHYSPROP	-8.0E+01	PHYSPROP	1.3E+00	CRC	8.8E-02	1.1E-05	WATER9 (U.S. EPA, 2001)					4.0E+01	EPI	1.9E+00	PHYSPROP	6.4E+03	PHYSPROP	4.2E-02	3.7E-01	8.8E-01	1.1E-02	EPI
Dichloroethylene, trans-1,2-	156-60-5	9.7E+01	PHYSPROP	3.8E-01	9.4E-03	PHYSPROP	3.3E+02	EPI	-5.0E+01	PHYSPROP	1.3E+00	CRC	8.8E-02	1.1E-05	WATER9 (U.S. EPA, 2001)					4.0E+01	EPI	2.1E+00	PHYSPROP	4.5E+03	PHYSPROP	4.2E-02	3.7E-01	8.8E-01	1.1E-02	EPI
Dichlorophenol, 2,4-	120-83-2	1.6E+02	PHYSPROP	1.8E-04	4.3E-06	EPI	9.0E-02	PHYSPROP	4.5E+01	PHYSPROP	1.4E+00	PERRY	4.9E-02	8.7E-06	WATER9 (U.S. EPA, 2001)					1.5E+02	SSL	3.1E+00	PHYSPROP	5.6E+03	PHYSPROP	1.0E-01	8.6E-01	2.1E+00	2.1E-02	EPI
Dichlorophenoxy Acetic Acid, 2,4-	94-75-7	2.2E+02	PHYSPROP	1.4E-06	3.5E-08	EPI	8.3E-05	PHYSPROP	1.4E+02	PHYSPROP	1.4E+00	PubChem	2.8E-02	7.3E-06	WATER9 (U.S. EPA, 2001)					3.0E+01	EPI	2.8E+00	PHYSPROP	6.8E+02	PHYSPROP	3.8E-02	1.8E+00	4.4E+00	6.6E-03	EPI
Dichloropropane, 1,2-	78-87-5	1.1E+02	PHYSPROP	1.2E-01	2.8E-03	PHYSPROP	5.3E+01	PHYSPROP	-1.0E+02	PHYSPROP	1.2E+00	PERRY	7.3E-02	9.7E-06	WATER9 (U.S. EPA, 2001)					6.1E+01	EPI	2.0E+00	PHYSPROP	2.8E+03	PHYSPROP	3.1E-02	4.5E-01	1.1E+00	7.5E-03	EPI
Dichloropropene. 1,3-	142-28-9	1.1E+02	PHYSPROP	4.0E-02	9.8E-04	PHYSPROP	1.8E+01	PHYSPROP	-1.0E+02	PHYSPROP	1.2E+00	CRC	7.4E-02	9.8E-06	WATER9 (U.S. EPA, 2001)					7.2E+01	EPI	2.0E+00	PHYSPROP	2.8E+03	PHYSPROP	3.2E-02	4.5E-01	1.1E+00	7.8E-03	EPI
Dichloropropanol, 2,3-	616-23-9	1.3E+02	PHYSPROP	1.5E-07	3.6E-09	PHYSPROP	1.8E-01	PHYSPROP	-2.5E+01	EPI	1.4E+00	CRC	6.8E-02	9.9E-06	WATER9 (U.S. EPA, 2001)					5.6E+00	EPI	7.8E-01	PHYSPROP	6.4E+04	PHYSPROP	4.3E-03	5.5E-01	1.3E+00	9.8E-04	EPI
Dichloropropene, 1,3-	542-75-6	1.1E+02	PHYSPROP	1.5E-01	3.6E-03	PHYSPROP	3.4E+01	PHYSPROP	-3.5E+01	PHYSPROP	1.2E+00	LANGE	7.6E-02	1.0E-05	WATER9 (U.S. EPA, 2001)					7.2E+01	EPI	2.0E+00	PHYSPROP	2.8E+03	PHYSPROP	3.4E-02	4.4E-01	1.1E+00	8.3E-03	EPI
Dichlorvos	62-73-7	2.2E+02	PHYSPROP	2.3E-05	5.7E-07	EPI	1.6E-02	PHYSPROP	-6.0E+01	PHYSPROP	1.4E+00	CRC	2.8E-02	7.3E-06	WATER9 (U.S. EPA, 2001)					5.4E+01	EPI	1.4E+00	PHYSPROP	8.0E+03	PHYSPROP	4.6E-03	1.8E+00	4.4E+00	8.0E-04	EPI
Dicrotophos	141-66-2	2.4E+02	PHYSPROP	2.1E-09	5.0E-11	PHYSPROP	1.6E-04	PHYSPROP	7.9E+01	EPI	1.2E+00	CRC	2.5E-02	6.4E-06	WATER9 (U.S. EPA, 2001)					1.7E+01	EPI	0.0E+00	PHYSPROP	1.0E+06	PHYSPROP	4.3E-04	2.2E+00	5.4E+00	7.3E-05	EPI
Dicyclopentadiene	77-73-6	1.3E+02	PHYSPROP	2.6E+00	6.3E-02	PHYSPROP	2.3E+00	EPI	-1.0E+00	PHYSPROP	9.3E-01	LANGE	5.6E-02	7.8E-06	WATER9 (U.S. EPA, 2001)					1.5E+03	EPI	3.2E+00	PHYSPROP	2.6E+01	PHYSPROP	1.6E-01	5.8E-01	1.4E+00	3.6E-02	EPI
Dieldrin	60-57-1	3.8E+02	PHYSPROP	4.1E-04	1.0E-05	PHYSPROP	5.9E-06	PHYSPROP	1.8E+02	PHYSPROP	1.8E+00	CRC	2.3E-02	6.0E-06	WATER9 (U.S. EPA, 2001)					2.0E+04	EPI	5.4E+00	PHYSPROP	2.0E-01	PHYSPROP	2.4E-01	1.4E+01	3.4E+01	3.3E-02	EPI
Diesel Engine Exhaust	E17136615																											
Diethanolamine	111-42-2	1.1E+02	PHYSPROP	1.6E-09	3.9E-11	EPI	2.8E-04	PHYSPROP	2.8E+01	PHYSPROP	1.1E+00	CRC	7.7E-02	9.8E-06	WATER9 (U.S. EPA, 2001)					1.0E+00	EPI	-1.4E+00	PHYSPROP	1.0E+06	PHYSPROP	1.8E-04	4.1E-01	9.8E-01	4.5E-05	EPI
Diethylene Glycol Monobutyl Ether	112-34-5	1.6E+02	PHYSPROP	2.9E-07	7.2E-09	PHYSPROP	2.2E-02	PHYSPROP	-6.8E+01	PHYSPROP	9.6E-01	CRC	4.1E-02	7.0E-06	WATER9 (U.S. EPA, 2001)					1.0E+01	EPI	5.6E-01	PHYSPROP	1.0E+06	PHYSPROP	2.2E-03	8.5E-01	2.0E+00	4.5E-04	EPI
Diethylene Glycol Monoethyl Ether	111-90-0	1.3E+02	PHYSPROP	9.1E-07	2.2E-08	EPI	1.3E-01	PHYSPROP	-7.6E+01	PHYSPROP	9.9E-01	CRC	5.6E-02	8.0E-06	WATER9 (U.S. EPA, 2001)					1.0E+00	EPI	-5.4E-01	PHYSPROP	1.0E+06	PHYSPROP	5.4E-04	5.9E-01	1.4E+00	1.2E-04	EPI
Diethylformamide	617-84-5	1.0E+02	PHYSPROP	5.3E-06	1.3E-07	PHYSPROP	1.2E+00	EPI	-7.6E+00	EPI	9.1E-01	CRC	7.3E-02	9.0E-06	WATER9 (U.S. EPA, 2001)					2.1E+00	EPI	5.0E-02	PHYSPROP	1.0E+06	PHYSPROP	1.8E-03	3.9E-01	9.3E-01	4.6E-04	EPI
Diethylstilbestrol	56-53-1	2.7E+02	PHYSPROP	2.4E-10	5.8E-12	PHYSPROP	1.4E-08	PHYSPROP	1.7E+02	PHYSPROP			4.6E-02	5.3E-06	WATER9 (U.S. EPA, 2001)					2.7E+05	EPI	5.1E+00	PHYSPROP	1.2E+01	PHYSPROP	7.2E-01	3.3E+00	1.3E+01	1.1E-01	EPI
Difenzoquat	43222-48-6	3.6E+02	PHYSPROP				4.1E-12	PHYSPROP	1.6E+02	PHYSPROP			3.8E-02	4.4E-06	WATER9 (U.S. EPA, 2001)					7.8E+04	EPI	6.5E-01	PHYSPROP	8.2E+05	PHYSPROP	2.9E-04	1.1E+01	2.6E+01	4.0E-05	EPI
Diflubenzuron	35367-38-5	3.1E+02	PHYSPROP	1.9E-07	4.6E-09	EPI	9.0E-10	PHYSPROP	2.4E+02	PHYSPROP			4.1E-02	4.8E-06	WATER9 (U.S. EPA, 2001)					4.6E+02	EPI	3.9E+00	PHYSPROP	8.0E-02	PHYSPROP	7.3E-02	5.8E+00	1.4E+01	1.1E-02	EPI
Diluoroethane, 1,1-	75-37-6	6.6E+01	PHYSPROP	8.3E-01	2.0E-02	PHYSPROP	4.6E+03	PHYSPROP	-1.2E+02	PHYSPROP	9.0E-01	CRC	1.0E-01	1.1E-05	WATER9 (U.S. EPA, 2001)					2.2E+01	EPI	7.5E-01	PHYSPROP	2.7E+03	PHYSPROP	6.6E-03	2.5E-01	5.9E-01	1.2E-03	EPI
Diluoropropane, 2,2-	420-45-1	8.0E+01	PHYSPROP	2.1E+01	5.1E-01	PHYSPROP	1.8E+03	PHYSPROP	-1.0E+02	PHYSPROP	9.2E-01	CRC	9.0E-02	1.0E-05	WATER9 (U.S. EPA, 2001)					4.4E+01	EPI	2.3E+00	PHYSPROP	1.6E+02	PHYSPROP	6.4E-02	3.0E-01	7.1E-01	1.9E-02	EPI
Dihydrosafrole	94-58-6	1.6E+02	PHYSPROP	5.0E-04	1.2E-05	PHYSPROP	5.6E-02	PHYSPROP	4.4E+01	EPI	1.1E+00	PubChem	4.3E-02	7.4E-06	WATER9 (U.S. EPA, 2001)					2.1E+02	EPI	3.6E+00	PHYSPROP	5.7E+01	PHYSPROP	2.2E-01	8.7E-01	2.1E+00	4.5E-02	EPI
Diisopropyl Ether	108-20-3	1.0E+02	PHYSPROP	1.0E-01	2.6E-03	PHYSPROP	1.5E+02	PHYSPROP	-8.7E+01	PHYSPROP	7.2E-01	CRC	6.5E-02	7.8E-06	WATER9 (U.S. EPA, 2001)					2.3E+01	EPI	1.5E+00	PHYSPROP	8.8E+03	PHYSPROP	1.7E-02	3.9E-01	9.4E-01	4.3E-03	EPI
Diisopropyl Methylphosphonate	1445-75-6	1.8E+02	PHYSPROP	1.8E-03	4.4E-05	EPI	2.3E-01	PHYSPROP	-2.4E+01	EPI	9.8E-01	ATSDR Profile	3.4E-02	6.6E-06	WATER9 (U.S. EPA, 2001)					4.2E+01	EPI	1.0E+00	PHYSPROP	1.5E+03	PHYSPROP	3.8E-03	1.1E+00	2.6E+00	7.4E-04	EPI
Dimethipin	55290-64-7	2.1E+02	PHYSPROP	9.4E-10	2.3E-11	EPI	3.8E-07	PHYSPROP	1.7E+02	PHYSPROP			5.4E-02	6.3E-06	WATER9 (U.S. EPA, 2001)					1.0E+01	EPI	-1.7E-01	PHYSPROP	4.6E+03	PHYSPROP	4.5E-04	1.6E+00	3.8E+00	8.0E-05	EPI
Dimethoate	60-51-5	2.3E+02	PHYSPROP	9.9E-09	2.4E-10	EPI	1.9E-05	PHYSPROP	5.2E+01	PHYSPROP	1.3E+00	CRC	2.6E-02	6.7E-06	WATER9 (U.S. EPA, 2001)					1.3E+01	EPI	7.8E-01	PHYSPROP	2.3E+04	PHYSPROP	1.6E-03	2.0E+00	4.9E+00	2.7E-04	EPI
Dimethoxybenzidine, 3,3'-	119-90-4	2.4E+02	PHYSPROP	1.9E-09	4.7E-11	PHYSPROP	1.3E-07	PHYSPROP	1.4E+02	PHYSPROP			4.9E-02	5.7E-06	WATER9 (U.S. EPA, 2001)					5.1E+02	EPI	1.8E+01	PHYSPROP	6.4E+03	PHYSPROP	5.2E-02	5.9E+00	1.5E+01	1.1E-03	EPI
Dimethyl methylphosphonate	756-79-6	1.2E+02	PHYSPROP	5.6E-06	1.4E-07	PHYSPROP	8.3E-01	PHYSPROP	-4.8E+01	EPI	1.2E+00	CRC	6.7E-02	9.2E-06	WATER9 (U.S. EPA, 2001)					5.4E+00	EPI	-6.1E-01	PHYSPROP	1.0E+06	PHYSPROP	5.3E-04	5.2E-01	1.2E+00	1.2E-04	EPI
Dimethylamino azobenzene [p-]	60-11-7	2.3E+02	PHYSPROP	1.6E-08	4.0E-10	PHYSPROP	7.0E-08	EPI	1.2E+02	PHYSPROP			5.1E-02	6.0E-06	WATER9 (U.S. EPA, 2001)					2.0E+03	EPI	4.6E+00	PHYSPROP	2.3E-01	PHYSPROP	5.4E-01	1.9E+00	4.6E+00	9.4E-02	EPI
Dimethylaniline HCl, 2,4-	21436-96-4	1.2E+02	PHYSPROP	9.5E-05	2.3E-06	PHYSPROP	1.8E-01	PHYSPROP	1.6E+02	EPI			7.8E-02	9.1E-06	WATER9 (U.S. EPA, 2001)					3.5E+02	EPI	2.2E+00	PHYSPROP	3.7E+03	PHYSPROP	8.6E-05	5.0E-01	1.2E+00	2.0E-05	EPI
Dimethylaniline. 2,4-	95-68-1	1.2E+02	PHYSPROP	1.0E-04	2.5E-06	PHYSPROP	1.3E-01	PHYSPROP	-1.4E+01	PHYSPROP	9.7E-01	CRC	6.3E-02	8.4E-06	WATER9 (U.S. EPA, 2001)					1.8E+02	EPI	1.7E+00	PHYSPROP	6.1E+03	PHYSPROP	1.8E-02	5.0E-01	1.2E+00	4.3E-03	EPI
Dimethylaniline. N,N-	121-69-7	1.2E+02	PHYSPROP	2.3E-03	5.7E-05	EPI	7.0E-01	PHYSPROP	2.5E+00	PHYSPROP	9.6E-01	CRC	6.3E-02	8.3E-06	WATER9 (U.S. EPA, 2001)					7.9E+01	EPI	2.3E+00	PHYSPROP	1.5E+03	PHYSPROP	4.7E-02	5.0E-01	1.2E+00	1.1E-02	EPI
Dimethylbenzidine, 3,3'-	119-93-7	2.4E+02	PHYSPROP	2.6E-09	6.3E-11	PHYSPROP	6.9E-07	PHYSPROP	1.3E+02	PHYSPROP			5.3E-02	6.2E-06	WATER9 (U.S. EPA, 2001)					3.2E+03	EPI	2.3E+00	PHYSPROP	1.3E+03	PHYSPROP	2.0E-02	1.6E+00	3.6E+00	3.6E-03	EPI
Dimethylformamide	68-12-2	7.3E+01	PHYSPROP	3.0E-06	7.4E-08	PHYSPROP	3.9E+00	PHYSPROP	-6.0E+01	PHYSPROP	9.4E-01	CRC	9.7E-02	1.1E-05	WATER9 (U.S. EPA, 2001)					1.0E+00	EPI	-1.0E+00	PHYSPROP	1.0E+06	PHYSPROP	4.3E-04	2.7E-01	6.5E-01	1.3E-04	EPI
Dimethylhydrazine, 1,1-	57-14-7	6.0E+01	PHYSPROP	5.3E-04	1.3E-05	PHYSPROP	1.6E+02	PHYSPROP	-5.8E+01	PHYSPROP	7.9E-01	CRC	1.0E-01	1.1E-05	WATER9 (U.S. EPA, 2001)					1.2E+01	EPI	-1.2E+00	PHYSPROP	1.0E+06	PHYSPROP	3.5E-04	2.3E-01	5.5E-01	1.2E-04	EPI
Dimethylhydrazine, 1,2-	540-73-8	6.0E+01	PHYSPROP	2.8E-06	7.0E-08	PHYSPROP	7.0E+01	PHYSPROP	-9.0E+00	PHYSPROP	8.3E-01	CRC	1.1E-01	1.2E-05	WATER9 (U.S. EPA, 2001)					1.5E+01	EPI	-5.4E-01	PHYSPROP	1.0E+06	PHYSPROP	9.5E-04	2.3E-01	5.5E-01	3.2E-04	EPI
Dimethylphenol, 2,4-	105-67-9	1.2E+02	PHYSPROP	3.9E-05	9.5E-07	PHYSPROP	1.0E-01	PHYSPROP	2.5E+01	PHYSPROP	9.7E-01	CRC	6.2E-02	8.3E-06	WATER9 (U.S. EPA, 2001)					4.9E+02	EPI	2.3E+00	PHYSPROP	7.9E+03	PHYSPROP	4.6E-02	5.1E-01	1.2E+00	1.1E-02	EPI
Dimethylphenol. 2,6-	576-26-1	1.2E+02	PHYSPROP	2.7E-04	6.7E-06	PHYSPROP	1.7E-01	EPI	4.6E+01	PHYSPROP			7.7E-02	9.0E-06	WATER9 (U.S. EPA, 2001)					5.0E+02	EPI	2.4E+00	PHYSPROP	6.1E+03	PHYSPROP	5.1E-02	5.1E-01	1.2E+00	1.2E-02	EPI
Dimethylphenol, 3,4-	95-65-8	1.2E+02	PHYSPROP	1.7E-05	4.2E-07	PHYSPROP	3.6E-02	EPI	6.1E+01	PHYSPROP	9.8E-01	CRC	6.3E-02	8.4E-06	WATER9 (U.S. EPA, 2001)					4.9E+02	EPI	2.2E+00	PHYSPROP	4.8E+03	PHYSPROP	4.2E-02	5.1E-01	1.2E+00	9.8E-03	EPI
Dimethylvinylchloride	513-37-1	9.1E+01	PHYSPROP	4.8E-02	1.2E-03	CRC	2.1E+02	PHYSPROP	-1.0E+02	EPI	9.2E-01	CRC	8.1E-02	9.7E-06	WATER9 (U.S. EPA, 2001)					6.1E+01	EPI	2.6E+00	PHYSPROP	1.0E+03	PHYSPROP	9.3E-02	3.4E-01	8.1E-01	2.5E-02	EPI
Dinitro-o-cresol, 4,6-	534-52-1	2.0E+02	PHYSPROP	5.7E-05	1.4E-06	PHYSPROP	1.2E-04	PHYSPROP	8.7E+01	PHYSPROP			5.6E-02	6.5E-06	WATER9 (U.S. EPA, 2001)					7.5E+02	EPI	2.1E+00	PHYSPROP	2.0E+02	PHYSPROP	1.7E-02	1.4E+00	3.2E+00	3.2E-03	EPI
Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5	2.7E+02	PHYSPROP	2.3E-06	5.5E-08	PHYSPROP	4.2E-08	PHYSPROP	1.1E+02	PHYSPROP			4.6E-02	5.4E-06	WATER9 (U.S. EPA, 2001)					1.7E+04	EPI	4.1E+00	PHYSPROP	1.5E+01	PHYSPROP	1.7E-01	3.3E+00	7.8E+00	2.8E-02	EPI
Dinitroaniline, 3,5-	618-87-1	1.8E+02	PHYSPROP	1.2E-09	3.0E-11	PHYSPROP	2.7E-05	PHYSPROP	1.6E+02	PHYSPROP	1.6E+00	CRC	4.0E-02	8.8E-06	WATER9 (U.S. EPA, 2001)					1.7E+02	EPI	1.9E+00	PHYSPROP	1.3E+03	PHYSPROP	1.4E-02	1.1E+00	2.7E+00	2.7E-03	EPI
Dinitrobenzene. 1,2-	528-29-0	1.7E+02	PHYSPROP	2.2E-06	5.3E-08	EPI	4.6E-05	EPI	1.2E+02	PHYSPROP	1.3E+00	CRC	4.5E-02	8.3E-06	WATER9 (U.S. EPA, 2001)					3.6E+02	EPI	1.7E+00	PHYSPROP	1.3E+02	PHYSPROP	1.2E-02	9.2E-01	2.2E+00	2.4E-03	EPI
Dinitrobenzene, 1,3-	99-65-0	1.7E+02	PHYSPROP	2.0E-06	4.9E-08	PHYSPROP	9.0E-04	EPI	9.0E+01	PHYSPROP	1.6E+00	CRC	4.8E-02	9.2E-06	WATER9 (U.S. EPA, 2001)					3.5E+02	EPI	1.5E+00	PHYSPROP	5.3E+02	PHYSPROP	8.7E-03	9.2E-01	2.2E+00	1.7E-03	EPI
Dinitrobenzene, 1,4-	100-25-4	1.7E+02	PHYSPROP	3.4E-06	8.4E-08	PHYSPROP	2.6E-05	PHYSPROP	1.7E+02	PHYSPROP	1.6E+00	CRC	4.9E-02	9.4E-06	WATER9 (U.S. EPA, 2001)					3.5E+02	EPI	1.5E+00	PHYSPROP	6.9E+01	PHYSPROP	8.3E-03	9.2E-01	2.2E+00	1.7E-03	EPI
Dinitrophenol, 2,4-	51-28-5	1.8E+02	PHYSPROP	3.5E-06	8.6E-08	PHYSPROP	3.9E-04	PHYSPROP	1.1E+02	PHYSPROP	1.7E+00	CRC	4.1E-02	9.1E-06	WATER9 (U.S. EPA, 2001)					4.6E+02	EPI	1.7E+00	PHYSPROP	2.8E+03	PHYSPROP	9.8E-03	1.1E+00	2.7E+00	1.9E-03	EPI
Dinitrotoluene Mixture, 2,4/2,6-	E1615210	1.8E+02	EPI	1.6E-05	4.0E-07	EPI	2.2E-03	EPI	6.0E+01	EPI			5.9E-02	6.9E-06	WATER9 (U.S. EPA, 2001)					5.9E+02	EPI	2.2E+00	EPI	2.7E+02	EPI	2.2E-02	1.1E+00	2.6E+00	4.2E-03	EPI
Dinitrotoluene, 2,4-	121-14-2	1.8E+02	PHYSPROP	2.2E-06	5.4E-08	PHYSPROP	1.5E-04	PHYSPROP	7.1E+01	PHYSPROP	1.3E+00	CRC	3.8E-02	7.9E-06	WATER9 (U.S. EPA, 2001)					5.8E+02	EPI	2.0E+00	PHYSPROP	2.0E+02	PHYSPROP	1.6E-02	1.1E+00	2.6E+00	3.1E-03	EPI
Dinitrotoluene, 2,6-	606-20-2	1.8E+02	PHYSPROP	3.1E-05	7.5E-07	EPI	5.7E-04	PHYSPROP	6.6E+01	PHYSPROP	1.3E+00	CRC	3.7E-02	7.8E-06	WATER9 (U.S. EPA, 2001)					5.9E+02	EPI	2.1E+00	PHYSPROP	1.8E+02	PHYSPROP	1.9E-02	1.1E+00	2.6E+00	3.7E-03	EPI
Dinitrotoluene. 2-Amino-4,6-	35572-78-2	2.0E+02	PHYSPROP	1.3E-09	3.3E-11	PHYSPROP	1.1E-05	PHYSPROP	1.7E+02	PHYSPROP			5.6E-02	6.6E-06	WATER9 (U.S. EPA, 2001)					2.8E+02	EPI	1.8E+00	PHYSPROP	1.2E+03	PHYSPROP	1.1E-02	1.3E+00	3.2E+00	2.0E-03	EPI
Dinitrotoluene, 4-Amino-2,6-	19406-51-0	2.0E+02	PHYSPROP	1.3E-09	3.3E-11	PHYSPROP	1.1E-05	PHYSPROP	1.7E+02	PHYSPROP			5.6E-02	6.6E-06	WATER9 (U.S. EPA, 2001)					2.8E+02	EPI	1.8E+00	PHYSPROP	1.2E+03	PHYSPROP	1.1E-02	1.3E+00	3.2E+00	2.0E-03	EPI
Dinitrotoluene, Technical grade	25321-14-6	2.0E+02	PHYSPROP	2.8E-06	9.3E-08	PHYSPROP	4.0E-04	PHYSPROP	6.0E+01	EPI			5.9E-02	6.9E-06	WATER9 (U.S. EPA, 2001)					2.8E+02	EPI	1.9E+00	PHYSPROP	3.9E+02	PHYSPROP

Contaminant		Molecular Weight		Volatility Parameters					Melting Point		Density		Diffusivity in Air and Water			Partition Coefficients						Water Solubility		Tapwater Dermal Parameters				
Analyte	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m^3/mole)	H' and HLC Ref	VP (mmHg)	VP Ref	MP (C)	MP Ref	Density (g/cm^3)	Density Ref	D_{oa} (cm^2/s)	D_{iw} (cm^2/s)	D_{ia} and D_{iw} Ref	K_d (L/kg)	K_d Ref	K_{oc} (L/kg)	K_{oc} Ref	log K_{ow} (unitless)	log K_{ow} Ref	S (mg/L)	S Ref	B (unitless)	τ_{event} (hr/event)	t^* (hr)	K_p (cm/hr)	K_p Ref
Ethylene Thiourea	96-45-7	1.0E+02	PHYSPROP	5.6E-10	4.4E-11	PHYSPROP	2.0E-06	PHYSPROP	2.0E+02	PHYSPROP			8.7E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.3E+01	EPI	-6.6E-01	PHYSPROP	2.0E+04	PHYSPROP	5.9E-04	3.9E-01	9.4E-01	1.5E-04	EPI
Ethyleneimine	151-56-4	4.3E+01	PHYSPROP	4.9E-04	1.2E-05	EPI	2.1E+02	PHYSPROP	-7.8E+01	PHYSPROP	8.3E-01	CRC	1.3E-01	1.4E-05	WATER9 (U.S. EPA, 2001)			9.0E+00	EPI	-2.8E-01	PHYSPROP	1.0E+06	PHYSPROP	1.5E-03	1.8E-01	4.4E-01	5.8E-04	EPI
Ethylphthalyl Ethyl Glycolate	84-72-0	2.8E+02	PHYSPROP	2.7E-07	6.6E-09	PHYSPROP	2.2E-04	PHYSPROP	2.3E+01	EPI			4.4E-02	5.2E-06	WATER9 (U.S. EPA, 2001)			1.0E+03	EPI	2.2E+00	PHYSPROP	2.2E+02	PHYSPROP	7.7E-03	3.9E+00	9.4E+00	1.2E-03	EPI
Fenamiphos	22224-92-6	3.0E+02	PHYSPROP	4.9E-08	1.2E-09	EPI	1.0E-06	PHYSPROP	4.9E+01	PHYSPROP	1.2E+00	CRC	2.1E-02	5.4E-06	WATER9 (U.S. EPA, 2001)			4.0E+02	EPI	3.2E+00	PHYSPROP	3.3E+02	PHYSPROP	2.9E-02	5.3E+00	1.3E+01	4.4E-03	EPI
Fenoroxathrin	39515-41-8	3.5E+02	PHYSPROP	3.1E-04	7.6E-06	EPI	5.5E-06	PHYSPROP	4.7E+01	PHYSPROP			3.8E-02	4.5E-06	WATER9 (U.S. EPA, 2001)			2.2E+04	EPI	5.7E+00	PHYSPROP	3.3E-01	PHYSPROP	1.2E+00	9.5E+00	3.7E+01	1.7E-01	EPI
Fenvalerate	51630-58-1	4.2E+02	PHYSPROP	1.4E-06	3.5E-08	EPI	1.5E-09	PHYSPROP	4.0E+01	PHYSPROP	1.2E+00	CRC	1.8E-02	4.4E-06	WATER9 (U.S. EPA, 2001)			3.2E+05	EPI	6.2E+00	PHYSPROP	2.4E-02	PHYSPROP	7.4E-01	2.4E+01	9.1E+01	9.4E-02	EPI
Fluometuron	2164-17-2	2.3E+02	PHYSPROP	1.1E-07	2.6E-09	EPI	9.4E-07	PHYSPROP	1.6E+02	PHYSPROP			5.0E-02	5.9E-06	WATER9 (U.S. EPA, 2001)			2.9E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	1.9E-02	2.1E+00	5.0E+00	3.2E-03	EPI
Fluoride	16984-48-8	3.8E+01	EPI						-2.2E+02	EPI						1.5E+02	BAES					1.7E+00	EPI	2.4E-03	1.7E-01	4.1E-01	1.0E-03	RAGSE
Fluorine (Soluble Fluoride)	7782-41-4	3.8E+01	PHYSPROP						-2.2E+02	PHYSPROP	1.6E+00	CRC				1.5E+02	BAES					1.7E+00	PHYSPROP	2.4E-03	1.7E-01	4.1E-01	1.0E-03	RAGSE
Fluridone	59756-60-4	3.3E+02	PHYSPROP	3.3E-07	8.1E-09	EPI	9.8E-08	PHYSPROP	1.5E+02	PHYSPROP			4.0E-02	4.7E-06	WATER9 (U.S. EPA, 2001)			5.7E+04	EPI	3.2E+00	PHYSPROP	1.2E+01	PHYSPROP	2.0E-02	7.3E+00	1.8E+01	2.8E-03	EPI
Flurprimidol	56425-91-3	3.1E+02	PHYSPROP	5.4E-08	1.3E-09	EPI	3.6E-07	PHYSPROP	9.5E+01	PHYSPROP			4.1E-02	4.8E-06	WATER9 (U.S. EPA, 2001)			2.2E+03	EPI	3.3E+00	PHYSPROP	1.1E+02	PHYSPROP	3.1E-02	5.9E+00	1.4E+01	4.6E-03	EPI
Flusilazole	85509-19-9	3.2E+02	PHYSPROP	9.2E-08	2.3E-09	PHYSPROP	2.9E-07	PHYSPROP	5.4E+01	PHYSPROP			4.1E-02	4.8E-06	WATER9 (U.S. EPA, 2001)			8.1E+04	EPI	3.7E+00	PHYSPROP	5.4E+01	PHYSPROP	5.2E-02	6.1E+00	1.5E+01	7.7E-03	EPI
Flutolanil	66332-96-5	3.2E+02	PHYSPROP	1.3E-07	3.2E-09	EPI	4.9E-08	PHYSPROP	1.0E+02	PHYSPROP			4.0E-02	4.7E-06	WATER9 (U.S. EPA, 2001)			2.6E+03	EPI	3.7E+00	PHYSPROP	6.5E+00	PHYSPROP	4.8E-02	6.8E+00	1.6E+01	6.9E-03	EPI
Fluvalinate	69409-94-5	5.0E+02	PHYSPROP	5.9E-07	1.5E-08	PHYSPROP	1.0E-07	PHYSPROP	1.6E+02	EPI			3.0E-02	3.5E-06	WATER9 (U.S. EPA, 2001)			7.3E+05	EPI	6.8E+00	PHYSPROP	5.0E-03	PHYSPROP	6.8E-01	6.9E+01	2.7E+02	7.9E-02	EPI
Folpet	133-07-3	3.0E+02	PHYSPROP	3.1E-06	7.7E-08	EPI	1.6E-07	PHYSPROP	1.8E+02	EPI			4.3E-02	5.0E-06	WATER9 (U.S. EPA, 2001)			1.8E+01	EPI	2.9E+00	PHYSPROP	8.0E-01	PHYSPROP	1.8E-02	4.8E+00	1.2E+01	2.7E-03	EPI
Fomesafen	72178-02-0	4.4E+02	PHYSPROP	3.1E-11	7.5E-13	PHYSPROP	7.5E-07	EPI	2.2E+02	PHYSPROP	1.3E+00	CRC	1.9E-02	4.6E-06	WATER9 (U.S. EPA, 2001)			1.5E+03	EPI	2.9E+00	PHYSPROP	5.0E+01	PHYSPROP	3.7E-03	3.0E+01	7.2E+01	4.6E-04	EPI
Fonofos	944-22-9	2.5E+02	PHYSPROP	2.9E-04	7.0E-06	EPI	3.4E-04	PHYSPROP	6.6E-01	EPI	1.2E+00	CRC	2.4E-02	6.1E-06	WATER9 (U.S. EPA, 2001)			8.6E+02	EPI	3.9E+00	PHYSPROP	1.6E+01	PHYSPROP	1.6E-01	2.5E+00	6.0E+00	2.7E-02	EPI
Formaldehyde	50-00-0	3.0E+01	PHYSPROP	1.4E-05	3.4E-07	PHYSPROP	3.9E+03	EPI	-9.2E+01	PHYSPROP	8.2E-01	CRC	1.7E-01	1.7E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	3.5E-01	PHYSPROP	4.0E+05	PHYSPROP	3.8E-03	1.5E-01	3.7E-01	1.8E-03	EPI
Formic Acid	64-18-6	4.6E+01	PHYSPROP	6.8E-06	1.7E-07	PHYSPROP	4.3E+01	PHYSPROP	8.3E+00	PHYSPROP	1.2E+00	CRC	1.5E-01	1.7E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-5.4E-01	PHYSPROP	1.0E+06	PHYSPROP	9.9E-04	1.9E-01	4.6E-01	3.8E-04	EPI
Fosetyl-AL	39148-24-8	3.5E+02	PHYSPROP	1.3E-12	3.2E-14	PHYSPROP	7.5E-11	PHYSPROP	2.2E+02	PHYSPROP			3.8E-02	4.4E-06	WATER9 (U.S. EPA, 2001)			6.5E+03	EPI	-2.4E+00	PHYSPROP	1.1E+05	PHYSPROP	3.0E-06	1.0E+01	2.4E+01	4.1E-07	EPI
Furans																												
~Dibenzofuran	132-64-9	1.7E+02	PHYSPROP	8.7E-03	2.1E-04	EPI	2.5E-03	PHYSPROP	8.7E+01	PHYSPROP	1.1E+00	CRC	6.5E-02	7.4E-06	WATER9 (U.S. EPA, 2001)			9.2E+03	EPI	4.1E+00	PHYSPROP	3.1E+00	PHYSPROP	4.9E-01	9.2E-01	2.2E+00	9.8E-02	EPI
~Furan	110-00-9	6.8E+01	PHYSPROP	2.2E-01	5.4E-03	EPI	6.0E+02	PHYSPROP	-8.6E+01	PHYSPROP	9.5E-01	CRC	1.0E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			8.0E+01	EPI	1.3E+00	PHYSPROP	1.0E+04	PHYSPROP	1.6E-02	2.5E-01	6.1E-01	5.1E-03	EPI
~Tetrahydrofuran	109-99-9	7.2E+01	PHYSPROP	2.9E-03	7.1E-05	PHYSPROP	1.6E+02	PHYSPROP	-1.1E+02	PHYSPROP	8.8E-01	CRC	9.9E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			1.1E+01	EPI	4.6E-01	PHYSPROP	1.0E+06	PHYSPROP	4.1E-03	2.7E-01	6.4E-01	1.3E-03	EPI
Furazolidone	67-45-8	2.3E+02	PHYSPROP	1.3E-09	3.3E-11	PHYSPROP	2.6E-06	PHYSPROP	2.6E+02	PHYSPROP			5.1E-02	6.0E-06	WATER9 (U.S. EPA, 2001)			8.6E+02	EPI	-4.0E-02	PHYSPROP	4.0E+01	PHYSPROP	4.6E-04	1.9E+00	4.6E+00	8.0E-05	EPI
Furfural	98-01-1	9.6E+01	PHYSPROP	1.5E-04	3.8E-06	EPI	2.2E+00	PHYSPROP	-3.8E+01	PHYSPROP	1.2E+00	CRC	8.5E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			6.1E+00	EPI	4.1E-01	PHYSPROP	7.4E+04	PHYSPROP	3.2E-03	3.6E-01	8.7E-01	8.5E-04	EPI
Furium	531-82-8	2.6E+02	EPI	5.4E-14	1.3E-15	EPI	8.8E-09	EPI	1.9E+02	EPI			4.7E-02	5.5E-06	WATER9 (U.S. EPA, 2001)			5.8E+02	EPI	1.8E+00	EPI	4.2E+03	EPI	5.7E-03	2.8E+00	6.6E+00	9.4E-04	EPI
Furmecyclox	60568-05-0	2.5E+02	PHYSPROP	2.8E-07	6.9E-09	PHYSPROP	8.4E-05	PHYSPROP	3.3E+01	PHYSPROP			4.8E-02	5.6E-06	WATER9 (U.S. EPA, 2001)			4.3E+02	EPI	4.4E+00	PHYSPROP	3.0E-01	PHYSPROP	3.0E-01	2.7E+00	6.4E+00	5.0E-02	EPI
Glufosinate, Ammonium	77182-82-2	2.0E+02	PHYSPROP	1.8E-12	4.4E-14	PHYSPROP	9.1E-12	PHYSPROP	2.2E+02	PHYSPROP			5.6E-02	6.5E-06	WATER9 (U.S. EPA, 2001)			1.0E+01	EPI	-4.8E+00	PHYSPROP	1.4E+06	PHYSPROP	1.9E-07	1.4E+00	3.2E+00	3.4E-08	EPI
Glutaraldehyde	111-30-8	1.0E+02	PHYSPROP	1.3E-06	3.3E-08	PHYSPROP	6.0E-01	PHYSPROP	-3.0E+01	EPI			8.8E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-3.3E-01	PHYSPROP	2.2E+05	PHYSPROP	1.3E-03	3.8E-01	9.2E-01	3.3E-04	EPI
Glycidaldehyde	765-34-4	7.2E+01	PHYSPROP	2.1E-05	5.1E-07	PHYSPROP	4.5E+01	PHYSPROP	-6.2E+01	PHYSPROP	1.1E+00	CRC	1.1E-01	1.3E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-1.2E-01	PHYSPROP	1.0E+06	PHYSPROP	1.7E-03	2.7E-01	6.4E-01	5.2E-04	EPI
Glyphosate	1071-83-6	1.7E+02	PHYSPROP	8.6E-11	2.1E-12	EPI	9.8E-08	PHYSPROP	1.9E+02	PHYSPROP			6.2E-02	7.3E-06	WATER9 (U.S. EPA, 2001)			2.1E+03	USDA ARS	-3.4E+00	PHYSPROP	1.1E+04	PHYSPROP	2.3E-07	9.3E-01	2.2E+00	4.5E-08	EPI
Guanidine	113-00-8	5.9E+01	PHYSPROP	9.6E-10	2.3E-11	PHYSPROP	2.2E+00	PHYSPROP	5.0E+01	PHYSPROP	1.6E+00	GuideChem	1.4E-01	1.7E-05	WATER9 (U.S. EPA, 2001)			1.2E+01	EPI	-1.6E+00	PHYSPROP	1.8E+03	PHYSPROP	1.8E-04	2.3E-01	5.4E-01	6.0E-05	EPI
Guanidine Chloride	50-01-1	9.6E+01	PHYSPROP	8.9E-17	2.2E-18	PHYSPROP	1.8E-06	PHYSPROP	1.8E+02	PHYSPROP	1.4E+00	CRC	9.2E-02	1.2E-05	WATER9 (U.S. EPA, 2001)					-3.6E+00	PHYSPROP	1.0E+06	PHYSPROP	1.5E-07	3.6E-01	8.7E-01	3.9E-08	EPI
Guanidine Nitrate	506-93-4	1.2E+02	PHYSPROP	3.7E-17	9.0E-19	PHYSPROP	1.3E-07	PHYSPROP	2.1E+02	PHYSPROP			7.7E-02	9.0E-06	WATER9 (U.S. EPA, 2001)			2.3E+01	EPI	-8.4E+00	PHYSPROP	1.0E+06	PHYSPROP	1.1E-07	5.1E-01	1.2E+00	2.7E-08	EPI
Haloxyfop, Methyl	69806-40-2	3.8E+02	PHYSPROP	1.3E-05	3.2E-07	EPI	6.0E-06	PHYSPROP	5.6E+01	PHYSPROP			3.6E-02	4.3E-06	WATER9 (U.S. EPA, 2001)			5.5E+03	EPI	4.1E+00	PHYSPROP	9.3E+00	PHYSPROP	4.5E-02	1.3E+01	3.2E+01	6.0E-03	EPI
Heptachlor	76-44-8	3.7E+02	PHYSPROP	1.2E-02	2.9E-04	PHYSPROP	4.0E-04	PHYSPROP	9.6E+01	PHYSPROP	1.6E+00	CRC	2.2E-02	5.7E-06	WATER9 (U.S. EPA, 2001)			4.1E+04	EPI	6.1E+00	PHYSPROP	1.8E-01	PHYSPROP	1.1E+00	1.3E+01	5.0E+01	1.4E-01	EPI
Heptachlor Epoxide	1024-57-3	3.9E+02	PHYSPROP	8.6E-04	2.1E-05	PHYSPROP	2.0E-05	PHYSPROP	1.6E+02	PHYSPROP	1.9E+00	LookChem	2.4E-02	6.2E-06	WATER9 (U.S. EPA, 2001)			1.0E+04	EPI	5.0E+00	PHYSPROP	2.0E-01	PHYSPROP	1.6E-01	1.6E+01	3.8E+01	2.1E-02	EPI
Heptanal, n-	111-71-7	1.1E+02	PHYSPROP	1.1E-02	2.7E-04	PHYSPROP	3.5E+00	PHYSPROP	-4.3E+01	PHYSPROP	8.1E-01	CRC	6.2E-02	7.8E-06	WATER9 (U.S. EPA, 2001)			1.1E+01	EPI	2.3E+00	PHYSPROP	1.3E+03	PHYSPROP	4.9E-02	4.6E-01	1.1E+00	1.2E-02	EPI
Heptane, N-	142-82-5	1.0E+02	PHYSPROP	8.2E+01	2.0E+00	EPI	4.6E+01	PHYSPROP	-9.1E+01	PHYSPROP	6.8E-01	CRC	6.5E-02	7.6E-06	WATER9 (U.S. EPA, 2001)			2.4E+02	EPI	4.7E+00	PHYSPROP	2.9E+00	PHYSPROP	2.1E+00	3.8E-01	1.6E+00	5.4E-01	EPI
Hexabromobenzene	87-82-1	5.5E+02	PHYSPROP	1.1E-03	2.8E-05	PHYSPROP	1.6E-08	PHYSPROP	3.3E+02	PHYSPROP	3.0E+00	LookChem	2.5E-02	6.6E-06	WATER9 (U.S. EPA, 2001)			2.8E+03	EPI	6.1E+00	PHYSPROP	1.6E-04	PHYSPROP	1.2E-01	1.3E+02	3.1E+02	1.4E-02	EPI
Hexabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-153)	68631-49-2	6.4E+02	PubChem				5.8E-06	IRIS Profile					2.5E-02	3.0E-06	WATER9 (U.S. EPA, 2001)							9.0E-04	IRIS Profile		4.2E+02	1.0E+03		
Hexachlorobenzene	118-74-1	2.8E+02	PHYSPROP	7.0E-02	1.7E-03	PHYSPROP	1.8E-05	PHYSPROP	2.3E+02	PHYSPROP	2.0E+00	CRC	2.9E-02	7.8E-06	WATER9 (U.S. EPA, 2001)			6.2E+03	EPI	5.7E+00	PHYSPROP	6.2E-03	PHYSPROP	1.6E+00	4.1E+00	1.7E+01	2.5E-01	EPI
Hexachlorobutadiene	87-68-3	2.6E+02	PHYSPROP	4.2E-01	1.0E-02	PHYSPROP	2.2E-01	PHYSPROP	-2.1E+01	PHYSPROP	1.6E+00	CRC	2.7E-02	7.0E-06	WATER9 (U.S. EPA, 2001)			8.5E+02	EPI	4.8E+00	PHYSPROP	3.2E+00	PHYSPROP	5.0E-01	3.0E+00	7.3E+00	8.1E-02	EPI
Hexachlorocyclohexane, Alpha-	319-84-6	2.9E+02	PHYSPROP	2.7E-04	6.7E-06	PHYSPROP	3.5E-05	EPI	1.6E+02	PHYSPROP			4.3E-02	5.1E-06	WATER9 (U.S. EPA, 2001)			2.8E+03	EPI	3.8E+00	PHYSPROP	2.0E+00	PHYSPROP	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclohexane, Beta-	319-85-7	2.9E+02	PHYSPROP	1.8E-05	4.4E-07	PHYSPROP	3.6E-07	PHYSPROP	1.1E+02	PHYSPROP	1.9E+00	CRC	2.8E-02	7.4E-06	WATER9 (U.S. EPA, 2001)			2.8E+03	EPI	3.8E+00	PHYSPROP	2.4E-01	PHYSPROP	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	2.9E+02	PHYSPROP	2.1E-04	5.1E-06	PHYSPROP	4.2E-05	PHYSPROP	1.1E+02	PHYSPROP			4.3E-02	5.1E-06	WATER9 (U.S. EPA, 2001)			2.8E+03	EPI	3.7E+00	PHYSPROP	7.3E+00	PHYSPROP	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclohexane, Technical	608-73-1	2.9E+02	PHYSPROP	2.1E-04	5.1E-06	EPI	3.5E-05	EPI	1.1E+02	EPI			4.3E-02	5.1E-06	WATER9 (U.S. EPA, 2001)			2.8E+03	EPI	4.1E+00	EPI	8.0E+00	PHYSPROP	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclopentadiene	77-47-4	2.7E+02	PHYSPROP	1.1E+00	2.7E-02	PHYSPROP	6.0E-02	PHYSPROP	-9.0E+00	PHYSPROP	1.7E+00	CRC	2.7E-02	7.2E-06	WATER9 (U.S. EPA, 2001)			1.4E+03	EPI	5.0E+00	PHYSPROP	1.8E+00	PHYSPROP	6.5E-01	3.5E+00	1.4E+01	1.0E-01	EPI
Hexachloroethane	67-72-1	2.4E+02	PHYSPROP	1.6E-01	3.9E-03	PHYSPROP	2.1E-01	PHYSPROP	1.9E+02	PHYSPROP	2.1E+00	CRC	3.2E-02	8.9E-06	WATER9 (U.S. EPA, 2001)			2.0E+02	EPI	4.1E+00	PHYSPROP	5.0E+01	PHYSPROP	2.5E-01	2.2E+00	5.3E+00	4.2E-02	EPI
Hexachlorophene	70-30-4	4.1E+02	PHYSPROP	2.2E-11	5.5E-13	PHYSPROP	1.0E-10	PHYSPROP	1.7E+02	PHYSPROP			3.5E-02	4.0E-06	WATER9 (U.S. EPA, 2001)			6.7E+05	EPI	7.5E+00	PHYSPROP	1.4E+02	PHYSPROP	6.5E+00	2.0E+01	8.9E+01	8.4E-01	EPI
Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	2.2E+02	PHYSPROP	8.2E-10	2.0E-11	EPI	4.1E-09	EPI	2.1E+02	PHYSPROP	1.8E+00	CRC	3.1E-02	8.5E-06	WATER9 (U.S. EPA, 2001)			8.9E+01	EPI	8.7E-01	PHYSPROP	6.0E+01	PHYSPROP	1.9E-03	1.8E+00	4.4E+00	3.4E-04	EPI
Hexamethylene Diisocyanate, 1,6-	822-06-0	1.7E+02	PHYSPROP	2.0E-03	4.8E-05	PHYSPROP	3.0E-02	PHYSPROP	-6.7E+01	PHYSPROP	1.1E+00	CRC	4.0E-02	7.2E-06	WATER9 (U.S. EPA, 2001)			4.8E+03	EPI	3.2E+00	PHYSPROP	1.2E+02	PHYSPROP	1.2E-01	9.2E-01	2.2E+00	2.4E-02	EPI
Hexamethylene diisocyanate biuret	4035-89-6	4.8E+02	EPI	1.2E-14	3.0E-16	EPI	2.5E-15	EPI	2.9E+02	EPI			3.1E-02	3.6E-06	WATER9 (U.S. EPA, 2001)			4.4E+06	EPI	7.6E+00	EPI	6.9E-07	EPI	3.0E+00	5.0E+01	2.1E+02	3.5E-01	EPI
Hexamethylene diisocyanate isocyanurate	3779-63-3	5.0E+02	EPI	5.3E-16	1.3E-17	EPI	9.8E-17	EPI	3.1E+02	EPI			3.0E-02	3.5E-06	WATER9 (U.S. EPA, 2001)			1.3E+07	EPI	9.8E+00	EPI	5.0E-07	EPI	6.6E+01	7.0E+01	3.3E+02	7.7E+00	EPI
Hexamethylphosphoramide	680-31-9	1.8E+02	PHYSPROP	8.2E-07	2.0E-08	PHYSPROP	4.6E-02	PHYSPROP	7.2E+00	PHYSPROP	1.0E+00	CRC	3.5E-02	6.9E-06	WATER9 (U.S. EPA, 2001)			1.0E+01	EPI	2.8E-01	PHYSPROP	1.0E+06	PHYSPROP	1.2E-03	1.1E+00	2.5E+00	2.4E-04	EPI
Hexane, Commercial	E5241997	8.6E+01	PHYSPROP	7.4E+01	1.8E+00	PHYSPROP	1.5E+02	PHYSPROP	-9.5E+01	PHYSPROP	6.6E-01	CRC	7.3E-02	8.2E-06	WATER9 (U.S. EPA, 2001)			1.3E+02	EPI	3.9E+00	PHYSPROP	9.5E+00	PHYSPROP	7.2E-01	3.2E-01	1.2E+00	2.0E-01	EPI
Hexane, N-	110-54-3	8.6E+01	PHYSPROP	7.4E+01	1.8E+00	EPI	1.5E+02	PHYSPROP	-9.5E+01	PHYSPROP	6.6E-01	CRC	7.3E-02	8.2E-06	WATER9 (U.S. EPA, 2001)			1.3E+02	EPI	3.9E+00	PHYSPROP	9.5E+00	PHYSPROP	7.2E-01	3.2E-01	1.2E+00	2.0E-01	EPI
Hexanedioic Acid	124-04-9	1.5E+02	PHYSPROP	1.9E-10	4.7E-12	EPI	3.2E-07	EPI	1.5E+02	PHYSPROP	1.4E+00	CRC	5.8E-02	9.2E-06	WATER9 (U.S. EPA, 2001)			2.4E+01	EPI	8.0E-02	PHYSPROP	3.1E+04	PHYSPROP	1.2E-03	6.9E-01	1.7E+00	2.7E-04	EPI
Hexanol, 1-,2-ethyl- (2-Ethyl-1-hexanol)	104-76-7	1.3E+02	PHYSPROP	1.1E-03	2.7E-05	EPI	1.4E-01	PHYSPROP	-7.0E+01	PHYSPROP	8.3E-01	CRC	5.4E-02	7.3E-06	WATER9 (U.S. EPA, 2001)			3.5E+01	EPI	2.7E+00	PHYSPROP	8.8E+02	PHYSPROP	8.3E-02	5.6E-01	1.4E+00	1.9E-02	EPI
Hexanone, 2-	591-78-6	1.0E+02	PHYSPROP	3.8E-03	9.3E-05	EPI	1.2E+01	PHYSPROP	-5.6E+01	PHYSPROP	8.1E-01	CRC	7.0E-02	8.4E-06	WATER9 (U.S. EPA, 2001)			1.5E+01	EPI	1.4E+00	PHYSPROP	1.7E+04	PHYSPROP	1.4E-02	3.8E-01	9.2E-01	3.6E-03	EPI
Hexazinone	51235-04-2	2.5E+02	PHYSPROP	9.2E-11	2.3E-12	EPI	2.3E-07	EPI	1.2E+02	PHYSPROP	1.3E+00	CRC	2.5E-02	6.3E-06	WATER9 (U.S. EPA, 2001)			1.3E+02	EPI	1.9E+00	PHYSPROP	3.3E+04	PHYSPROP	6.2E-03	2.7E+00	6.5E+00	1.0E-03	EPI
Hexythiazox	78587-05-0	3.5E+02	PHYSPROP	9.7E-07	2.4E-08	EPI	2.6E-08	PHYSPROP	1.1E+02	PHYSPROP			3.8E-02	4.4E-06	WATER9 (U.S. EPA, 2001)			2.1E+03	EPI	5.6E+00	PHYSPROP	5.0E-01	PHYSPROP	6.0E-01	1.0E+01	2.4E+01	8.3E-02	EPI
Hydramethylnon	67485-29-4	4.9E+02	PHYSPROP	9.0E-05	2.2E-06	EPI	2.0E-08	PHYSPROP	1.9E+02	PHYSPROP			3.0E-02	3.6E-06	WATER9 (U.S. EPA, 2001)			1.8E+08	EPI	2.3E+00	PHYSPROP	6.0E-03	PHYSPROP	7.7E-04	6.2E+01	1.5E+02	9.0E-05	EPI
Hydrazine	302-01-2	3.2E+01	PHYSPROP	2.5E-05	6.1E-07	PHYSPROP	1.4E+01	PHYSPROP	2.0E+00	PHYSPROP	1.0E+00	CRC	1.7E-01	1.9E-05	WATER9 (U.S. EPA, 2001)			2.0E+00	PubChem	-2.1E+00	PHYSPROP	1.0E+06	PHYSPROP	9.5E-05	1.6E-01	3.8E-01	4.4E-05	RAGSE
Hydrazine Sulfate	10034-93-2	1.3E+02	EPI						2.5E+02	CRC	1.4E+00	CRC										3.1E+05	PERRY	4.4E-03	5.5E-01	1.3E+00	1.0E-03	RAGSE
Hydrogen Chloride	7647-01-0	3.5E+01	EPI	2.0E-08	4.9E-10	HSDB	3.5E+04	HSDB	-1.1E+02	CRC	1.5E+00	CRC	1.9E-01	2.3E-05	WATER9 (U.S. EPA, 2001)							6.7E+05	HSDB	2.3E-03	1.7E-01	4.0E-01	1.0E-03	RAGSE
Hydrogen Fluoride	7664-39-3	2.0E+01	PHYSPROP	4.3E-03	1.0E-04	PHYSPROP	9.2E+02	PHYSPROP	-8.4E+01	PHYSPROP	8.2E-01	CRC	2.2E-01	2.2E-05	WATER9 (U.S. EPA, 2001)					2.3E-01	PHYSPROP	1.0E+06	PHYSPROP	1.7E-03	1.4E-01	3.3E-01	1.0E-03	RAGSE
Hydrogen Sulfide	7783-06-4	3.4E+01	PHYSPROP	3.5E-01	8.6E-03	PHYSPROP	1.6E+04	PHYSPROP	-8.5E+01	PHYSPROP	1.4E+00	CRC	1.9E-01	2.2E-05	WATER9 (U.S. EPA, 2001)					2.3E-01	PhysProp	3.7E+03	PHYSPROP	2.2E-03	1.6E-01	3.9E-01	1.0E-03	RAGSE
Hydroquinone	123-31-9	1.1E+02	PHYSPROP	1.9E-09	4.7E-11	EPI	2.4E-05	EPI	1.7E+02	PHYSPROP	1.3E+00	CRC	8.0E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			2.4E+02	EPI	5.9E-01	PHYSPROP	7.2E+04	PHYSPROP	3.8E-03	4.3E-01	1.0E+00	9.3E-04

Contaminant		Molecular Weight		Volatility Parameters					Melting Point		Density		Diffusivity in Air and Water			Partition Coefficients					Water Solubility		Tapwater Dermal Parameters				
Analyte	CAS No.	MW	MW Ref	H^+ (unitless)	HLC (atm-m^3/mole)	H^+ and HLC Ref	VP (mmHg)	VP Ref	MP (C)	MP Ref	Density (g/cm^3)	Density Ref	D_{ia} (cm^2/s)	D_{iw} (cm^2/s)	D_{ia} and D_{iw} Ref	K_d (L/kg)	K_d Ref	K_{oc} (L/kg)	K_{oc} Ref	log K_{ow} (unitless)	log K_{ow} Ref	S (mg/L)	S Ref	B (unitless)	t_{event} (hr/event)	t^* (hr)	K_p (cm/hr)	K_p Ref
Manganese (Non-diet)	7439-96-5	5.5E+01	PHYSPROP				0.0E+00	NIOSH	1.2E+03	PHYSPROP	7.3E+00	CRC				6.5E+01	BAES							2.9E-03	2.1E-01	5.1E-01	1.0E-03	RAGSE
Mephosfolan	950-10-7	2.7E+02	PHYSPROP	4.9E-09	1.2E-10	PHYSPROP	3.2E-05	PHYSPROP	8.4E+01	EPI			4.6E-02	5.3E-06	WATER9 (U.S. EPA, 2001)			6.4E+02	EPI	1.0E+00	PHYSPROP	5.7E+01	PHYSPROP	1.5E-03	3.4E+00	8.1E+00	2.4E-04	EPI
Mepiquat Chloride	24307-26-4	1.5E+02	PHYSPROP	1.8E-10	4.3E-12	PHYSPROP	3.7E-07	PHYSPROP	2.2E+02	PHYSPROP			6.7E-02	7.9E-06	WATER9 (U.S. EPA, 2001)			6.6E+01	EPI	-2.8E+00	PHYSPROP	5.0E+05	PHYSPROP	1.4E-05	7.2E-01	1.7E+00	3.0E-06	EPI
Mercaptobenzothiazole, 2-	149-30-4	1.7E+02	PHYSPROP	1.5E-06	3.6E-08	PHYSPROP	4.6E-04	PHYSPROP	1.8E+02	PHYSPROP	1.4E+00	CRC	4.7E-02	8.7E-06	WATER9 (U.S. EPA, 2001)			1.4E+03	EPI	2.4E+00	PHYSPROP	1.2E+02	PHYSPROP	3.6E-02	9.1E-01	2.2E+00	7.3E-03	EPI
Mercury Compounds																												
~Mercuric Chloride (and other Mercury salts)	7487-94-7	2.7E+02	PHYSPROP						2.8E+02	PHYSPROP	5.6E+00	CRC								-2.2E-01	PHYSPROP	6.9E+04	PHYSPROP	6.3E-03	3.5E+00	8.4E+00	1.0E-03	RAGSE
~Mercury (elemental)	7439-97-6	2.0E+02	PHYSPROP	3.5E-01	8.6E-03	PHYSPROP VP/S	2.0E-03	PHYSPROP	-3.9E+01	PHYSPROP	1.4E+01	CRC	3.1E-02	6.3E-06	WATER9 (U.S. EPA, 2001)	5.2E+01	SSL			6.2E-01	PHYSPROP	6.0E-02	PHYSPROP	5.4E-03	1.4E+00	3.4E+00	1.0E-03	RAGSE
~Methyl Mercury	22967-92-6	2.2E+02	ChemID													7.0E+03	EPA							5.7E-03	1.7E+00	4.1E+00	1.0E-03	RAGSE
~Phenylmercuric Acetate	62-38-4	3.4E+02	PHYSPROP	2.3E-08	5.7E-10	EPI	6.0E-06	PHYSPROP	1.5E+02	PHYSPROP			3.9E-02	4.6E-06	WATER9 (U.S. EPA, 2001)			5.6E+01	EPI	7.1E-01	PHYSPROP	4.4E+03	PHYSPROP	4.2E-04	8.1E+00	1.9E+01	6.0E-05	EPI
Merphos	150-50-5	3.0E+02	PHYSPROP	9.3E-04	2.3E-05	PHYSPROP	2.0E-05	PHYSPROP	1.0E+02	PHYSPROP	1.0E+00	CRC	2.0E-02	5.0E-06	WATER9 (U.S. EPA, 2001)			4.9E+04	EPI	7.7E+00	PHYSPROP	3.5E-03	PHYSPROP	2.8E+01	4.9E+00	2.3E+01	4.2E+00	EPI
Metalaxyl	57837-19-1	2.8E+02	PHYSPROP	1.2E-07	3.0E-09	EPI	5.6E-06	PHYSPROP	7.1E+01	PHYSPROP			4.4E-02	5.2E-06	WATER9 (U.S. EPA, 2001)			3.9E+01	EPI	1.7E+00	PHYSPROP	8.4E+03	PHYSPROP	3.7E-03	3.9E+00	9.3E+00	5.8E-04	EPI
Methacrylonitrile	126-98-7	6.7E+01	PHYSPROP	1.0E-02	2.5E-04	EPI	7.1E+01	PHYSPROP	-3.6E+01	PHYSPROP	8.0E-01	CRC	9.6E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			1.3E+01	EPI	6.8E-01	PHYSPROP	2.5E+04	PHYSPROP	5.9E-03	2.5E-01	6.0E-01	1.9E-03	EPI
Methamidophos	10265-92-6	1.4E+02	PHYSPROP	3.5E-08	8.7E-10	PHYSPROP	3.5E-05	PHYSPROP	4.6E+01	PHYSPROP	1.3E+00	CRC	6.0E-02	9.2E-06	WATER9 (U.S. EPA, 2001)			5.4E+00	EPI	-8.0E-01	PHYSPROP	1.0E+06	PHYSPROP	3.4E-04	6.5E-01	1.6E+00	7.4E-05	EPI
Methanol	67-56-1	3.2E+01	PHYSPROP	1.9E-04	4.6E-06	PHYSPROP	1.3E+02	PHYSPROP	-9.8E+01	PHYSPROP	7.9E-01	CRC	1.6E-01	1.6E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-7.7E-01	PHYSPROP	1.0E+06	PHYSPROP	6.9E-04	1.6E-01	3.8E-01	3.2E-04	EPI
Methidathion	950-37-8	3.0E+02	PHYSPROP	2.9E-07	7.2E-09	EPI	3.4E-06	PHYSPROP	3.9E+01	PHYSPROP			4.2E-02	4.9E-06	WATER9 (U.S. EPA, 2001)			2.1E+01	EPI	2.2E+00	PHYSPROP	1.9E+02	PHYSPROP	6.1E-03	5.2E+00	1.2E+01	9.1E-04	EPI
Methomyl	16752-77-5	1.6E+02	PHYSPROP	8.1E-10	2.0E-11	EPI	5.4E-06	PHYSPROP	7.8E+01	PHYSPROP	1.3E+00	CRC	4.8E-02	8.4E-06	WATER9 (U.S. EPA, 2001)			1.0E+01	EPI	6.0E-01	PHYSPROP	5.8E+04	PHYSPROP	2.4E-03	8.5E-01	2.0E+00	4.8E-04	EPI
Methoxy-5-nitroaniline, 2-	99-59-2	1.7E+02	PHYSPROP	5.1E-07	1.3E-08	PHYSPROP	3.2E-04	PHYSPROP	1.2E+02	PHYSPROP	1.2E+00	CRC	4.3E-02	7.8E-06	WATER9 (U.S. EPA, 2001)			7.1E+01	EPI	1.5E+00	PHYSPROP	1.2E+02	PHYSPROP	8.4E-03	9.2E-01	2.2E+00	1.7E-03	EPI
Methoxychlor	72-43-5	3.5E+02	PHYSPROP	8.3E-06	2.0E-07	PHYSPROP	2.6E-06	PHYSPROP	8.7E+01	PHYSPROP	1.4E+00	CRC	2.2E-02	5.6E-06	WATER9 (U.S. EPA, 2001)			2.7E+04	EPI	5.1E+00	PHYSPROP	1.0E-01	PHYSPROP	3.1E-01	9.1E+00	2.2E+01	4.3E-02	EPI
Methoxyethanol Acetate, 2-	110-49-6	1.2E+02	PHYSPROP	1.3E-05	3.1E-07	EPI	7.0E+00	PHYSPROP	-6.5E+01	PHYSPROP	1.0E+00	CRC	6.6E-02	8.7E-06	WATER9 (U.S. EPA, 2001)			2.5E+00	EPI	1.0E-01	PHYSPROP	1.0E+06	PHYSPROP	1.7E-03	4.8E-01	1.2E+00	4.0E-04	EPI
Methoxyethanol, 2-	109-86-4	7.6E+01	PHYSPROP	1.3E-05	3.3E-07	PHYSPROP	9.5E+00	PHYSPROP	-9.5E+01	PHYSPROP	9.6E-01	CRC	9.5E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-7.7E-01	PHYSPROP	1.0E+06	PHYSPROP	6.0E-04	2.8E-01	6.7E-01	1.8E-04	EPI
Methyl Acetate	79-20-9	7.4E+01	PHYSPROP	4.7E-03	1.2E-04	PHYSPROP	2.2E+02	PHYSPROP	9.3E-01	CRC	9.6E-02	1.1E-05	WATER9 (U.S. EPA, 2001)					3.1E+00	EPI	1.8E-01	PHYSPROP	2.4E+05	PHYSPROP	2.6E-03	2.7E-01	6.6E-01	7.9E-04	EPI
Methyl Acrylate	96-33-3	8.6E+01	PHYSPROP	8.1E-03	2.0E-04	EPI	8.7E+01	PHYSPROP	-7.7E+01	PHYSPROP	9.5E-01	CRC	8.6E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			5.8E+00	EPI	8.0E-01	PHYSPROP	4.9E+04	PHYSPROP	6.2E-03	3.2E-01	7.7E-01	1.8E-03	EPI
Methyl Ethyl Ketone (2-Butanone)	78-93-3	7.2E+01	PHYSPROP	2.3E-03	5.7E-05	PHYSPROP	9.1E+01	PHYSPROP	-8.7E+01	PHYSPROP	8.0E-01	CRC	9.1E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			4.5E+00	EPI	2.9E-01	PHYSPROP	2.2E+05	PHYSPROP	3.1E-03	2.7E-01	6.4E-01	9.6E-04	EPI
Methyl Hydrazine	60-34-4	4.6E+01	PHYSPROP	1.2E-04	3.0E-06	PHYSPROP	5.0E+01	PHYSPROP	-5.2E+01	PHYSPROP	8.7E-01	LANGE	1.3E-01	1.4E-05	WATER9 (U.S. EPA, 2001)			1.3E+01	EPI	-1.1E+00	PHYSPROP	1.0E+06	PHYSPROP	4.5E-04	1.9E-01	4.6E-01	1.7E-04	EPI
Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1	1.0E+02	PHYSPROP	5.6E-03	1.4E-04	EPI	2.0E+01	PHYSPROP	-8.4E+01	PHYSPROP	8.0E-01	CRC	7.0E-02	8.3E-06	WATER9 (U.S. EPA, 2001)			1.3E+01	EPI	1.3E+00	PHYSPROP	1.9E+04	PHYSPROP	1.2E-02	3.8E-01	9.2E-01	3.2E-03	EPI
Methyl Isocyanate	624-83-9	5.7E+01	PHYSPROP	3.8E-02	9.3E-04	PHYSPROP	3.5E+02	PHYSPROP	-4.5E+01	PHYSPROP	9.6E-01	CRC	1.2E-01	1.3E-05	WATER9 (U.S. EPA, 2001)			4.0E+01	EPI	7.9E-01	PHYSPROP	2.9E+04	PHYSPROP	7.3E-03	2.2E-01	5.3E-01	2.5E-03	EPI
Methyl Methacrylate	80-62-6	1.0E+02	PHYSPROP	1.3E-02	3.2E-04	EPI	3.9E+01	PHYSPROP	-4.8E+01	PHYSPROP	9.4E-01	CRC	7.5E-02	9.2E-06	WATER9 (U.S. EPA, 2001)			9.1E+00	EPI	1.4E+00	PHYSPROP	1.5E+04	PHYSPROP	1.4E-02	3.8E-01	9.2E-01	3.6E-03	EPI
Methyl Parathion	298-00-0	2.6E+02	PHYSPROP	4.1E-06	1.0E-07	PHYSPROP	3.5E-06	PHYSPROP	3.6E+01	PHYSPROP	1.4E+00	CRC	2.5E-02	6.4E-06	WATER9 (U.S. EPA, 2001)			7.3E+02	EPI	2.9E+00	PHYSPROP	3.8E+01	PHYSPROP	2.6E-02	3.1E+00	7.5E+00	4.2E-03	EPI
Methyl Phosphonic Acid	993-13-5	9.6E+01	PHYSPROP	5.0E-10	1.2E-11	PHYSPROP	3.3E-04	EPI	1.1E+02	PHYSPROP			9.1E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			1.4E+00	EPI	-7.0E-01	PHYSPROP	2.0E+04	PHYSPROP	3.7E-04	3.6E-01	8.7E-01	9.8E-05	EPI
Methyl Styrene (Mixed Isomers)	25013-15-4	3.5E+02	PHYSPROP	1.1E-01	2.6E-03	PHYSPROP	1.5E+00	PHYSPROP	-8.6E+01	EPI	8.9E-01	HSDB	1.7E-02	4.2E-06	WATER9 (U.S. EPA, 2001)			7.2E+02	EPI	3.4E+00	PHYSPROP	8.9E+01	PHYSPROP	4.8E-01	1.0E+01	2.4E+01	6.6E-02	EPI
Methyl methanesulfonate	66-27-3	1.1E+02	PHYSPROP	1.6E-04	4.0E-06	PHYSPROP	3.1E-01	PHYSPROP	2.0E+01	PHYSPROP	1.3E+00	CRC	7.9E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			4.3E+00	EPI	-6.6E-01	PHYSPROP	2.0E+05	LANGE	5.6E-04	4.4E-01	1.0E+00	1.4E-04	EPI
Methyl tert-Butyl Ether (MTBE)	1634-04-4	8.8E+01	PHYSPROP	2.4E-02	5.9E-04	PHYSPROP	2.5E+02	PHYSPROP	-1.1E+02	PHYSPROP	7.4E-01	CRC	7.5E-02	8.6E-06	WATER9 (U.S. EPA, 2001)			1.2E+01	EPI	9.4E-01	PHYSPROP	5.1E+04	PHYSPROP	7.6E-03	3.3E-01	7.9E-01	2.1E-03	EPI
Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2	2.0E+02	PHYSPROP	2.6E-16	6.4E-18	PHYSPROP	4.1E-12	EPI	2.4E+02	EPI			5.6E-02	6.5E-06	WATER9 (U.S. EPA, 2001)			2.0E+02	EPI	-2.1E+00	PHYSPROP	1.0E+06	PHYSPROP	2.9E-05	1.3E+00	3.1E+00	5.4E-06	EPI
Methyl-2-Pentanol, 4-	108-11-2	1.0E+02	PHYSPROP	1.8E-03	4.5E-05	PHYSPROP	5.3E+00	PHYSPROP	-9.0E+01	PHYSPROP	8.1E-01	CRC	8.9E-02	9.3E-06	WATER9 (U.S. EPA, 2001)			8.2E+00	EPI	1.4E+00	PHYSPROP	1.6E+04	PHYSPROP	2.1E-02	3.9E-01	9.4E-01	5.4E-03	EPI
Methyl-5-Nitroaniline, 2-	99-55-8	1.5E+02	PHYSPROP	3.4E-07	8.3E-09	PHYSPROP	9.8E-04	PHYSPROP	1.1E+02	PHYSPROP			6.7E-02	7.8E-06	WATER9 (U.S. EPA, 2001)			1.8E+02	EPI	1.9E+00	PHYSPROP	1.0E+04	PHYSPROP	1.8E-02	7.5E-01	1.8E+00	3.8E-03	EPI
Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	1.5E+02	PHYSPROP	5.0E-11	1.2E-12	PHYSPROP	1.2E-04	PHYSPROP	1.2E+02	EPI			6.8E-02	8.0E-06	WATER9 (U.S. EPA, 2001)			7.2E+01	EPI	-9.2E-01	PHYSPROP	2.7E+05	PHYSPROP	2.7E-04	7.0E-01	1.7E+00	5.7E-05	EPI
Methylaniline Hydrochloride, 2-	636-21-5	1.4E+02	PHYSPROP	8.6E-05	2.1E-06	PHYSPROP	2.9E-01	PHYSPROP	2.2E+02	PHYSPROP			6.9E-02	8.1E-06	WATER9 (U.S. EPA, 2001)			1.2E+02	EPI	1.6E+00	PHYSPROP	8.3E+03	PHYSPROP	4.8E-05	6.7E-01	1.6E+00	1.1E-05	EPI
Methylarsonic acid	124-58-3	1.4E+02	PHYSPROP				1.6E-03	PHYSPROP	1.6E+02	PHYSPROP			7.0E-02	8.2E-06	WATER9 (U.S. EPA, 2001)			4.4E+01	EPI	-1.2E+00	PHYSPROP	2.6E+05	PHYSPROP	1.9E-04	6.4E-01	1.5E+00	4.2E-05	EPI
Methylbenzene,1-4-diamine monohydrochloride, 2-	74612-12-7	1.6E+02	PubChem										6.5E-02	7.6E-06	WATER9 (U.S. EPA, 2001)										8.1E-01	2.0E+00		
Methylbenzene-1,4-diamine sulfate, 2-	615-50-9	2.2E+02	ChemicalBook										5.2E-02	6.1E-06	WATER9 (U.S. EPA, 2001)										1.8E+00	4.3E+00		
Methylcholanthrene, 3-	56-49-5	2.7E+02	PHYSPROP	2.1E-04	5.2E-06	EPI	4.3E-08	EPI	1.8E+02	PHYSPROP	1.3E+00	CRC	2.4E-02	6.1E-06	WATER9 (U.S. EPA, 2001)			9.6E+05	EPI	6.4E+00	PHYSPROP	2.9E-03	PHYSPROP	5.7E+00	3.3E+00	1.5E+01	9.0E-01	EPI
Methylene Chloride	75-09-2	8.5E+01	PHYSPROP	1.3E-01	3.3E-03	PHYSPROP	4.4E+02	PHYSPROP	-9.5E+01	PHYSPROP	1.3E+00	CRC	1.0E-01	1.3E-05	WATER9 (U.S. EPA, 2001)			2.2E+01	EPI	1.3E+00	PHYSPROP	1.3E+04	PHYSPROP	1.3E-02	3.1E-01	7.5E-01	3.5E-03	EPI
Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	2.7E+02	PHYSPROP	1.7E-09	4.1E-11	PHYSPROP	2.9E-07	PHYSPROP	1.1E+02	PHYSPROP			4.6E-02	5.4E-06	WATER9 (U.S. EPA, 2001)			5.7E+03	EPI	3.9E+00	PHYSPROP	1.4E+01	PHYSPROP	1.2E-01	3.3E+00	7.9E+00	2.0E-02	EPI
Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	2.5E+02	PHYSPROP	4.4E-08	1.1E-09	PHYSPROP	1.8E-05	PHYSPROP	9.2E+01	PHYSPROP			4.7E-02	5.5E-06	WATER9 (U.S. EPA, 2001)			2.7E+03	EPI	4.4E+00	PHYSPROP	4.1E+00	PHYSPROP	2.9E-01	2.8E+00	6.7E+00	4.7E-02	EPI
Methylenebisbenzenamine, 4,4'-	101-77-9	2.0E+02	PHYSPROP	2.2E-09	5.3E-11	PHYSPROP	2.0E-07	PHYSPROP	9.3E+01	PHYSPROP			5.6E-02	6.5E-06	WATER9 (U.S. EPA, 2001)			2.1E+03	EPI	1.6E+00	PHYSPROP	1.0E+03	PHYSPROP	7.5E-03	1.4E+00	3.3E+00	1.4E-03	EPI
Methylenediphenyl Diisocyanate	101-68-8	2.5E+02	PHYSPROP	3.7E-05	9.0E-07	PHYSPROP	5.0E-06	PHYSPROP	3.8E+01	PHYSPROP	1.2E+00	CRC	2.4E-02	6.2E-06	WATER9 (U.S. EPA, 2001)			2.8E+05	EPI	5.2E+00	PHYSPROP	8.3E-01	PHYSPROP	1.1E+00	2.7E+00	1.0E+01	1.8E-01	EPI
Methylstyrene, Alpha-	98-83-9	1.2E+02	PHYSPROP	1.0E-01	2.6E-03	EPI	1.9E+00	EPI	-2.3E+01	PHYSPROP	9.1E-01	CRC	6.3E-02	8.2E-06	WATER9 (U.S. EPA, 2001)			7.0E+02	EPI	3.5E+00	PHYSPROP	1.2E+02	PHYSPROP	2.9E-01	4.8E-01	1.2E+00	7.0E-02	EPI
Metolachlor	51218-45-2	2.8E+02	PHYSPROP	3.7E-07	9.0E-09	PHYSPROP	3.1E-05	PHYSPROP	-6.2E+01	PHYSPROP	1.1E+00	CRC	2.2E-02	5.5E-06	WATER9 (U.S. EPA, 2001)			4.9E+02	EPI	3.1E+00	PHYSPROP	5.3E+02	PHYSPROP	2.2E-02	4.1E+00	9.8E+00	3.4E-03	EPI
Metribuzin	21087-64-9	2.1E+02	PHYSPROP	4.8E-09	1.2E-10	EPI	4.4E-07	PHYSPROP	1.3E+02	PHYSPROP	1.3E+00	CRC	2.7E-02	7.1E-06	WATER9 (U.S. EPA, 2001)			5.3E+01	EPI	1.7E+00	PHYSPROP	1.1E+03	PHYSPROP	7.4E-03	1.7E+00	4.0E+00	1.3E-03	EPI
Metsulfuron-methyl	74223-64-6	3.8E+02	PHYSPROP	5.4E-15	1.3E-16	EPI	2.5E-12	PHYSPROP	1.6E+02	PHYSPROP			3.6E-02	4.2E-06	WATER9 (U.S. EPA, 2001)			9.3E+01	EPI	2.2E+00	PHYSPROP	9.5E+03	PHYSPROP	2.5E-03	1.4E+01	3.4E+01	3.3E-04	EPI
Midrange Aliphatic Hydrocarbon Streams	E1790669	1.3E+02	PHYSPROP	1.4E+02	3.4E+00	EPI	4.5E+00	PHYSPROP	-5.4E+01	PHYSPROP	7.2E-01	CRC	5.1E-02	6.8E-06	WATER9 (U.S. EPA, 2001)			8.0E+02	EPI	5.7E+00	PHYSPROP	2.2E-01	PHYSPROP	7.4E+00	5.5E-01	2.5E+00	1.7E+00	EPI
Mineral oils	8012-95-1	1.7E+02	EPI	3.3E+02	8.2E+00	EPI	1.4E-01	EPI	-9.6E+00	EPI	8.8E-01	ChemNet	3.6E-02	6.4E-06	WATER9 (U.S. EPA, 2001)			4.8E+03	EPI	6.1E+00	EPI	3.7E-03	EPI	9.8E+00	9.5E-01	4.3E+00	2.0E+00	EPI
Mirex	2385-85-5	5.5E+02	PHYSPROP	3.3E-02	8.1E-04	PHYSPROP	8.0E-07	PHYSPROP	4.9E+02	CRC	2.3E+00	ChemNet	2.2E-02	5.5E-06	WATER9 (U.S. EPA, 2001)			3.6E+05	EPI	6.9E+00	PHYSPROP	8.5E-02	PHYSPROP	4.6E-01	1.2E+02	2.9E+02	5.2E-02	EPI
Molinate	2212-67-1	1.9E+02	PHYSPROP	1.7E-04	4.1E-06	PHYSPROP	5.6E-03	PHYSPROP	7.0E+01	EPI	1.1E+00	CRC	3.2E-02	6.8E-06	WATER9 (U.S. EPA, 2001)			1.8E+02	EPI	3.2E+00	PHYSPROP	9.7E+02	PHYSPROP	9.9E-02	1.2E+00	2.8E+00	1.9E-02	EPI
Molybdenum	7439-98-7	9.6E+01	PHYSPROP				0.0E+00	NIOSH	2.6E+03	PHYSPROP	1.0E+01	CRC				2.0E+01	BAES							3.8E-03	3.6E-01	8.7E-01	1.0E-03	RAGSE
Monochloramine	10599-90-3	5.1E+01	EPI						-6.6E+01	CRC														2.8E-03	2.0E-01	4.9E-01	1.0E-03	RAGSE
Monomethylaniline	100-61-8	1.1E+02	PHYSPROP	3.6E-04	8.9E-06	PHYSPROP	4.5E-01	PHYSPROP	-5.7E+01	PHYSPROP	9.9E-01	CRC	7.2E-02	9.1E-06	WATER9 (U.S. EPA, 2001)			8.2E+01	EPI	1.7E+00	PHYSPROP	5.6E+03	PHYSPROP	2.0E-02	4.2E-01	1.0E+00	5.0E-03	EPI
Myclobutanil	88671-89-0	2.7E+02	PHYSPROP	1.7E-07	4.3E-09	EPI	1.6E-06	PHYSPROP	6.6E+01	PHYSPROP			4.5E-02	5.3E-06	WATER9 (U.S. EPA, 2001)			6.1E+03	EPI	2.9E+00	PHYSPROP	1.4E+02	PHYSPROP	2.1E-02	3.6E+00	8.7E+00	3.4E-03	EPI
N,N-Diphenyl-1,4-benzenediamine	74-31-7	2.6E+02	PHYSPROP	8.4E-09	2.1E-10	PHYSPROP	6.4E-09	EPI	1.4E+02	PHYSPROP			4.7E-02	5.4E-06	WATER9 (U.S. EPA, 2001)			5.2E+04	EPI	4.0E+00	PHYSPROP	7.4E+00	PHYSPROP	1.6E-01	3.0E+00	7.2E+00	2.6E-02	EPI
Naled	300-76-5	3.8E+02	PHYSPROP	2.7E-03	6.5E-05	EPI	2.0E-04	PHYSPROP	2.7E+01	PHYSPROP	2.0E+00	CRC	2.5E-02	6.4E-06	WATER9 (U.S. EPA, 2001)			1.3E+02	EPI	1.4E+00	PHYSPROP	1.5E+00	PHYSPROP	7.1E-04	1.4E+01	3.4E+01	9.4E-05	EPI
Naphtha, High Flash Aromatic (HFAN)	64742-95-6			1.8E-02	4.4E-04	EPI	8.5E-02	EPI														3.1E+01	EPI					
Naphthylamine, 2-	91-59-8	1.4E+02	PHYSPROP	3.3E-06	8.1E-08	PHYSPROP	2.6E-04	PHYSPROP	1.1E+02	PHYSPROP	1.6E+00	CRC	6.4E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			2.5E+03	EPI	2.3E+00	PHYSPROP	1.9E+02	PHYSPROP	3.7E-02	6.7E-01	1.6E+00	8.1E-03	EPI
Napropamide	15299-99-7	2.7E+02	PHYSPROP	3.4E-08	8.4E-10	EPI	1.7E-07	PHYSPROP	7.5E+01	PHYSPROP			4.5E-02	5.3E-06	WATER9 (U.S. EPA, 2001)			3.2E+03	EPI	3.4E+00	PHYSPROP	7.3E+01	PHYSPROP	5.1E-02	3.5E+00	8.3E+00	8.0E-03	EPI
Nickel Acetate	373-02-4	1.8E+02	PHYSPROP				1.8E-05	PHYSPROP			1.8E+00	PERRY	4.6E-02	9.7E-06	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-1.4E+00	PHYSPROP	1.7E+05	PHYSPROP	9.9E-05	1.0E+00	2.5E+00	1.9E-05	EPI
Nickel Carbonate	3333-67-3	1.2E+02	PHYSPROP				3.6E-06	PHYSPROP					7.9E-02	9.2E-06	WATER9 (U.S. EPA, 2001)					-2.1E+00	PHYSPROP	9.3E+01	PERRY	5.5E-05	4.9E-01	1.2E+00	1.3E-05	EPI
Nickel Carbonyl	13463-39-3	1.7E+02	CRC	2.0E+01	5.0E-01	Matheson Gas MSDS	3.2E+02	NIOSH	-1.9E+01	CRC	1.3E+00	CRC	4.3E-02	8.2E-06	WATER9 (U.S. EPA, 2001)							1.8E+02	PERRY	9.5E-01	2.3E+00			
Nickel Hydroxide	12054-48-7	9.3E+01	WebBook								3.7E+03	CRC												3.7E-03	3.5E-01	8.3E-01	1.0E-03	RAGSE
Nickel Oxide	1313-99-1	7.5E+01	EPI						2.0E+03	CRC	6.7E+00	CRC												3.3E-03	2.8E-01	6.6E-01	1.0E-03	RAGSE
Nickel Refinery Dust	E715532															1.5E+02	BAES										2.0E-04	RAGSE
Nickel Soluble Salts	7440-02-0	5.9E+01	PHYSPROP				0.0E+00	NIOSH	1.5E+03	CRC	8.9E+00	CRC				6.5E+01	SSL							5.9E-04	2.2E-01	5.4E-01	2.0E-04	RAGSE
Nickel Subsulfide	12035-72-2	2.4E+02	CRC						7.9E+02	CRC	5.9E+00	CRC												1.2E-03	2.3E+00	5.6E+00	2.0E-04	RAGSE
Nickelocene	1271-28-9	1.9E+02	CRC						1.7E+02	CRC			5.8E-02	6.7E-06	WATER9 (U.S. EPA, 2001)										1.2E+00	2.9E+00		
Nitrate (measured as nitrogen)	14797-55-8	6.2E+01	EPI																					3.0E-03	2.3E-01	5.6E-01	1.0E-03	RAGSE
Nitrate + Nitrite (measured as nitrogen)	E701177																										1.0E-03	RAGSE
Nitrite (measured as nitrogen)	14797-65-0	4.7E+01	EPI																					2.6E-03	1.9E-01	4.6E-01	1.0E-03	RAGSE
Nitroaniline, 2-	88-74-4	1.4E+02	PHYSPROP	2.4E-06	5.9E-08	PHYSPROP	2.8E-03	PHYSPROP	7.1E+01	PHYSPROP	9.0E-01	CRC	5.2E-02	7.4E-06	WATER9 (U.S. EPA, 2001)			1.1E+02	EPI	1.9E+00	PHYSPROP	1.5E+03	PHYSPROP	2.0E-02	6.2E-01	1.5E+00	4.5E-03	EPI
Nitroaniline, 4-	100-01-6	1.4E+02	PHYSPROP	5.2E-08	1.3E-09	PHY

Contaminant		Molecular Weight		Volatility Parameters					Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients						Water Solubility		Tapwater Dermal Parameters				
Analyte	CAS No.	MW	MW Ref	H^+ (unitless)	HLC (atm-m^3/mole)	H^+ and HLC Ref	VP (mmHg)	VP Ref	MP (C)	MP Ref	Density (g/cm^3)	Density Ref	D_{oa} (cm^2/s)	D_{iw} (cm^2/s)	D_{ia} and D_{iw} Ref	K_d (L/kg)	K_d Ref	K_{oc} (L/kg)	K_{oc} Ref	log K_{ow} (unitless)	log K_{ow} Ref	S (mg/L)	S Ref	B (unitless)	τ_{event} (hr/event)	t^* (hr)	K_p (cm/hr)	K_p Ref
Paclobutrazol	76738-62-0	2.9E+02	PHYSPROP	3.4E-09	8.3E-11	EPI	7.5E-09	PHYSPROP	1.7E+02	PHYSPROP	1.2E+00	CRC	2.2E-02	5.7E-06	WATER9 (U.S. EPA, 2001)			9.2E+02	EPI	3.2E+00	PHYSPROP	2.6E+01	PHYSPROP	3.1E-02	4.6E+00	1.1E+01	4.7E-03	EPI
Paraquat Dichloride	1910-42-5	2.6E+02	PHYSPROP	1.3E-11	3.2E-13	PHYSPROP	7.5E-08	PHYSPROP	3.0E+02	EPI			4.7E-02	5.5E-06	WATER9 (U.S. EPA, 2001)			6.8E+03	EPI	-4.5E+00	PHYSPROP	6.2E+05	PHYSPROP	3.6E-07	2.9E+00	7.0E+00	5.8E-08	EPI
Parathion	56-38-2	2.9E+02	PHYSPROP	1.2E-05	3.0E-07	PHYSPROP	6.7E-06	PHYSPROP	6.1E+00	PHYSPROP	1.3E+00	CRC	2.3E-02	5.8E-06	WATER9 (U.S. EPA, 2001)			2.4E+03	EPI	3.8E+00	PHYSPROP	1.1E+01	PHYSPROP	8.4E-02	4.5E+00	1.1E+01	1.3E-02	EPI
Pebulate	1114-71-2	2.0E+02	PHYSPROP	9.7E-03	2.4E-04	EPI	8.9E-02	PHYSPROP	7.1E+01	EPI	9.5E-01	CRC	2.4E-02	6.1E-06	WATER9 (U.S. EPA, 2001)			3.0E+02	EPI	3.8E+00	PHYSPROP	1.0E+02	PHYSPROP	2.2E-01	1.4E+00	3.5E+00	4.0E-02	EPI
Pendimethalin	40487-42-1	2.8E+02	PHYSPROP	3.5E-05	8.6E-07	EPI	1.5E-05	PHYSPROP	5.6E+01	PHYSPROP	1.2E+00	CRC	2.3E-02	5.7E-06	WATER9 (U.S. EPA, 2001)			5.6E+03	EPI	5.2E+00	PHYSPROP	3.3E-01	PHYSPROP	7.4E-01	4.0E+00	1.5E+01	1.2E-01	EPI
Pentabromodiphenyl Ether	32534-81-9	5.6E+02	PHYSPROP	4.4E-03	1.1E-04	PHYSPROP	3.1E-08	EPI	-5.0E+00	PHYSPROP			2.8E-02	3.2E-06	WATER9 (U.S. EPA, 2001)			2.4E+03	EPI	6.8E+00	PHYSPROP	2.4E-03	PHYSPROP	3.4E-01	1.5E+02	3.7E+02	3.7E-02	EPI
Pentabromodiphenyl ether, 2,2',4,4',5- (BDE-99)	60348-60-9	5.6E+02	PHYSPROP	4.8E-05	1.2E-06	PHYSPROP	3.1E-08	EPI	-5.0E+00	EPI	2.3E+00	IRIS Profile	2.2E-02	5.6E-06	WATER9 (U.S. EPA, 2001)			2.2E+04	EPI	7.7E+00	PHYSPROP	3.4E-01	1.5E+02	3.7E+02	3.7E-02	EPI
Pentachlorobenzene	608-93-5	2.5E+02	PHYSPROP	2.9E-02	7.0E-04	PHYSPROP	1.0E-03	EPI	8.6E+01	PHYSPROP	1.8E+00	CRC	2.9E-02	7.9E-06	WATER9 (U.S. EPA, 2001)			3.7E+03	EPI	5.2E+00	PHYSPROP	8.3E-01	PHYSPROP	1.0E+00	2.7E+00	1.0E+01	1.7E-01	EPI
Pentachloroethane	76-01-7	2.0E+02	PHYSPROP	7.9E-02	1.9E-03	EPI	3.5E+00	PHYSPROP	-2.9E+01	PHYSPROP	1.7E+00	CRC	3.2E-02	8.6E-06	WATER9 (U.S. EPA, 2001)			1.4E+02	EPI	3.2E+00	PHYSPROP	4.9E+02	PHYSPROP	8.6E-02	1.4E+00	3.4E+00	1.6E-02	EPI
Pentachloronitrobenzene	82-68-8	3.0E+02	PHYSPROP	1.8E-03	4.4E-05	EPI	5.0E-05	PHYSPROP	1.4E+02	PHYSPROP	1.7E+00	CRC	2.6E-02	6.9E-06	WATER9 (U.S. EPA, 2001)			6.0E+03	EPI	4.6E+00	PHYSPROP	4.4E-01	PHYSPROP	2.8E-01	4.7E+00	1.1E+01	4.2E-02	EPI
Pentachlorophenol	87-86-5	2.7E+02	PHYSPROP	1.0E-06	2.5E-08	PHYSPROP	1.1E-04	PHYSPROP	1.7E+02	PHYSPROP	2.0E+00	CRC	3.0E-02	8.0E-06	WATER9 (U.S. EPA, 2001)			5.9E+02	SSL	5.1E+00	PHYSPROP	1.4E+01	PHYSPROP	8.0E-01	3.3E+00	1.3E+01	1.3E-01	EPI
Pentaerythritol tetranitrate (PETN)	78-11-5	3.2E+02	PHYSPROP	5.4E-08	1.3E-09	PHYSPROP	5.5E-09	EPI	1.4E+02	PHYSPROP	1.8E+00	CRC	2.6E-02	6.8E-06	WATER9 (U.S. EPA, 2001)			6.5E+02	EPI	2.4E+00	PHYSPROP	4.3E+01	PHYSPROP	6.9E-03	6.2E+00	1.5E+01	1.0E-03	EPI
Pentamethylphosphoramide (PMPA)	10159-46-3	1.7E+02	EPA Dashboard	3.0E-05	7.3E-07	EPA Dashboard	1.5E-01	EPA Dashboard	2.2E+01	EPI	1.1E+00	EPA Dashboard	4.2E-02	7.4E-06	WATER9 (U.S. EPA, 2001)			3.1E+00	EPI	-7.7E-01	EPI	1.8E+00	EPI	2.8E-04	8.8E-01	2.1E+00	5.7E-05	RAGSE
Pentane, n-	109-66-0	7.2E+01	PHYSPROP	5.1E+01	1.3E+00	PHYSPROP	5.1E+02	PHYSPROP	-1.3E+02	PHYSPROP	6.3E-01	CRC	8.2E-02	8.8E-06	WATER9 (U.S. EPA, 2001)			7.2E+01	EPI	3.4E+00	PHYSPROP	3.8E+01	PHYSPROP	3.6E-01	2.7E-01	6.4E-01	1.1E-01	EPI
Per- and Polyfluoroalkyl Substances (PFAS)																													
~Ammonium perfluoro-2-methyl-3-oxahexanoate	62037-80-3	3.5E+02	EPA OW										3.8E-02	4.5E-06	WATER9 (U.S. EPA, 2001)			1.2E+01	EPA OW			7.4E+05	EPA OW		9.2E+00	2.2E+01		
~Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6	3.3E+02	3M	1.0E-02	2.5E-04	EPA OW	2.3E+00	3M	-4.0E+01	EPA OW			4.0E-02	4.6E-06	WATER9 (U.S. EPA, 2001)							7.6E+05	3M		7.4E+00	1.8E+01		
~Perfluorobutanesulfonate	45187-15-3	3.0E+02	EPA SRS								1.8E+00	LookChem	2.7E-02	7.2E-06	WATER9 (U.S. EPA, 2001)			6.2E+01	Guelfo and Higgins			5.7E+04	Australian CHR		5.0E+00	1.2E+01		
~Perfluorobutanesulfonic acid (PFBS)	375-73-5	3.0E+02	3M				5.0E-02	3M	-2.1E+01	3M	1.8E+00	3M	2.7E-02	7.1E-06	WATER9 (U.S. EPA, 2001)			6.2E+01	Guelfo and Higgins	-3.4E-01	3M	2.6E+05	3M	1.3E-04	5.0E+00	1.2E+01	1.9E-05	RAGSE
~Perfluorohexanesulfonate	108427-53-8	4.0E+02	3M						4.1E+01	3M	1.8E+00	3M	2.3E-02	6.0E-06	WATER9 (U.S. EPA, 2001)			1.1E+02	Guelfo and Higgins	2.2E+00	3M			2.0E-03	1.8E+01	4.3E+01	2.6E-04	RAGSE
~Perfluorohexanesulfonic acid (PFHxS)	355-46-4	4.0E+02	3M						4.1E+01	3M	1.8E+00	3M	2.3E-02	6.0E-06	WATER9 (U.S. EPA, 2001)			1.1E+02	Guelfo and Higgins	2.2E+00	3M			2.0E-03	1.8E+01	4.4E+01	2.6E-04	RAGSE
~Perfluorononanoate	72007-68-2	4.6E+02	3M				9.5E-03	3M	6.8E+01	3M	1.8E+00	3M	2.1E-02	5.4E-06	WATER9 (U.S. EPA, 2001)	4.0E+00	3M	2.5E+02	Higgins and Luthy	2.6E+00	3M			1.6E-03	4.1E+01	9.9E+01	2.0E-04	RAGSE
~Perfluorononanoic acid (PFNA)	375-95-1	4.6E+02	3M				9.5E-03	3M	6.8E+01	3M	1.8E+00	3M	2.1E-02	5.4E-06	WATER9 (U.S. EPA, 2001)	4.0E+00	3M	2.5E+02	Higgins and Luthy	2.6E+00	3M			1.6E-03	4.2E+01	1.0E+02	2.0E-04	RAGSE
~Perfluorooctanesulfonate	45298-90-6	5.0E+02	EPA SRS				2.0E-03	HSDB			1.8E+00	LookChem	2.1E-02	5.3E-06	WATER9 (U.S. EPA, 2001)			3.7E+02	Higgins and Luthy			6.8E+02	OECD		6.6E+01	1.6E+02		
~Perfluorooctanesulfonic acid (PFOS)	1763-23-1	5.0E+02	3M	1.8E-05	4.4E-07	3M	6.5E-04	3M			1.8E+00	LookChem	2.1E-02	5.3E-06	WATER9 (U.S. EPA, 2001)			3.7E+02	Higgins and Luthy	-1.1E+00	3M	6.8E+02	3M	4.0E-06	6.6E+01	1.6E+02	4.7E-07	RAGSE
~Perfluorooctanoate	45285-51-6	4.1E+02	3M	1.5E-04	3.6E-06	ATSDR Profile	3.0E-02	3M	5.3E+01	3M	1.8E+00	3M	2.3E-02	5.8E-06	WATER9 (U.S. EPA, 2001)	1.5E+01	3M	1.1E+02	Higgins and Luthy			9.5E+03	3M		2.2E+01	5.2E+01		
~Perfluorooctanoic acid (PFOA)	335-67-1	4.1E+02	3M	1.5E-04	3.6E-06	ATSDR Profile	3.0E-02	3M	5.3E+01	3M	1.8E+00	3M	2.3E-02	5.8E-06	WATER9 (U.S. EPA, 2001)	1.5E+01	3M	1.1E+02	Higgins and Luthy			9.5E+03	3M		2.2E+01	5.3E+01		
~Potassium perfluorobutanesulfonate	29420-49-3	3.4E+02	3M	3.6E-11	8.8E-13	3M	9.2E-08	3M	2.8E+02	3M	9.5E-01	3M	1.8E-02	4.5E-06	WATER9 (U.S. EPA, 2001)	3.0E-01	3M			-1.8E+00	3M	4.6E+04	3M	8.9E-06	8.2E+00	2.0E+01	1.3E-06	RAGSE
~Potassium perfluorooctanesulfonate	2795-39-3	5.4E+02	Sax's	8.2E-05	2.0E-06	Beach (2005)	2.5E-06	3M	4.0E+02	3M			2.9E-02	3.4E-06	WATER9 (U.S. EPA, 2001)					-1.1E+00	3M	6.8E+02	3M	2.6E-06	1.1E+02	2.6E+02	2.9E-07	RAGSE
Perchlorates																													
~Ammonium Perchlorate	7790-98-9	1.2E+02	PHYSPROP								2.0E+00	CRC										2.5E+05	PHYSPROP	4.2E-03	4.8E-01	1.1E+00	1.0E-03	RAGSE
~Lithium Perchlorate	7791-03-9	1.1E+02	CRC						2.4E+02	CRC	2.4E+00	CRC										5.9E+05	CRC	4.0E-03	4.1E-01	1.0E+00	1.0E-03	RAGSE
~Perchlorate and Perchlorate Salts	14797-73-0	1.2E+02	CRC																			2.5E+05	CRC	4.2E-03	4.8E-01	1.1E+00	1.0E-03	RAGSE
~Potassium Perchlorate	7778-74-7	1.4E+02	PHYSPROP						5.3E+02	PHYSPROP	2.5E+00	CRC										1.5E+04	PHYSPROP	9.1E-03	6.3E-01	1.5E+00	2.0E-03	RAGSE
~Sodium Perchlorate	7601-89-0	1.2E+02	PHYSPROP						4.8E+02	EPI	2.5E+00	CRC										2.1E+06	PHYSPROP	4.3E-03	5.1E-01	1.2E+00	1.0E-03	RAGSE
Permethrin	52645-53-1	3.9E+02	PHYSPROP	7.6E-05	1.9E-06	EPI	2.2E-08	PHYSPROP	3.4E+01	PHYSPROP	1.2E+00	CRC	1.9E-02	4.8E-06	WATER9 (U.S. EPA, 2001)			1.2E+05	EPI	6.5E+00	PHYSPROP	6.0E-03	PHYSPROP	1.6E+00	1.6E+01	6.5E+01	2.1E-01	EPI
Phenacetin	62-44-2	1.8E+02	PHYSPROP	8.7E-09	2.1E-10	EPI	6.9E-07	PHYSPROP	1.4E+02	PHYSPROP			6.0E-02	7.0E-06	WATER9 (U.S. EPA, 2001)			4.1E+01	EPI	1.6E+00	PHYSPROP	7.7E+02	PHYSPROP	8.9E-03	1.1E+00	2.5E+00	1.7E-03	EPI
Phenmedipham	13684-63-4	3.0E+02	PHYSPROP	3.4E-11	8.4E-13	EPI	1.0E-11	PHYSPROP	1.4E+02	PHYSPROP			4.2E-02	5.0E-06	WATER9 (U.S. EPA, 2001)			2.6E+03	EPI	3.6E+00	PHYSPROP	4.7E+00	PHYSPROP	5.2E-02	5.1E+00	1.2E+01	7.9E-03	EPI
Phenol	108-95-2	9.4E+01	PHYSPROP	1.4E-05	3.3E-07	PHYSPROP	3.5E-01	PHYSPROP	4.1E+01	PHYSPROP	1.1E+00	CRC	8.3E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.9E+02	EPI	1.5E+00	PHYSPROP	8.3E+04	PHYSPROP	1.6E-02	3.5E-01	8.5E-01	4.3E-03	EPI
Phenol, 2,2-(1-methylethoxy)-, methylcarbamate	114-26-1	2.1E+02	PHYSPROP	5.8E-08	1.4E-09	EPI	2.1E-05	PHYSPROP	9.0E+01	PHYSPROP	1.1E+00	CRC	2.0E-02	6.6E-06	WATER9 (U.S. EPA, 2001)			6.0E+01	EPI	1.5E+00	PHYSPROP	1.9E+05	PHYSPROP	6.0E-03	1.6E+01	3.7E+01	1.1E-03	EPI
Phenothiazine	92-84-2	2.0E+02	PHYSPROP	1.1E-06	2.8E-08	PHYSPROP	8.9E-07	PHYSPROP	1.9E+02	PHYSPROP	1.3E+00	PubChem	2.9E-02	7.5E-06	WATER9 (U.S. EPA, 2001)			1.5E+03	EPI	4.2E+00	PHYSPROP	1.6E+00	PHYSPROP	3.7E-01	1.4E+00	3.3E+00	6.8E-02	EPI
Phenyl Isothiocyanate	103-72-0	1.4E+02	PHYSPROP	1.2E-01	3.0E-03	EPI	1.5E+00	PHYSPROP	-2.1E+01	PHYSPROP	1.1E+00	CRC	5.9E-02	8.6E-06	WATER9 (U.S. EPA, 2001)			2.2E+02	EPI	3.3E+00	PHYSPROP	9.0E+01	PHYSPROP	1.8E-01	6.0E-01	1.4E+00	4.1E-02	EPI
Phenylenediamine, m-	108-45-2	1.1E+02	PHYSPROP	5.1E-08	1.3E-09	EPI	2.1E-03	EPI	6.4E+01	PHYSPROP	1.0E+00	CRC	7.2E-02	9.2E-06	WATER9 (U.S. EPA, 2001)			3.4E+01	EPI	-3.3E-01	PHYSPROP	2.4E+05	PHYSPROP	9.4E-04	4.2E-01	1.0E+00	2.3E-04	EPI
Phenylenediamine, o-	95-54-5	1.1E+02	PHYSPROP	2.9E-07	7.2E-09	EPI	2.1E-03	EPI	1.0E+02	PHYSPROP			8.4E-02	9.8E-06	WATER9 (U.S. EPA, 2001)			3.5E+01	EPI	1.5E-01	PHYSPROP	4.0E+04	PHYSPROP	1.9E-03	4.2E-01	1.0E+00	4.9E-04	EPI
Phenylenediamine, p-	106-50-3	1.1E+02	PHYSPROP	2.8E-08	6.7E-10	PHYSPROP	5.0E-03	PHYSPROP	1.5E+02	PHYSPROP			8.4E-02	9.8E-06	WATER9 (U.S. EPA, 2001)			3.4E+01	EPI	-3.0E-01	PHYSPROP	3.7E+04	PHYSPROP	9.8E-04	4.2E-01	1.0E+00	2.5E-04	EPI
Phenylphenol, 2-	90-43-7	1.7E+02	PHYSPROP	4.3E-05	1.1E-06	EPI	2.0E-03	EPI	5.9E+01	PHYSPROP	1.2E+00	CRC	4.2E-02	7.8E-06	WATER9 (U.S. EPA, 2001)			6.7E+03	EPI	3.1E+00	PHYSPROP	7.0E+02	PHYSPROP	9.8E-02	9.4E-01	2.3E+00	2.0E-02	EPI
Phorate	298-02-2	2.6E+02	PHYSPROP	1.8E-04	4.4E-06	EPI	6.4E-04	PHYSPROP	-1.5E+01	CRC	1.2E+00	CRC	2.3E-02	5.9E-06	WATER9 (U.S. EPA, 2001)			4.6E+02	EPI	3.6E+00	PHYSPROP	5.0E+01	PHYSPROP	7.8E-02	3.0E+00	7.2E+00	1.3E-02	EPI
Phosgene	75-44-5	9.9E+01	PHYSPROP	6.8E-01	1.7E-02	PHYSPROP	1.4E+03	PHYSPROP	-1.2E+02	PHYSPROP	1.4E+00	CRC	8.9E-02	1.2E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-7.1E-01	PHYSPROP	6.8E+03	YAWS	5.6E-04	3.8E-01	9.0E-01	1.5E-04	EPI
Phosmet	732-11-6	3.2E+02	PHYSPROP	3.4E-07	8.4E-09	EPI	4.9E-07	PHYSPROP	7.2E+01	PHYSPROP			4.1E-02	4.8E-06	WATER9 (U.S. EPA, 2001)			1.0E+01	EPI	2.8E+00	PHYSPROP	2.4E+01	PHYSPROP	1.3E-02	6.3E+00	1.5E+01	1.8E-03	EPI
Phosphine	7803-51-2	3.4E+01	PHYSPROP	1.0E+00	2.4E-02	PHYSPROP	2.9E+04	PHYSPROP	-1.3E+02	PHYSPROP	1.4E+00	CRC	1.9E-01	2.2E-05	WATER9 (U.S. EPA, 2001)					-2.7E-01	PHYSPROP	2.6E+05	PERRY	2.2E-03	1.6E-01	3.9E-01	1.0E-03	RAGSE
Phosphoric Acid	7664-38-2	9.8E+01	PHYSPROP				3.0E-02	NIOSH	4.2E+01	PHYSPROP	1.8E+00	PERRY										5.5E+06	CRC	3.8E-03	3.7E-01	8.9E-01	1.0E-03	RAGSE
Phosphorus, White	7723-14-0	3.1E+01	YAWS	8.6E-02	2.1E-03	ATSDR Profile	2.5E-02	ATSDR Profile	4.4E+01	ATSDR Profile	1.8E+00	ATSDR Profile	2.2E-01	2.8E-05	WATER9 (U.S. EPA, 2001)	3.5E+00	BAES	1.1E+03	ATSDR Profile	3.1E+00	ATSDR Profile	3.0E+00	ATSDR Profile	2.1E-03	1.6E-01	3.8E-01	1.0E-03	RAGSE
Phthalates																													
~Bis(2-ethylhexyl)phthalate	117-81-7	3.9E+02	PHYSPROP	1.1E-05	2.7E-07	EPI	1.4E-07	PHYSPROP	-5.5E+01	PHYSPROP	9.8E-01	CRC	1.7E-02	4.2E-06	WATER9 (U.S. EPA, 2001)			1.2E+05	EPI	7.6E+00	PHYSPROP	2.7E-01	PHYSPROP	8.6E+00	1.6E+01	7.3E+01	1.1E+00	EPI
~Butyl Benzyl Phthalate	85-68-7	3.1E+02	PHYSPROP	5.2E-05	1.3E-06	EPI	8.3E-06	PHYSPROP	6.1E+01	EPI	1.1E+00	CRC	2.1E-02	5.2E-06	WATER9 (U.S. EPA, 2001)			7.2E+03	EPI	4.7E+00	PHYSPROP	2.7E+00	PHYSPROP	2.6E-01	5.9E+00	1.4E+01	3.9E-02	EPI
~Butylphthalyl Butylglycolate	85-70-1	3.4E+02	PHYSPROP	8.4E-07	2.1E-08	PHYSPROP	7.1E-06	PHYSPROP	-3.5E+01	PHYSPROP	1.1E+00	LANGE	2.0E-02	4.9E-06	WATER9 (U.S. EPA, 2001)			1.1E+04	EPI	4.2E+00	PHYSPROP	8.8E+00	PHYSPROP	8.2E-02	8.0E+00	1.9E+01	1.2E-02	EPI
~Dibutyl Phthalate	84-74-2	2.8E+02	PHYSPROP	7.4E-05	1.8E-06	PHYSPROP	2.0E-05	PHYSPROP	-3.5E+01	PHYSPROP	1.0E+00	CRC	2.1E-02	5.3E-06	WATER9 (U.S. EPA, 2001)			1.2E+03	EPI	4.5E+00	PHYSPROP	1.1E+01	PHYSPROP	2.7E-01	3.8E+00	9.1E+00	4.2E-02	EPI
~Diethyl Phthalate	84-66-2	2.2E+02	PHYSPROP	2.5E-05	6.1E-07	EPI	2.1E-03	PHYSPROP	-4.1E+01	PHYSPROP	1.2E+00	CRC	2.6E-02	6.7E-06	WATER9 (U.S. EPA, 2001)			1.0E+02	EPI	2.4E+00	PHYSPROP	1.1E+03	PHYSPROP	2.1E-02	1.8E+00	4.4E+00	3.6E-03	EPI
~Dimethylterephthalate	120-61-6	1.9E+02	PHYSPROP	5.5E-03	1.3E-04	EPI	1.0E-02	PHYSPROP	1.4E+02	PHYSPROP	1.1E+00	CRC	2.0E-02	6.7E-06	WATER9 (U.S. EPA, 2001)			3.1E+01	EPI	2.2E+00	PHYSPROP	1.9E+01	PHYSPROP	2.1E-02	1.3E+00	3.1E+00	4.0E-03	EPI
~Octyl Phthalate, di-N-	117-84-0	3.9E+02	PHYSPROP	1.1E-04	2.6E-06	EPI	1.0E-07	PHYSPROP	2.5E+01	PHYSPROP			3.6E-02	4.2E-06	WATER9 (U.S. EPA, 2001)			1.4E+05	EPI	8.1E+00	PHYSPROP	2.2E-02	PHYSPROP	1.8E+01	1.6E+01	7.5E+01	2.4E+00	EPI
~Phthalic Acid, p-	100-21-0	1.7E+02	PHYSPROP	1.6E-11	3.9E-13	PHYSPROP	9.2E-06	EPI	1.3E+02	EPI	1.5E+00	CRC	4.9E-02	9.1E-06	WATER9 (U.S. EPA, 2001)			7.9E+01	EPI	2.0E+00	PHYSPROP	1.5E+01	PHYSPROP	1.9E-02	9.0E-01	2.1E+00	3.9E-03	EPI
~Phthalic Anhydride	85-44-9	1.5E+02	PHYSPROP	6.7E-07	1.6E-08	EPI	5.2E-04	EPI	1.3E+02	PHYSPROP	1.5E+00	CRC	5.9E-02	9.8E-06	WATER9 (U.S. EPA, 2001)			1.0E+01	EPI	1.6E+00	PHYSPROP	6.2E+03	PHYSPROP	1.2E-02	7.1E-01	1.7E+00	2.7E-03	EPI
Picloram	1918-02-1	2.4E+02	PHYSPROP	2.2E-12	5.3E-14	EPI	7.2E-11	PHYSPROP	2.2E+02	PHYSPROP			4.9E-02	5.7E-06	WATER9 (U.S. EPA, 2001)			3.9E+01	EPI	1.9E+00	PHYSPROP	4.3E+02	PHYSPROP	7.6E-03	2.4E+00	5.7E+00	1.3E-03	EPI
Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3	2.0E+02	PHYSPROP	4.0E-10	9.8E-12	PHYSPROP	4.2E-07	PHYSPROP	1.7E+02	PHYSPROP			5.6E-02	6.5E-06	WATER9 (U.S. EPA, 2001)			2.3E+02	EPI	9.3E-01	PHYSPROP	1.4E+03	PHYSPROP	2.7E-03	1.4E+00	3.3E+00	5.0E-04	EPI
Picric Acid (2,4,6-Trinitrophenol)	88-89-1	2.3E+02	PHYSPROP	7.0E-10	1.7E-11	EPI	7.5E-07	PHYSPROP	1.2E+02	PHYSPROP	1.8E+00	PERRY	3.0E-02	8.2E-06	WATER9 (U.S. EPA, 2001)			2.3E+03	EPI	1.4E+00	PHYSPROP	1.3E+04	PHYSPROP	3.6E-03	2.0E+00	4.8E+00	6.2E-04	EPI
Pirimiphos, Methyl	29232-93-7	3.1E+02	PHYSPROP	2.9E-05	7.0E-07	EPI	1.5E-05	PHYSPROP	1.5E+01	PHYSPROP	1.2E+00	CRC	2.2E-02	5.4E-06	WATER9 (U.S. EPA, 2001)			3.7E+02	EPI	4.2E+00	PHYSPROP	8.6E+00	PHYSPROP	1.3E-01	5.4E+00	1.3E+01	1.9E-02	EPI
Polybrominated Biphenyls	36355-01-8																												
Polychlorinated Biphenyls (PCBs)																													
~Aroclor 1016	12674-11-2	2.6E+02	EPI	8.2E-03	2.0E-04	EPI	4.0E-04	PHYSPROP	1.0E+02	EPI	1.4E+00	ATSDR Profile	2.5E-02	6.6E-06	WATER9 (U.S. EPA, 2001)			4.8E+04	EPI	5.7E+00	PHYSPROP	4.2E-01	PHYSPROP	1.9E+00	2.9E+00	1.2E+01	3.1E-01	EPI
~Aroclor 1221	11104-28-2	1.9E+02	PHYSPROP	9.3E-03	2.3E-04	PHYSPROP	6.7E-03	PHYSPROP	3.4E+01	EPI	1.2E+00	ATSDR Profile	3.2E-02	7.2E-06	WATER9 (U.S. EPA, 2001)			8.4E+03	EPI	4.7E+00	PHYSPROP	1.5E+01	PHYSPROP	8.9E-01	1.2E+00	4.6E+00	1.7E-01	EPI
~Aroclor 1232	11141-16-5	1.9E+02	PHYSPROP	3.0E-02	7.4E-04	EPI	4.1E-03	PHYSPROP	3.4E+01	EPI	1.3E+00	ATSDR Profile	3.3E-02	7.5E-06	WATER9 (U.S. EPA, 2001)			8.4E+03	EPI	4.4E+00	PHYSPROP	1.5E+00	PHYSPROP	8.9E-01	1.2E+00	4.6E+00	1.7E-01	EPI
~Aroclor 1242	53469-21-9	2.3E+02	PHYSPROP	1.4E-02	3.4E-04	PHYSPROP	8.6E-05	EPI	1.4E+02	EPI	1.4E+00	ATSDR Profile	2.4E-02	6.1E-06	WATER9 (U.S. EPA, 2001)			7.8E+04	EPI	6.3E+00	PHYSPROP	2.8E-01	PHYSPROP	3.6E+00	4.5E+00	1.9E+01	5.5E-01	EPI
~Aroclor 1248	12672-29-6	2.9E+02	EPI	1.8E-02	4.4E-04	PHYSPROP	4.9E-04	PHYSPROP	1.2E+02	EPI	1.4E+00	HSDB	2.4E-02	6.2E-06	WATER9 (U.S. EPA, 2001)			7.7E+04	EPI	6.2E+00	PHYSPROP	1.0E-01	PHYSPROP	3.1E+00	4.5E+00	1.9E+01	4.8E-01	EPI
~Aroclor 1254	11097-69-1	3.3E+02	PHYSPROP	1.2E-02	2.8E-04	PHYSPROP	7.7E-05	PHYSPROP	1.3E+02	EPI	1.5E+00	ATSDR Profile	2.4E-02	6.1E-06	WATER9 (U.S. EPA, 2001)			1.3E+05	EPI	6.5E+00	PHYSPROP	4.3E-02	PHYSPROP	5.2E+00	7.1E+00	3.1E+01	7.5E-01	EPI
~Aroclor 1260	11096-82-5	4.0E+02	PHYSPROP	1.4E-02	3.4E-04	PHYSPROP	4.1E-05	PHYSPROP	1.6E+02	EPI	1.6E+00	ATSDR Profile	2.2E-02	5.6E-06	WATER9 (U.S. EPA, 2001)			3.5E+05	EPI	7.6E+00	PHYSPROP	1.4E-02	PHYSPROP	7.5E+00

Contaminant		Molecular Weight		Volatility Parameters					Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients						Water Solubility		Tapwater Dermal Parameters				
Analyte	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m^3/mole)	H' and HLC Ref	VP (mmHg)	VP Ref	MP (C)	MP Ref	Density (g/cm^3)	Density Ref	D_{ia} (cm^2/s)	D_{iw} (cm^2/s)	D_{ia} and D_{iw} Ref	K_d (L/kg)	K_d Ref	K_{oc} (L/kg)	K_{oc} Ref	log K_{ow} (unitless)	log K_{ow} Ref	S (mg/L)	S Ref	B (unitless)	t_{event} (hr/event)	t^* (hr)	K_p (cm/hr)	K_p Ref
~Methylnaphthalene, 1-	90-12-0	1.4E+02	PHYSPROP	2.1E-02	5.1E-04	PHYSPROP	6.7E-02	PHYSPROP	-3.0E+01	PHYSPROP	1.0E+00	CRC	5.3E-02	7.8E-06	WATER9 (U.S. EPA, 2001)			2.5E+03	EPI	3.9E+00	PHYSPROP	2.6E+01	PHYSPROP	4.3E-01	6.6E-01	1.6E+00	9.3E-02	EPI
~Methylnaphthalene, 2-	91-57-6	1.4E+02	PHYSPROP	2.1E-02	5.2E-04	PHYSPROP	5.5E-02	PHYSPROP	3.4E+01	PHYSPROP	1.0E+00	CRC	5.2E-02	7.8E-06	WATER9 (U.S. EPA, 2001)			2.5E+03	EPI	3.9E+00	PHYSPROP	2.5E+01	PHYSPROP	4.2E-01	6.6E-01	1.6E+00	9.2E-02	EPI
~Naphthalene	91-20-3	1.3E+02	PHYSPROP	1.8E-02	4.4E-04	PHYSPROP	8.5E-02	PHYSPROP	8.0E+01	PHYSPROP	1.0E+00	CRC	6.0E-02	8.4E-06	WATER9 (U.S. EPA, 2001)			1.5E+03	EPI	3.3E+00	PHYSPROP	3.1E+01	PHYSPROP	2.0E-01	5.5E-01	1.3E+00	4.7E-02	EPI
~Nitropyrene, 4-	57835-92-4	2.5E+02	PHYSPROP	1.0E-06	2.5E-08	PHYSPROP	5.5E-08	PHYSPROP	1.9E+02	PHYSPROP			4.8E-02	5.6E-06	WATER9 (U.S. EPA, 2001)			8.6E+04	EPI	4.8E+00	PHYSPROP	6.8E-02	PHYSPROP	5.6E-01	2.6E+00	6.3E+00	9.2E-02	EPI
~Pyrene	129-00-0	2.0E+02	PHYSPROP	4.9E-04	1.2E-05	PHYSPROP	4.5E-06	PHYSPROP	1.5E+02	PHYSPROP	1.3E+00	CRC	2.8E-02	7.2E-06	WATER9 (U.S. EPA, 2001)			5.4E+04	EPI	4.9E+00	PHYSPROP	1.4E-01	PHYSPROP	1.1E+00	1.4E+00	5.5E+00	2.0E-01	EPI
Prochloraz	67747-09-5	3.8E+02	PHYSPROP	6.7E-07	1.6E-08	EPI	1.1E-06	PHYSPROP	4.8E+01	PHYSPROP			3.6E-02	4.3E-06	WATER9 (U.S. EPA, 2001)			2.4E+03	EPI	4.1E+00	PHYSPROP	3.4E+01	PHYSPROP	4.8E-02	1.4E+01	3.2E+01	6.4E-03	EPI
Profluralin	26399-36-0	3.5E+02	PHYSPROP	1.2E-02	2.9E-04	EPI	6.3E-05	PHYSPROP	3.2E+01	PHYSPROP	1.4E+00	HSDB	2.2E-02	5.5E-06	WATER9 (U.S. EPA, 2001)			3.1E+04	EPI	5.6E+00	PHYSPROP	1.0E-01	PHYSPROP	6.5E-01	9.3E-00	3.7E+01	9.0E-02	EPI
Prometon	1610-18-0	2.3E+02	PHYSPROP	3.7E-08	9.1E-10	EPI	2.3E-06	PHYSPROP	9.1E+01	PHYSPROP			5.1E-02	6.0E-06	WATER9 (U.S. EPA, 2001)			1.4E+02	EPI	3.0E+00	PHYSPROP	7.5E+02	PHYSPROP	4.8E-02	1.9E+00	4.6E+00	8.3E-03	EPI
Prometryn	7287-19-6	2.4E+02	PHYSPROP	4.9E-07	1.2E-08	EPI	1.2E-06	PHYSPROP	1.2E+02	PHYSPROP	1.2E+00	CRC	2.4E-02	6.2E-06	WATER9 (U.S. EPA, 2001)			6.6E+02	EPI	3.5E+00	PHYSPROP	3.3E+01	PHYSPROP	8.9E-02	2.4E+00	5.7E+00	1.5E-02	EPI
Pronamide	23950-58-5	2.6E+02	PHYSPROP	4.0E-07	9.8E-09	EPI	4.4E-07	PHYSPROP	1.6E+02	PHYSPROP			4.7E-02	5.5E-06	WATER9 (U.S. EPA, 2001)			4.0E+02	EPI	3.4E+00	PHYSPROP	1.5E+01	PHYSPROP	6.7E-02	2.9E+00	6.9E+00	1.1E-02	EPI
Propachlor	1918-16-7	2.1E+02	PHYSPROP	1.5E-05	3.6E-07	EPI	2.3E-04	PHYSPROP	7.7E+01	PHYSPROP	1.2E+00	CRC	2.7E-02	7.0E-06	WATER9 (U.S. EPA, 2001)			2.0E+02	EPI	2.2E+00	PHYSPROP	5.8E+02	PHYSPROP	1.6E-02	1.6E+00	3.9E+00	2.9E-03	EPI
Propanil	709-98-8	2.2E+02	PHYSPROP	7.0E-08	1.7E-09	EPI	9.1E-07	PHYSPROP	9.2E+01	PHYSPROP	1.3E+00	CRC	2.7E-02	6.9E-06	WATER9 (U.S. EPA, 2001)			1.8E+02	EPI	3.1E+00	PHYSPROP	1.5E+02	PHYSPROP	5.9E-02	1.8E+00	4.2E+00	1.0E-02	EPI
Propargite	2312-35-8	3.5E+02	PHYSPROP	2.6E-05	6.4E-07	EPI	3.0E-07	PHYSPROP	1.7E+02	EPI	1.1E+00	CRC	1.9E-02	4.8E-06	WATER9 (U.S. EPA, 2001)			3.7E+04	EPI	5.0E+00	PHYSPROP	2.2E-01	PHYSPROP	2.6E-01	9.7E+00	2.3E+01	3.6E-02	EPI
Propargyl Alcohol	107-19-7	5.6E+01	PHYSPROP	4.7E-05	1.2E-06	EPI	1.6E+01	PHYSPROP	-5.0E+01	PHYSPROP	9.5E-01	CRC	1.2E-01	1.3E-05	WATER9 (U.S. EPA, 2001)			1.9E+00	EPI	-3.8E-01	PHYSPROP	1.0E+06	PHYSPROP	1.2E-03	2.2E-01	5.2E-01	4.2E-04	EPI
Propazine	139-40-2	2.3E+02	PHYSPROP	1.9E-07	4.6E-09	EPI	1.3E-07	PHYSPROP	2.1E+02	PHYSPROP	1.2E+00	CRC	2.5E-02	6.4E-06	WATER9 (U.S. EPA, 2001)			3.4E+02	EPI	2.9E+00	PHYSPROP	8.6E+00	PHYSPROP	4.2E-02	2.0E+00	4.9E+00	7.1E-03	EPI
Propham	122-42-9	1.8E+02	PHYSPROP	7.5E-06	1.8E-07	EPI	1.4E-04	PHYSPROP	8.7E+01	PHYSPROP	1.1E+00	CRC	3.6E-02	7.1E-06	WATER9 (U.S. EPA, 2001)			2.2E+02	EPI	2.6E+00	PHYSPROP	1.8E+02	PHYSPROP	4.3E-02	1.1E+00	2.5E+00	8.3E-03	EPI
Propiconazole	60207-90-1	3.4E+02	PHYSPROP	7.0E-08	1.7E-09	EPI	4.2E-07	PHYSPROP	1.7E+02	EPI	1.3E+00	CRC	2.1E-02	5.3E-06	WATER9 (U.S. EPA, 2001)			1.6E+03	EPI	3.7E+00	PHYSPROP	1.1E+02	PHYSPROP	4.0E-02	8.7E+00	2.1E+01	5.6E-03	EPI
Propionaldehyde	123-38-6	5.8E+01	PHYSPROP	3.0E-03	7.3E-05	PHYSPROP	3.2E+02	PHYSPROP	-8.0E+01	PHYSPROP	8.7E-01	CRC	1.1E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	5.9E-01	PHYSPROP	3.1E+05	PHYSPROP	5.3E-03	2.2E-01	5.3E-01	1.8E-03	EPI
Propyl benzene	103-65-1	1.2E+02	PHYSPROP	4.3E-01	1.1E-02	PHYSPROP	3.4E+00	PHYSPROP	-1.0E+02	PHYSPROP	8.6E-01	CRC	6.0E-02	7.8E-06	WATER9 (U.S. EPA, 2001)			8.1E+02	EPI	3.7E+00	PHYSPROP	5.2E+01	PHYSPROP	4.0E-01	5.0E-01	1.2E+00	9.4E-02	EPI
Propylene	115-07-1	4.2E+01	PHYSPROP	8.0E+00	2.0E-01	PHYSPROP	8.7E+03	PHYSPROP	2.9E+02	PHYSPROP	5.1E-01	CRC	1.1E-01	1.1E-05	WATER9 (U.S. EPA, 2001)			2.2E+01	EPI	1.8E+00	PHYSPROP	2.0E+02	PHYSPROP	3.4E-02	1.8E-01	4.3E-01	1.4E-02	EPI
Propylene Glycol	57-55-6	7.6E+01	PHYSPROP	5.3E-07	1.3E-08	EPI	1.3E-01	PHYSPROP	-6.0E+01	PHYSPROP	1.0E+00	CRC	9.8E-02	1.2E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-9.2E-01	PHYSPROP	1.0E+06	PHYSPROP	4.8E-04	2.8E-01	6.7E-01	1.4E-04	EPI
Propylene Glycol Dinitrate	6423-43-4	1.7E+02	PHYSPROP	3.9E-05	9.4E-07	PHYSPROP	3.8E-01	PHYSPROP	-9.6E+00	EPI			6.3E-02	7.3E-06	WATER9 (U.S. EPA, 2001)			6.1E+01	EPI	1.8E+00	PHYSPROP	3.3E+03	EPI	1.0E-02	9.0E-01	2.1E+00	2.1E-03	EPI
Propylene Glycol Monomethyl Ether	107-98-2	9.0E+01	PHYSPROP	3.8E-05	9.2E-07	PHYSPROP	1.3E+01	PHYSPROP	-9.5E+01	PHYSPROP	9.6E-01	CRC	8.3E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.0E+00	EPI	-4.9E-01	PHYSPROP	1.0E+06	PHYSPROP	8.4E-04	3.4E-01	8.1E-01	2.3E-04	EPI
Propylene Oxide	75-56-9	5.8E+01	PHYSPROP	2.8E-03	7.0E-05	EPI	5.4E+02	PHYSPROP	-1.1E+02	PHYSPROP	8.3E-01	PERRY	1.1E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			5.2E+00	EPI	3.0E-02	PHYSPROP	5.9E+05	PHYSPROP	2.3E-03	2.2E-01	5.3E-01	7.7E-04	EPI
Pyridine	110-86-1	7.9E+01	PHYSPROP	4.5E-04	1.1E-05	PHYSPROP	2.1E+01	PHYSPROP	-4.2E+01	PHYSPROP	9.8E-01	CRC	9.3E-02	1.1E-05	WATER9 (U.S. EPA, 2001)			7.2E+01	EPI	6.5E-01	PHYSPROP	1.0E+06	PHYSPROP	5.2E-03	2.9E-01	7.0E-01	1.5E-03	EPI
Quinalphos	13593-03-8	3.0E+02	PHYSPROP	1.9E-06	4.6E-08	EPI	2.6E-06	PHYSPROP	3.2E+01	PHYSPROP			4.3E-02	5.0E-06	WATER9 (U.S. EPA, 2001)			4.2E+03	EPI	4.4E+00	PHYSPROP	2.2E+01	PHYSPROP	2.0E-01	4.9E+00	1.2E+01	3.0E-02	EPI
Quinoline	91-22-5	1.3E+02	PHYSPROP	6.8E-05	1.7E-06	EPI	6.0E-02	PHYSPROP	-1.5E+01	PHYSPROP	1.1E+00	CRC	6.2E-02	8.7E-06	WATER9 (U.S. EPA, 2001)			1.5E+03	EPI	2.0E+00	PHYSPROP	6.1E+03	PHYSPROP	2.9E-02	5.6E-01	1.3E+00	6.6E-03	EPI
Quizalofop-ethyl	76578-14-8	3.7E+02	PHYSPROP	4.3E-07	1.1E-08	EPI	6.5E-09	PHYSPROP	9.2E+01	PHYSPROP			3.7E-02	4.3E-06	WATER9 (U.S. EPA, 2001)			7.7E+03	EPI	4.3E+00	PHYSPROP	3.0E-01	PHYSPROP	6.6E-02	1.3E+01	3.1E+01	8.9E-03	EPI
Refractory Ceramic Fibers (units in fibers)	E715557																										1.0E-03	RAGSE
Resmethrin	10453-86-8	3.4E+02	PHYSPROP	5.4E-06	1.3E-07	EPI	1.1E-08	PHYSPROP	5.7E+01	PHYSPROP			3.9E-02	4.6E-06	WATER9 (U.S. EPA, 2001)			3.1E+05	EPI	6.1E+00	PHYSPROP	3.8E-02	PHYSPROP	1.7E+00	8.3E+00	3.3E+01	2.4E-01	EPI
Ronnel	299-84-3	3.2E+02	PHYSPROP	1.3E-03	3.2E-05	EPI	7.5E-05	PHYSPROP	4.1E+01	PHYSPROP	1.4E+00	CRC	2.3E-02	5.9E-06	WATER9 (U.S. EPA, 2001)			4.5E+03	EPI	4.9E+00	PHYSPROP	1.0E+00	PHYSPROP	3.0E-01	6.6E+00	1.6E+01	4.3E-02	EPI
Rotenone	83-79-4	3.9E+02	PHYSPROP	4.6E-12	1.1E-13	PHYSPROP	6.9E-10	PHYSPROP	1.8E+02	PHYSPROP			3.5E-02	4.1E-06	WATER9 (U.S. EPA, 2001)			2.6E+05	EPI	4.1E+00	PHYSPROP	2.0E-01	PHYSPROP	3.9E-02	1.7E+01	4.1E+01	5.1E-03	EPI
Safrole	94-59-7	1.6E+02	PHYSPROP	3.7E-04	9.1E-06	PHYSPROP	7.5E-02	CRC	1.1E+01	PHYSPROP	1.1E+00	CRC	4.4E-02	7.5E-06	WATER9 (U.S. EPA, 2001)			2.1E+02	EPI	3.5E+00	PHYSPROP	1.2E+02	PHYSPROP	5.5E-02	8.5E-01	2.0E+00	1.1E-02	RAGSE
Selenious Acid	7783-00-8	1.3E+02	PHYSPROP						7.0E+01	EPI	3.0E+00	CRC										9.0E+05	PERRY	4.4E-03	5.5E-01	1.3E+00	1.0E-03	RAGSE
Selenium	7782-49-2	7.9E+01	PHYSPROP				1.4E-10	EPI	2.2E+02	PHYSPROP	4.8E+00	CRC				5.0E+00	SSL							3.4E-03	2.9E-01	7.0E-01	1.0E-03	RAGSE
Selenium Sulfide	7446-34-6	1.1E+02	EPI																					4.1E-03	4.4E-01	1.1E+00	1.0E-03	RAGSE
Sethoxydim	74051-80-2	3.3E+02	PHYSPROP	8.8E-10	2.2E-11	PHYSPROP	1.6E-07	PHYSPROP	1.6E+02	EPI	1.0E+00	CRC	2.0E-02	4.8E-06	WATER9 (U.S. EPA, 2001)			4.4E+03	EPI	4.4E+00	PHYSPROP	2.5E+01	PHYSPROP	1.3E-01	7.2E+00	1.7E+01	1.9E-02	EPI
Silica (crystalline. respirable)	7631-86-9	6.0E+01	EPI						1.7E+03	PERRY	2.3E+00	PERRY												3.0E-03	2.3E-01	5.5E-01	1.0E-03	RAGSE
Silver	7440-22-4	1.1E+02	EPI				0.0E+00	NIOSH	9.6E+02	PHYSPROP	1.1E+01	CRC				8.3E+00	SSL							2.4E-03	4.2E-01	1.0E+00	6.0E-04	RAGSE
Simazine	122-34-9	2.0E+02	PHYSPROP	3.9E-08	9.4E-10	EPI	2.2E-08	PHYSPROP	2.3E+02	PHYSPROP	1.3E+00	CRC	2.8E-02	7.4E-06	WATER9 (U.S. EPA, 2001)			1.5E+02	EPI	2.2E+00	PHYSPROP	6.2E+00	PHYSPROP	1.8E-02	1.4E+00	3.4E+00	3.3E-03	EPI
Sodium Acifluorfen	62476-59-9	3.8E+02	PHYSPROP	2.5E-09	6.1E-11	PHYSPROP	9.8E-09	EPI	2.8E+02	EPI			3.6E-02	4.2E-06	WATER9 (U.S. EPA, 2001)			3.9E+03	EPI	3.7E-01	PHYSPROP	2.5E+05	PHYSPROP	1.5E-04	1.5E+01	3.6E+01	2.0E-05	EPI
Sodium Azide	26628-22-8	6.5E+01	EPI						3.0E+02	CRC	1.8E+00	CRC										4.1E+05	CRC	3.1E-03	2.4E-01	5.8E-01	1.0E-03	RAGSE
Sodium Diethyldithiocarbamate	148-18-5	1.7E+02	PHYSPROP				8.2E-10	PHYSPROP	9.4E+01	PHYSPROP			6.1E-02	7.2E-06	WATER9 (U.S. EPA, 2001)			2.0E+02	EPI	-1.4E+00	PHYSPROP	3.6E+05	PHYSPROP	9.7E-05	9.7E-01	2.3E+00	1.9E-05	EPI
Sodium Fluoride	7681-49-4	4.2E+01	PHYSPROP				0.0E+00	NIOSH	9.9E+02	PHYSPROP	2.8E+00	CRC				1.5E+02	BAES					4.2E+04	PHYSPROP	2.5E-03	1.8E-01	4.3E-01	1.0E-03	RAGSE
Sodium Fluoroacetate	62-74-8	1.0E+02	PHYSPROP	4.5E-05	1.1E-06	PHYSPROP	6.5E-07	PHYSPROP	2.0E+02	PHYSPROP			8.8E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.4E+00	EPI	-3.8E+00	PHYSPROP	1.1E+06	PHYSPROP	5.1E-06	3.8E-01	9.2E-01	1.3E-06	EPI
Sodium Metavanadate	13718-26-8	1.2E+02	CRC						6.3E+02	CRC												2.1E+05	CRC	4.2E-03	5.1E-01	1.2E+00	1.0E-03	RAGSE
Sodium Tungstate	13472-45-2	2.9E+02	CRC						7.0E+02	CRC	4.2E+00	CRC										7.4E+05	CRC	6.6E-03	4.6E+00	1.1E+01	1.0E-03	RAGSE
Sodium Tungstate Dihydrate	10213-10-2	3.3E+02	CRC						1.0E+02	CRC	3.3E+00	CRC										7.4E+05	CRC	7.0E-03	7.4E+00	1.8E+01	1.0E-03	RAGSE
Stirofos (Tetrachlorovinphos)	961-11-5	3.7E+02	PHYSPROP	7.5E-08	1.8E-09	EPI	4.2E-08	PHYSPROP	9.8E+01	PHYSPROP			3.7E-02	4.3E-06	WATER9 (U.S. EPA, 2001)			1.4E+03	EPI	3.5E+00	PHYSPROP	1.1E+01	PHYSPROP	2.3E-02	1.2E+01	2.8E+01	3.1E-03	EPI
Strontium, Stable	7440-24-6	8.8E+01	PHYSPROP						7.8E+02	PHYSPROP	2.6E+00	CRC				3.5E+01	BAES							3.6E-03	3.3E-01	7.8E-01	1.0E-03	RAGSE
Strychnine	57-24-9	3.3E+02	PHYSPROP	3.1E-12	7.6E-14	PHYSPROP	2.9E-09	PHYSPROP	2.9E+02	PHYSPROP	1.4E+00	CRC	2.2E-02	5.6E-06	WATER9 (U.S. EPA, 2001)			5.4E+03	EPI	1.9E+00	PHYSPROP	1.6E+02	PHYSPROP	2.8E-03	7.8E+00	1.9E+01	4.0E-04	EPI
Styrene	100-42-5	1.0E+02	PHYSPROP	1.1E-01	2.8E-03	PHYSPROP	6.4E+00	PHYSPROP	-3.1E+01	PHYSPROP	9.0E-01	CRC	7.1E-02	8.8E-06	WATER9 (U.S. EPA, 2001)			4.5E+02	EPI	3.0E+00	PHYSPROP	3.1E+02	PHYSPROP	1.5E-01	4.0E-01	9.7E-01	3.7E-02	EPI
Styrene-Acrylonitrile (SAN) Trimer (THNA isomer)	57964-39-3	2.1E+02	PPRTV						1.1E+00	PPRTV	1.1E+00	PPRTV	2.6E-02	6.5E-06	WATER9 (U.S. EPA, 2001)					2.8E+00	EPI	4.4E-04	EPI	6.6E-02	1.6E+00	3.8E+00	1.2E-02	RAGSE
Styrene-Acrylonitrile (SAN) Trimer (THNP isomer)	57964-40-6	2.1E+02	PPRTV				1.1E-07	ChemIndex	1.1E+00	PPRTV	2.6E-02	6.5E-06	WATER9 (U.S. EPA, 2001)							3.1E+00	US NTP	8.5E+01	PPRTV	6.6E-02	1.6E+00	3.8E+00	1.2E-02	RAGSE
Sulfolane	126-33-0	1.2E+02	PHYSPROP	2.0E-04	4.9E-06	PHYSPROP	4.1E-03	EPI	2.8E+01	PHYSPROP	1.3E+00	CRC	7.2E-02	9.9E-06	WATER9 (U.S. EPA, 2001)			9.1E+00	EPI	-7.7E-01	PHYSPROP	1.0E+06	PHYSPROP	4.3E-04	5.0E-01	1.2E+00	1.0E-04	EPI
Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9	2.9E+02	PHYSPROP	5.6E-06	1.4E-07	PHYSPROP	8.1E-07	PHYSPROP	1.5E+02	PHYSPROP			4.4E-02	5.1E-06	WATER9 (U.S. EPA, 2001)			2.9E+03	EPI	3.9E+00	PHYSPROP	2.4E+00	PHYSPROP	9.7E-02	4.3E+00	1.0E+01	1.5E-02	EPI
Sulfur Trioxide	7446-11-9	8.0E+01	PHYSPROP				2.6E+02	PHYSPROP	1.7E+01	PHYSPROP	1.9E+00	CRC	1.2E-01	1.6E-05	WATER9 (U.S. EPA, 2001)									3.4E-03	3.0E-01	7.1E-01	1.0E-03	RAGSE
Sulfuric Acid	7664-93-9	9.8E+01	PHYSPROP				5.9E-05	PHYSPROP	1.0E+01	PHYSPROP	1.8E+00	CRC										1.0E+06	PHYSPROP	3.8E-03	3.7E-01	8.9E-01	1.0E-03	RAGSE
Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl ester	140-57-8	3.3E+02	PHYSPROP	7.8E-06	1.9E-07	PHYSPROP	2.2E-07	PHYSPROP	-3.0E+01	PHYSPROP	1.1E+00	CRC	2.0E-02	5.0E-06	WATER9 (U.S. EPA, 2001)			5.6E+03	EPI	4.8E+00	PHYSPROP	5.9E-01	PHYSPROP	2.3E-01	7.9E+00	1.9E+01	3.3E-02	EPI
TCMTB	21564-17-0	2.4E+02	PHYSPROP	2.7E-10	6.5E-12	PHYSPROP	3.1E-07	PHYSPROP	1.5E+02	EPI			4.9E-02	5.8E-06	WATER9 (U.S. EPA, 2001)			3.4E+03	EPI	3.3E+00	PHYSPROP	1.3E+02	PHYSPROP	6.7E-02	2.3E+00	5.5E+00	1.1E-02	EPI
Tebuthiuron	34014-18-1	2.3E+02	PHYSPROP	4.9E-09	1.2E-10	PHYSPROP	3.0E-07	PHYSPROP	1.6E+02	PHYSPROP			5.1E-02	5.9E-06	WATER9 (U.S. EPA, 2001)			4.2E+01	EPI	1.8E+00	PHYSPROP	2.5E+03	PHYSPROP	7.4E-03	2.0E+00	4.8E+00	1.3E-03	EPI
Temephos	3383-96-8	4.7E+02	PHYSPROP	8.0E-08	2.0E-09	PHYSPROP	7.9E-08	PHYSPROP	3.0E+01	PHYSPROP	1.3E+00	CRC	1.8E-02	4.5E-06	WATER9 (U.S. EPA, 2001)			9.5E+04	EPI	6.0E+00	PHYSPROP	2.7E-01	PHYSPROP	2.9E-01	4.3E+01	1.0E+02	3.5E-02	EPI
Terbacil	5902-51-2	2.2E+02	PHYSPROP	4.9E-09	1.2E-10	EPI	4.7E-07	PHYSPROP	1.8E+02	PHYSPROP	1.3E+00	CRC	2.7E-02	7.2E-06	WATER9 (U.S. EPA, 2001)			5.0E+01	EPI	1.9E+00	PHYSPROP	7.1E+02	PHYSPROP	9.7E-03	1.7E+00	4.1E+00	1.7E-03	EPI
Terbufos	13071-79-9	2.9E+02	PHYSPROP	9.8E-04	2.4E-05	EPI	3.2E-04	PHYSPROP	-2.9E+01	PHYSPROP	1.1E+00	CRC	2.2E-02	5.4E-06	WATER9 (U.S. EPA, 2001)			1.0E+03	EPI	4.5E+00	PHYSPROP	5.1E+00	PHYSPROP	2.3E-01	4.3E+00	1.0E+01	3.6E-02	EPI
Terbutryn	886-50-0	2.4E+02	PHYSPROP	8.8E-07	2.2E-08	EPI	1.7E-06	PHYSPROP	1.0E+02	PHYSPROP	1.1E+00	CRC	2.4E-02	6.0E-06	WATER9 (U.S. EPA, 2001)			6.1E+02	EPI	3.7E+00	PHYSPROP	2.5E+01	PHYSPROP	1.3E-01	2.4E+00	5.7E+00	2.1E-02	EPI
Tert-Butyl Acetate	540-88-5	1.2E+02	PHYSPROP	3.5E-02	8.6E-04	EPI	4.7E+01	PHYSPROP	2.5E+01	PHYSPROP			8.6E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.2E+01	EPI	1.6E+00	PHYSPROP	9.3E+04	PHYSPROP	2.4E-02	4.7E-01	1.1E+00	5.2E-03	EPI
Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1	4.9E+02	PHYSPROP	1.2E-04	3.0E-06	PHYSPROP	7.0E-08	EPI	1.6E+02	EPI			3.1E-02	3.6E-06	WATER9 (U.S. EPA, 2001)			1.3E+04	EPI	6.8E+00	PHYSPROP	1.5E-03	PHYSPROP	7.9E-01	5.5E+01	2.1E+02	9.3E-02	EPI
Tetrachlorobenzene, 1,2,4,5-	95-94-3	2.2E+02	PHYSPROP	4.1E-02	1.0E-03	PHYSPROP	5.4E-03	EPI	1.4E+02	PHYSPROP	1.9E+00	CRC	3.2E-02	8.8E-06	WATER9 (U.S. EPA, 2001)			2.2E+03	EPI	4.6E+00	PHYSPROP	6.0E-01	PHYSPROP	6.6E-01	1.7E+00	6.7E+00	1.2E-01	EPI
Tetrachloroethane, 1,1,1,2-	630-20-6	1.7E+02	PHYSPROP	1.0E-01	2.5E-03	PHYSPROP	1.2E+01	PHYSPROP	-7.0E+01	PHYSPROP	1.5E+00	CRC	4.8E-02	9.1E-06	WATER9 (U.S. EPA, 2001)			8.6E+01	EPI	2.9E+00	PHYSPROP	1.1E+03	PHYSPROP	7.9E-02	9.2E-01	2.2E+00	1.6E-02	EPI
Tetrachloroethane. 1.1.2.2-	79-34-5	1.7E+02	PHYSPROP	1.5E-02	3.7E-04	PHYSPROP	4.6E+00	PHYSPROP	-4.4E+01	PHYSPROP	1.6E+00	CRC	4.9E-02	9.3E-06	WATER9 (U.S. EPA, 2001)			9.5E+01	EPI	2.4E+00	PHYSPROP	2.8E+03	PHYSPROP	3.5E-02	9.2E-01	2.2E+00	6.9E-03	EPI
Tetrachloroethylene	127-18-4	1.7E+02	PHYSPROP	7.2E-01	1.8E-02	PHYSPROP	1.9E+01	PHYSPROP	-2.2E+01	PHYSPROP	1.6E+00	CRC	5.0E-02	9.5E-06	WATER9 (U.S. EPA, 2001)			9.5E+01	EPI	3.4E+00	PHYSPROP	2.1E+02	PHYSPROP	1.7E-01	8.9E-01	2.1E+00	3.3E-02	EPI
Tetrachlorophenol, 2,3,4,6-	58-90-2	2.3E+02	PHYSPROP	3.6E-04	8.8E-06	EPI	6.7E-04	EPI	7.0E+01	PHYSPROP			5.6E-02	8.1E-06	WATER9 (U.S. EPA, 2001)			5.0E+02	EPI	4.5E+00	PHYSPROP	4.2E+01	PHYSPROP	4.2E-01	2.1E+00	5.0E+00	7.1E-02	EPI
Tetrachlorotoluene, 1- alpha, alpha, alpha-	5216-25-1	2.3E+02	PHYSPROP	7.9E-03	1.9E-04	PHYSPROP	3.8E-02	PHYSPROP	4.0E+01	EPI	1.4E+00	CRC	2.8E-02	7.3E-06	WATER9 (U.S. EPA, 2001)			1.6E+03	EPI	4.5E+00	PHYSPROP	4.0E+00	PHYSPROP	4.9E-01	2.0E+00	4.9E+00	8.4E-02	EPI
Tetraethyl Dithiopyrophosphate	3689-24-5	3.2E+02	PHYSPRO

Contaminant		Molecular Weight		Volatility Parameters					Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients						Water Solubility		Tapwater Dermal Parameters				
Analyte	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m^3/mole)	H' and HLC Ref	VP (mmHg)	VP Ref	MP (C)	MP Ref	Density (g/cm^3)	Density Ref	D_{oa} (cm^2/s)	D_{iw} (cm^2/s)	D_{ia} and D_{iw} Ref	K_d (L/kg)	K_d Ref	K_{oc} (L/kg)	K_{oc} Ref	log K_{ow} (unitless)	log K_{ow} Ref	S (mg/L)	S Ref	B (unitless)	τ_{event} (hr/event)	t^* (hr)	K_p (cm/hr)	K_p Ref
Toxaphene, Weathered	E1841606	4.5E+02	PHYSPROP	2.5E-04	6.0E-06	PHYSPROP	6.7E-06	PHYSPROP	7.7E+01	PHYSPROP	1.7E+00	SURROGATE	2.1E-02	5.3E-06	WATER9 (U.S. EPA, 2001)			7.7E+04	EPI	5.9E+00	PHYSPROP	5.5E-01	PHYSPROP	4.2E-01	3.4E+01	8.2E+01	5.2E-02	EPI
Tralomethrin	66841-25-6	6.7E+02	PHYSPROP	1.6E-08	3.9E-10	EPI	3.6E-11	PHYSPROP	1.4E+02	PHYSPROP			2.5E-02	2.9E-06	WATER9 (U.S. EPA, 2001)			1.9E+05	EPI	7.6E+00	PHYSPROP	8.0E-02	PHYSPROP	3.0E-01	5.6E+02	1.3E+03	3.1E-02	EPI
Tri-n-butyltin	688-73-3	2.9E+02	PHYSPROP	6.2E+01	1.5E+00	PHYSPROP	4.0E-02	PHYSPROP	2.9E+01	EPI	1.1E+00	CRC	2.1E-02	5.4E-06	WATER9 (U.S. EPA, 2001)			8.1E+03	EPI	4.1E+00	PHYSPROP	7.3E-03	PHYSPROP	1.3E-01	4.5E+00	1.1E+01	1.9E-02	EPI
Triacetin	102-76-1	2.2E+02	PHYSPROP	5.0E-07	1.2E-08	EPI	2.5E-03	PHYSPROP	7.8E+01	PHYSPROP	1.2E+00	CRC	2.6E-02	6.5E-06	WATER9 (U.S. EPA, 2001)			4.1E+01	EPI	2.5E-01	PHYSPROP	5.8E+04	PHYSPROP	7.8E-04	1.8E+00	4.2E+00	1.4E-04	EPI
Triadimefon	43121-43-3	2.9E+02	PHYSPROP	3.3E-09	8.1E-11	EPI	1.5E-08	PHYSPROP	8.2E+01	PHYSPROP	1.2E+00	CRC	2.2E-02	5.7E-06	WATER9 (U.S. EPA, 2001)			3.0E+02	EPI	2.8E+00	PHYSPROP	7.2E+01	PHYSPROP	1.6E-02	4.6E+00	1.1E+01	2.4E-03	EPI
Triallate	2303-17-5	3.0E+02	PHYSPROP	4.9E-04	1.2E-05	EPI	1.2E-04	PHYSPROP	2.9E+01	PHYSPROP	1.3E+00	CRC	2.2E-02	5.7E-06	WATER9 (U.S. EPA, 2001)			1.0E+03	EPI	4.6E+00	PHYSPROP	4.0E+00	PHYSPROP	2.3E-01	5.3E+00	1.3E+01	3.5E-02	EPI
Triasulfuron	82097-50-5	4.0E+02	PHYSPROP	1.3E-11	3.2E-13	PHYSPROP	5.5E-12	PHYSPROP	1.9E+02	PHYSPROP			3.5E-02	4.1E-06	WATER9 (U.S. EPA, 2001)			4.3E+02	EPI	1.1E+00	PHYSPROP	3.2E+01	PHYSPROP	3.6E-04	1.9E+01	4.5E+01	4.7E-05	EPI
Tribenuron-methyl	101200-48-0	4.0E+02	PHYSPROP	4.2E-12	1.0E-13	PHYSPROP	3.9E-10	PHYSPROP	1.4E+02	PHYSPROP			3.5E-02	4.1E-06	WATER9 (U.S. EPA, 2001)			9.5E+01	EPI	7.8E-01	PHYSPROP	5.0E+01	PHYSPROP	3.6E-03	1.7E+01	4.1E+01	4.7E-04	EPI
Tribromobenzene, 1,2,4-	615-54-3	3.1E+02	PHYSPROP	1.4E-02	3.4E-04	PHYSPROP	5.5E-03	PHYSPROP	4.5E+01	PHYSPROP	2.3E+00	ChemNet	2.9E-02	7.9E-06	WATER9 (U.S. EPA, 2001)			6.1E+02	EPI	4.7E+00	PHYSPROP	4.9E+00	PHYSPROP	2.3E-01	6.1E+00	1.5E+01	3.4E-02	EPI
Tribromophenol, 2,4,6-	118-79-6	3.3E+02	PHYSPROP	1.5E-06	3.6E-08	PHYSPROP	3.0E-04	PHYSPROP	9.6E+01	PHYSPROP	2.6E+00	CRC	3.0E-02	8.2E-06	WATER9 (U.S. EPA, 2001)			8.1E+02	EPI	4.1E+00	PHYSPROP	7.0E+01	PHYSPROP	8.5E-02	7.5E+00	1.8E+01	1.2E-02	EPI
Tribufos	78-48-8	3.1E+02	PHYSPROP	1.2E-05	2.9E-07	PHYSPROP	5.3E-06	PHYSPROP	-2.5E+01	CRC	1.1E+00	CRC	2.0E-02	5.0E-06	WATER9 (U.S. EPA, 2001)			2.4E+03	EPI	5.7E+00	PHYSPROP	2.3E+00	PHYSPROP	1.1E+00	6.1E+00	2.4E+01	1.7E-01	EPI
Tributyl Phosphate	126-73-8	2.7E+02	PHYSPROP	5.8E-05	1.4E-06	EPI	1.1E-03	PHYSPROP	-7.9E+01	PHYSPROP	9.7E-01	CRC	2.1E-02	5.2E-06	WATER9 (U.S. EPA, 2001)			2.4E+03	EPI	4.0E+00	PHYSPROP	2.8E+02	PHYSPROP	1.4E-01	3.3E+00	7.8E+00	2.3E-02	EPI
Tributyltin Compounds	E1790679																												
Tributyltin Oxide	56-35-9	6.0E+02	PHYSPROP	1.2E-05	3.0E-07	EPI	7.5E-06	PHYSPROP	-4.5E+01	PHYSPROP	1.2E+00	CRC	1.5E-02	3.6E-06	WATER9 (U.S. EPA, 2001)			2.6E+07	EPI	4.1E+00	PHYSPROP	2.0E+01	PHYSPROP	2.4E-03	2.3E+02	5.5E+02	2.5E-04	EPI
Trichloramine	10025-85-1																												
Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1	1.9E+02	PHYSPROP	2.2E+01	5.3E-01	EPI	3.6E+02	PHYSPROP	-3.5E+01	PHYSPROP	1.6E+00	CRC	3.8E-02	8.6E-06	WATER9 (U.S. EPA, 2001)			2.0E+02	EPI	3.2E+00	PHYSPROP	1.7E+02	PHYSPROP	9.2E-02	1.2E+00	2.8E+00	1.8E-02	EPI
Trichloroacetic Acid	76-03-9	1.6E+02	PHYSPROP	5.5E-07	1.4E-08	PHYSPROP	6.0E-02	EPI	5.8E+01	PHYSPROP	1.6E+00	CRC	5.2E-02	9.5E-06	WATER9 (U.S. EPA, 2001)			3.2E+00	EPI	1.3E+00	PHYSPROP	5.5E+04	PHYSPROP	7.1E-03	8.6E-01	2.1E+00	1.5E-03	EPI
Trichloroaniline HCl, 2,4,6-	33663-50-2	2.3E+02	EPI	2.9E-12	7.2E-14	EPI	6.1E-08	EPI	1.8E+02	EPI			5.0E-02	5.9E-06	WATER9 (U.S. EPA, 2001)			1.3E+03	EPI	-6.7E-01	EPI	2.1E+01	EPI	1.6E-04	2.1E+00	5.1E+00	2.8E-05	EPI
Trichloroaniline, 2,4,6-	634-93-5	2.0E+02	PHYSPROP	5.5E-05	1.3E-06	PHYSPROP	4.4E-03	PHYSPROP	7.9E+01	PHYSPROP			5.6E-02	6.6E-06	WATER9 (U.S. EPA, 2001)			4.4E+03	EPI	3.5E+00	PHYSPROP	4.0E+01	PHYSPROP	1.5E-01	1.3E+00	3.2E+00	2.7E-02	EPI
Trichlorobenzene, 1,2,3-	87-61-6	1.8E+02	PHYSPROP	5.1E-02	1.3E-03	PHYSPROP	2.1E-01	PHYSPROP	5.4E+01	PHYSPROP	1.5E+00	CRC	4.0E-02	8.4E-06	WATER9 (U.S. EPA, 2001)			1.4E+03	EPI	4.1E+00	PHYSPROP	1.8E+01	PHYSPROP	3.8E-01	1.1E+00	2.6E+00	7.4E-02	EPI
Trichlorobenzene, 1,2,4-	120-82-1	1.8E+02	PHYSPROP	5.8E-02	1.4E-03	PHYSPROP	4.6E-01	PHYSPROP	1.7E+01	PHYSPROP	1.5E+00	CRC	4.0E-02	8.4E-06	WATER9 (U.S. EPA, 2001)			1.4E+03	EPI	4.0E+00	PHYSPROP	4.9E+01	PHYSPROP	3.7E-01	1.1E+00	2.6E+00	7.1E-02	EPI
Trichloroethane, 1,1,1-	71-55-6	1.3E+02	PHYSPROP	7.0E-01	1.7E-02	PHYSPROP	1.2E+02	PHYSPROP	-3.0E+01	PHYSPROP	1.3E+00	CRC	6.5E-02	9.6E-06	WATER9 (U.S. EPA, 2001)			4.4E+01	EPI	2.5E+00	PHYSPROP	1.3E+03	PHYSPROP	5.6E-02	5.9E-01	1.4E+00	1.3E-02	EPI
Trichloroethane, 1,1,2-	79-00-5	1.3E+02	PHYSPROP	3.4E-02	8.2E-04	PHYSPROP	2.3E+01	PHYSPROP	-3.7E+01	PHYSPROP	1.4E+00	CRC	6.7E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			6.1E+01	EPI	1.9E+00	PHYSPROP	4.6E+03	PHYSPROP	2.2E-02	5.9E-01	1.4E+00	5.0E-03	EPI
Trichloroethylene	79-01-6	1.3E+02	PHYSPROP	4.0E-01	9.9E-03	PHYSPROP	6.9E+01	PHYSPROP	-8.5E+01	PHYSPROP	1.5E+00	CRC	6.9E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			6.1E+01	EPI	2.4E+00	PHYSPROP	1.3E+03	PHYSPROP	5.1E-02	5.7E-01	1.4E+00	1.2E-02	EPI
Trichlorofluoromethane	75-69-4	1.4E+02	PHYSPROP	4.0E+00	9.7E-02	PHYSPROP	8.0E+02	PHYSPROP	-1.1E+02	PHYSPROP	1.5E+00	CRC	6.5E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			4.4E+01	EPI	2.5E+00	PHYSPROP	1.1E+03	PHYSPROP	5.7E-02	6.2E-01	1.5E+00	1.3E-02	EPI
Trichlorophenol, 2,4,5-	95-95-4	2.0E+02	PHYSPROP	6.6E-05	1.6E-06	EPI	7.5E-03	EPI	6.9E+01	PHYSPROP	1.5E+00	PERRY	3.1E-02	8.1E-06	WATER9 (U.S. EPA, 2001)			1.6E+03	SSL	3.7E+00	PHYSPROP	1.2E+03	PHYSPROP	2.0E-01	1.3E+00	3.2E+00	3.6E-02	EPI
Trichlorophenol, 2,4,6-	88-06-2	2.0E+02	PHYSPROP	1.1E-04	2.6E-06	EPI	8.0E-03	EPI	6.9E+01	PHYSPROP	1.5E+00	CRC	3.1E-02	8.1E-06	WATER9 (U.S. EPA, 2001)			3.8E+02	SSL	3.7E+00	PHYSPROP	8.0E+02	PHYSPROP	1.9E-01	1.3E+00	3.2E+00	3.5E-02	EPI
Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5	2.6E+02	PHYSPROP	3.5E-07	8.7E-09	PHYSPROP	3.8E-05	EPI	1.5E+02	PHYSPROP	1.8E+00	PubChem	2.9E-02	7.8E-06	WATER9 (U.S. EPA, 2001)			1.1E+02	EPI	3.3E+00	PHYSPROP	2.8E+02	PHYSPROP	5.6E-02	2.8E+00	6.8E+00	9.1E-03	EPI
Trichlorophenoxypropionic acid, -2,4,5	93-72-1	2.7E+02	PHYSPROP	3.7E-07	9.1E-09	PHYSPROP	1.0E-05	PHYSPROP	1.8E+02	PHYSPROP	1.2E+00	PubChem	2.3E-02	5.9E-06	WATER9 (U.S. EPA, 2001)			1.8E+02	EPI	3.8E+00	PHYSPROP	7.1E+01	PHYSPROP	1.0E-01	3.4E+00	8.2E+00	1.6E-02	EPI
Trichloropropane, 1,1,2-	598-77-6	1.5E+02	PHYSPROP	1.3E-02	3.2E-04	EPI	3.1E+00	PHYSPROP	-6.5E+01	EPI	1.4E+00	CRC	5.7E-02	9.2E-06	WATER9 (U.S. EPA, 2001)			9.5E+01	EPI	2.4E+00	PHYSPROP	1.9E+03	PHYSPROP	4.5E-02	7.0E-01	1.7E+00	9.6E-03	EPI
Trichlorocorooane, 1.2.3-	96-18-4	1.5E+02	PHYSPROP	1.4E-02	3.4E-04	PHYSPROP	3.7E+00	PHYSPROP	-1.5E+01	PHYSPROP	1.4E+00	CRC	5.7E-02	9.2E-06	WATER9 (U.S. EPA, 2001)			1.2E+02	EPI	2.3E+00	PHYSPROP	1.8E+03	PHYSPROP	3.5E-02	7.0E-01	1.7E+00	7.5E-03	EPI
Trichloropropene, 1,2,3-	96-19-5	1.5E+02	PHYSPROP	7.2E-01	1.8E-02	PHYSPROP	4.4E+00	PHYSPROP	-5.6E+01	EPI	1.4E+00	CRC	5.9E-02	9.4E-06	WATER9 (U.S. EPA, 2001)			1.2E+02	EPI	2.8E+00	PHYSPROP	3.3E+02	PHYSPROP	7.8E-02	6.9E-01	1.6E+00	1.7E-02	EPI
Tricresyl Phosphate (TCP)	1330-78-5	3.7E+02	PHYSPROP	3.3E-05	8.1E-07	EPI	6.0E-07	EPI	-3.3E+01	PHYSPROP	1.2E+00	Yaws	1.9E-02	4.8E-06	WATER9 (U.S. EPA, 2001)			4.7E+04	EPI	5.1E+00	PHYSPROP	3.6E-01	PHYSPROP	2.5E-01	1.2E+01	2.9E+01	3.3E-02	EPI
Tridiphane	58138-08-2	3.2E+02	PHYSPROP	1.7E-05	4.1E-07	PHYSPROP	3.9E-04	PHYSPROP	4.3E+01	PHYSPROP			4.1E-02	4.7E-06	WATER9 (U.S. EPA, 2001)			3.4E+03	EPI	5.2E+00	PHYSPROP	1.1E+00	PHYSPROP	4.7E-01	6.6E+00	1.6E+01	6.9E-02	EPI
Triethylamine	121-44-8	1.0E+02	PHYSPROP	6.1E-03	1.5E-04	PHYSPROP	5.7E+01	PHYSPROP	-1.1E+02	PHYSPROP	7.3E-01	CRC	6.6E-02	7.9E-06	WATER9 (U.S. EPA, 2001)			5.1E+01	EPI	1.5E+00	PHYSPROP	6.9E+04	PHYSPROP	1.5E-02	3.9E-01	9.3E-01	3.9E-03	EPI
Triethylene Glycol	112-27-6	1.5E+02	PHYSPROP	1.3E-09	3.2E-11	PHYSPROP	1.3E-03	PHYSPROP	-7.0E+00	PHYSPROP	1.1E+00	CRC	5.1E-02	8.1E-06	WATER9 (U.S. EPA, 2001)			1.0E+01	EPI	-1.8E+00	PHYSPROP	1.0E+06	PHYSPROP	7.3E-05	7.3E-01	1.8E+00	1.6E-05	EPI
Trifluoroethane, 1,1,1-	420-46-2	8.4E+01	PHYSPROP	3.1E+01	7.7E-01	PHYSPROP	9.5E+03	PHYSPROP	-1.1E+02	PHYSPROP			9.9E-02	1.2E-05	WATER9 (U.S. EPA, 2001)			4.4E+01	EPI	1.7E+00	PHYSPROP	7.6E+02	PHYSPROP	2.7E-02	3.1E-01	7.5E-01	7.6E-03	EPI
Trifluralin	1582-09-8	3.4E+02	PHYSPROP	4.2E-03	1.0E-04	PHYSPROP	4.6E-05	PHYSPROP	4.9E+01	PHYSPROP	1.4E+00	PubChem	2.2E-02	5.6E-06	WATER9 (U.S. EPA, 2001)			1.6E+04	EPI	5.3E+00	PHYSPROP	1.8E-01	PHYSPROP	5.1E-01	7.9E+00	1.9E+01	7.3E-02	EPI
Trimethyl Phosphate	512-56-1	1.4E+02	PHYSPROP	2.9E-07	7.2E-09	PHYSPROP	8.5E-01	EPI	-4.6E+01	PHYSPROP	1.2E+00	CRC	5.8E-02	8.8E-06	WATER9 (U.S. EPA, 2001)			1.1E+01	EPI	-6.5E-01	PHYSPROP	5.0E+05	PHYSPROP	4.3E-04	6.4E-01	1.5E+00	9.5E-05	EPI
Trimethylbenzene, 1,2,3-	526-73-8	1.2E+02	PHYSPROP	1.8E-01	4.4E-03	PHYSPROP	1.7E+00	PHYSPROP	-2.5E+01	PHYSPROP	8.9E-01	CRC	6.1E-02	8.0E-06	WATER9 (U.S. EPA, 2001)			6.3E+02	EPI	3.7E+00	PHYSPROP	7.5E+01	PHYSPROP	3.8E-01	5.0E-01	1.2E+00	9.0E-02	EPI
Trimethylbenzene, 1,2,4-	95-63-6	1.2E+02	PHYSPROP	2.5E-01	6.2E-03	PHYSPROP	2.1E+00	PHYSPROP	-4.4E+01	PHYSPROP	8.8E-01	CRC	6.1E-02	7.9E-06	WATER9 (U.S. EPA, 2001)			6.1E+02	EPI	3.6E+00	PHYSPROP	5.7E+01	PHYSPROP	3.6E-01	5.0E-01	1.2E+00	8.6E-02	EPI
Trimethylbenzene, 1,3,5-	108-67-8	1.2E+02	PHYSPROP	3.6E-01	8.8E-03	PHYSPROP	2.5E+00	PHYSPROP	-4.5E+01	PHYSPROP	8.6E-01	CRC	6.0E-02	7.8E-06	WATER9 (U.S. EPA, 2001)			6.0E+02	EPI	3.4E+00	PHYSPROP	4.8E+01	PHYSPROP	2.6E-01	5.0E-01	1.2E+00	6.2E-02	EPI
Trimethylpentene, 2,4,4-	25167-70-8	1.1E+02	PHYSPROP	3.0E+01	7.5E-01	PHYSPROP	7.1E+01	PHYSPROP	-8.4E+01	EPI	7.2E-01	PubChem	6.0E-02	7.3E-06	WATER9 (U.S. EPA, 2001)			2.4E+02	EPI	4.1E+00	PHYSPROP	4.0E+00	PHYSPROP	1.6E+00	4.5E-01	1.8E+00	3.9E-01	EPI
Trinitrobenzene, 1,3,5-	99-35-4	2.1E+02	PHYSPROP	2.7E-07	6.5E-09	EPI	6.4E-06	EPI	1.2E+02	PHYSPROP	1.5E+00	CRC	2.9E-02	7.7E-06	WATER9 (U.S. EPA, 2001)			1.7E+03	EPI	1.2E+00	PHYSPROP	2.8E+02	PHYSPROP	3.4E-03	1.6E+00	3.9E+00	6.1E-04	EPI
Trinitrotoluene, 2,4,6-	118-96-7	2.3E+02	PHYSPROP	8.5E-07	2.1E-08	EPI	8.0E-06	PHYSPROP	8.0E+01	PHYSPROP	1.7E+00	CRC	3.0E-02	7.9E-06	WATER9 (U.S. EPA, 2001)			2.8E+03	EPI	1.6E+00	PHYSPROP	1.2E+02	PHYSPROP	5.6E-03	2.0E+00	4.7E+00	9.6E-04	EPI
Triphenylphosphine Oxide	791-28-6	2.8E+02	PHYSPROP	2.2E-08	5.3E-10	PHYSPROP	2.6E-09	EPI	1.6E+02	PHYSPROP	1.2E+00	CRC	2.3E-02	5.8E-06	WATER9 (U.S. EPA, 2001)			2.0E+03	EPI	2.8E+00	PHYSPROP	6.3E+01	PHYSPROP	2.1E-02	3.8E+00	9.1E+00	3.3E-03	EPI
Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8	4.3E+02	PHYSPROP	1.1E-07	2.6E-09	PHYSPROP	7.4E-08	PHYSPROP	2.7E+01	PHYSPROP			3.3E-02	3.9E-06	WATER9 (U.S. EPA, 2001)			1.1E+04	EPI	3.7E+00	PHYSPROP	7.0E+00	PHYSPROP	1.3E-02	2.7E+01	6.5E+01	1.6E-03	EPI
Tris(1-chloro-2-propyl)phosphate	13674-84-5	3.3E+02	PHYSPROP	2.4E-06	6.0E-08	PHYSPROP	2.0E-05	PHYSPROP	-4.0E+01	PHYSPROP			4.0E-02	4.7E-06	WATER9 (U.S. EPA, 2001)			1.6E+03	EPI	2.6E+00	PHYSPROP	1.2E+03	PHYSPROP	8.4E-03	7.2E+00	1.7E+01	1.2E-03	EPI
Tris(2,3-dibromopropyl)phosphate	126-72-7	7.0E+02	PHYSPROP	8.9E-04	2.2E-05	EPI	1.9E-04	PHYSPROP	5.5E+00	PHYSPROP	2.3E+00	PubChem	1.9E-02	4.9E-06	WATER9 (U.S. EPA, 2001)			9.7E+03	EPI	4.3E+00	PHYSPROP	8.0E+00	PHYSPROP	1.4E-03	8.5E+02	2.0E+03	1.4E-04	EPI
Tris(2-chloroethyl)phosphate	115-96-8	2.9E+02	PHYSPROP	1.3E-04	3.3E-06	EPI	6.1E-02	PHYSPROP	-5.5E+01	PHYSPROP	1.4E+00	CRC	2.4E-02	6.2E-06	WATER9 (U.S. EPA, 2001)			3.9E+02	EPI	1.4E+00	PHYSPROP	7.0E+03	PHYSPROP	2.3E-03	4.2E+00	1.0E+01	3.6E-04	EPI
Tris(2-ethylhexyl)phosphate	78-42-2	4.3E+02	PHYSPROP	3.2E-06	7.9E-08	EPI	8.3E-08	PHYSPROP	-7.4E+01	PHYSPROP	9.9E-01	CRC	1.6E-02	3.9E-06	WATER9 (U.S. EPA, 2001)			2.5E+06	EPI	9.5E+00	PHYSPROP	6.0E-01	PHYSPROP	9.3E+01	2.9E+01	1.3E+02	1.2E+01	EPI
Tungsten	7440-33-7	1.8E+02	PHYSPROP				0.0E+00	NIOSH	3.4E+03	PHYSPROP	1.9E+01	CRC				1.5E+02	BAES							5.2E-03	1.1E+00	2.7E+00	1.0E-03	RAGSE
Uranium	7440-61-1	2.4E+02	CRC				0.0E+00	NIOSH	1.1E+03	CRC	1.9E+01	CRC				4.5E+02	BAES							5.9E-03	2.3E+00	5.4E+00	1.0E-03	RAGSE
Urethane	51-79-6	8.9E+01	PHYSPROP	2.6E-06	6.4E-08	EPI	2.6E-01	EPI	4.9E+01	PHYSPROP	9.9E-01	CRC	8.5E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.2E+01	EPI	-1.5E-01	PHYSPROP	4.8E+05	PHYSPROP	1.4E-03	3.3E-01	8.0E-01	3.9E-04	EPI
Vanadium Pentoxide	1314-62-1	1.8E+02	EPI				0.0E+00	NIOSH	6.8E+02	CRC	3.4E+00	CRC										7.0E+02	CRC	5.2E-03	1.1E+00	2.6E+00	1.0E-03	RAGSE
Vanadium and Compounds	7440-62-2	5.1E+01	EPI						1.9E+03	CRC	6.0E+00	CRC				1.0E+03	SSL							2.7E-03	2.0E-01	4.9E-01	1.0E-03	RAGSE
Vernolate	1929-77-7	2.0E+02	PHYSPROP	1.3E-03	3.1E-05	EPI	1.0E-02	PHYSPROP	7.1E+01	EPI	9.5E-01	CRC	2.4E-02	6.1E-06	WATER9 (U.S. EPA, 2001)			3.0E+02	EPI	3.8E+00	PHYSPROP	9.0E+01	PHYSPROP	2.2E-01	1.4E+00	3.5E+00	4.0E-02	EPI
Vinclozolin	50471-44-8	2.9E+02	PHYSPROP	7.1E-07	1.7E-08	EPI	1.2E-07	PHYSPROP	1.1E+02	PHYSPROP	1.5E+00	CRC	2.5E-02	6.5E-06	WATER9 (U.S. EPA, 2001)			2.8E+02	EPI	3.1E+00	PHYSPROP	2.6E+00	PHYSPROP	2.9E-02	4.2E+00	1.0E+01	4.5E-03	EPI
Vinyl Acetate	108-05-4	8.6E+01	PHYSPROP	2.1E-02	5.1E-04	EPI	9.0E+01	PHYSPROP	-9.3E+01	PHYSPROP	9.3E-01	CRC	8.5E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			5.6E+00	EPI	7.3E-01	PHYSPROP	2.0E+04	PHYSPROP	5.6E-03	3.2E-01	7.7E-01	1.6E-03	EPI
Vinyl Bromide	593-60-2	1.1E+02	PHYSPROP	5.0E-01	1.2E-02	PHYSPROP	1.0E+03	PHYSPROP	-1.4E+02	PHYSPROP	1.5E+00	CRC	8.6E-02	1.2E-05	WATER9 (U.S. EPA, 2001)			2.2E+01	EPI	1.6E+00	PHYSPROP	7.6E+03	PHYSPROP	1.7E-02	4.2E-01	1.0E+00	4.4E-03	EPI
Vinyl Chloride	75-01-4	6.2E+01	PHYSPROP	1.1E+00	2.8E-02	PHYSPROP	3.0E+03	EPI	-1.5E+02	PHYSPROP	9.1E-01	CRC	1.1E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			2.2E+01	EPI	1.4E+00	CRC	8.8E+03	PHYSPROP	2.5E-02	2.4E-01	5.7E-01	8.4E-03	EPI
Warfarin	81-81-2	3.1E+02	PHYSPROP	1.1E-07	2.8E-09	EPI	1.2E-07	PHYSPROP	1.6E+02	PHYSPROP			4.2E-02	4.9E-06	WATER9 (U.S. EPA, 2001)			4.3E+02	EPI	2.7E+00	PHYSPROP	1.7E+01	PHYSPROP	1.2E-02	5.6E+00	1.3E+01	1.8E-03	EPI
Xylene, m-	108-38-3	1.1E+02	PHYSPROP	2.9E-01	7.2E-03	PHYSPROP	8.3E+00	PHYSPROP	-4.8E+01	PHYSPROP	8.6E-01	CRC	6.9E-02	8.5E-06	WATER9 (U.S. EPA, 2001)			3.8E+02	EPI	3.2E+00	PHYSPROP	1.6E+02	PHYSPROP	2.1E-01	4.1E-01	9.9E-01	5.3E-02	EPI
Xylene, o-	95-47-6	1.1E+02	PHYSPROP	2.1E-01	5.2E-03	PHYSPROP	6.6E+00	PHYSPROP	-2.5E+01	PHYSPROP	8.8E-01	CRC	6.9E-02	8.6E-06	WATER9 (U.S. EPA, 2001)			3.8E+02	EPI	3.1E+00	PHYSPROP	1.8E+02	PHYSPROP	1.9E-01	4.1E-01	9.9E-01	4.7E-02	EPI
Xylene, p-	106-42-3	1.1E+02	PHYSPROP	2.8E-01	6.9E-03	PHYSPROP	8.8E+00	PHYSPROP	1.3E+01	PHYSPROP	8.6E-01	CRC	6.8E-02	8.4E-06	WATER9 (U.S. EPA, 2001)			3.8E+02	EPI	3.2E+00	PHYSPROP	1.6E+02	PHYSPROP	2.0E-01	4.1E-01	9.9E-01	4.9E-02	EPI
Xylenes	1330-20-7	1.1E+02	PHYSPROP	2.7E-01	6.6E-03	PHYSPROP	8.0E+00	PHYSPROP	-2.5E+01	EPI	8.6E-01	ATSDR Profile	6.9E-02	8.5E-06	WATER9 (U.S. EPA, 2001)			3.8E+02	EPI	3.2E+00	PHYSPROP	1.1E+02	PHYSPROP	2.0E-01	4.1E-01	9.9E-01	5.0E-02	EPI
Zinc Phosphide	1314-84-7	2.6E+02	CRC						4.2E+02	PERRY	4.6E+00	CRC												3.7E-03	2.9E+00	7.0E+00	6.0E-04	RAGSE
Zinc and Compounds	7440-66-6	6.5E+01	PHYSPROP						4.2E+02	PHYSPROP	7.1E+00	CRC				6.2E+01	SSL							1.9E-03	2.4E-01	5.9E-01	6.0E-04	RAGSE
Zineb	12122-67-7	2.8E+02	PHYSPROP	1.1E-07	2.7E-09	PHYSPROP	7.5E-08	PHYSPROP	1.6E+02	EPI			4.5E-02	5.2E-06	WATER9 (U.S. EPA, 2001)			1.3E+03	EPI	1.3E+00	PHYSPROP	1.0E+01	PHYSPROP	2.1E-03	3.7E+00	8.8E+00	3.3E-04	EPI
Zirconium	7440-67-7	9.1E+01	EPI				0.0E+00	NIOSH	1.9E+03	CRC	6.5E+00	CRC				3.0E+03	BAES							3.7E-03	3.4E-01	8.2E-01	1.0E-03	RAGSE