

Regional Screening Level (RSL) Summary Table (TR-1E-6, HQ=1) June 2015 (revised)

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Screening Levels								Protection of Ground Water SSLs								
SFO (mg/kg-day) ⁻¹	Key	IUR (ug/m ³) ⁻¹	Key	RfD _o (mg/kg-day)	Key	RfC ₁ (mg/m ³)	Key	muta-gen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)	
1.8E-02	C	5.1E-06	C	1.5E-01	I					1	0.1	ALAR	1596-84-5	3.0E+01	c	1.3E+02	c	5.5E-01	c	2.4E+00	c	4.3E+00	c		9.5E-04	c		
8.7E-03	I			4.0E-03	I					1	0.1	Acephate	30560-19-1	6.2E+01	c**	2.6E+02	c*					8.9E+00	c**		2.0E-03	c**		
		2.2E-06	I			9.0E-03	I	V		1	1.1E+05	Acetaldehyde	75-07-0	1.1E+01	c**	4.9E+01	c**	1.3E+00	c**	5.6E+00	c**	2.6E+00	c**		5.2E-04	c**		
				2.0E-02	I					1	0.1	Acetochlor	34256-82-1	1.3E+03	n	1.6E+04	n					3.5E+02	n		2.8E-01	n		
				9.0E-01	I	3.1E+01	A	V		1	1.1E+05	Acetone	67-64-1	6.1E+04	n	6.7E+05	nms	3.2E+04	n	1.4E+05	n	1.4E+04	n		2.9E+00	n		
				2.0E-03	X	V				1	1.1E+05	Acetone Cyanohydrin	75-86-5	5.0E+01	n	2.1E+02	n	2.1E+00	n	8.8E+00	n	4.2E+00	n		8.4E-04	n		
				6.0E-02	I	V				1	1.3E+05	Acetonitrile	75-05-8	8.1E+02	n	3.4E+03	n	6.3E+01	n	2.6E+02	n	1.3E+02	n		2.6E-02	n		
3.8E+00	C	1.3E-03	C	1.0E-01	I			V		1	2.5E+03	Acetophenone	98-86-2	7.8E+03	ns	1.2E+05	nms					1.9E+03	n		5.8E-01	n		
										1	0.1	Acetylaminofluorene, 2-	53-96-3	1.4E-01	c	6.0E-01	c	2.2E-03	c	9.4E-03	c	1.6E-02	c		7.2E-05	c		
				5.0E-04	I	2.0E-05	I	V		1	2.3E+04	Acrolein	107-02-8	1.4E-01	n	6.0E-01	n	2.1E-02	n	8.8E-02	n	4.2E-02	n		8.4E-06	n		
5.0E-01	I	1.0E-04	I	2.0E-03	I	6.0E-03	I		M	1	0.1	Acrylamide	79-06-1	2.4E-01	c	4.6E+00	c	1.0E-02	c	1.2E-01	c	5.0E-02	c		1.1E-05	c		
				5.0E-01	I	1.0E-03	I	V		1	1.1E+05	Acrylic Acid	79-10-7	9.9E+01	n	4.2E+02	n	1.0E+00	n	4.4E+00	n	2.1E+00	n		4.2E-04	n		
5.4E-01	I	6.8E-05	I	4.0E-02	A	2.0E-03	I	V		1	1.1E+04	Acrylonitrile	107-13-1	2.5E-01	c*	1.1E+00	c*	4.1E-02	c*	1.8E-01	c*	5.2E-02	c*		1.1E-05	c*		
5.6E-02	C			1.0E-02	I	6.0E-03	P			1	0.1	Adiponitrile	111-69-3	8.5E+06	nm	3.6E+07	nm	6.3E+00	n	2.6E+01	n							
										1	0.1	Alachlor	15972-80-8	9.7E+00	c*	4.1E+01	c*					1.0E+00	c	2.0E+00		8.6E-04	c	1.6E-03
				1.0E-03	I					1	0.1	Aldicarb	116-06-3	6.3E+01	n	8.2E+02	n					2.0E+01	n	3.0E+00	4.9E-03	n	7.5E-04	
				1.0E-03	I					1	0.1	Aldicarb Sulfone	1646-88-4	6.3E+01	n	8.2E+02	n					2.0E+01	n	3.0E+00	4.4E-03	n	4.4E-04	
										1	0.1	Aldicarb sulfoxide	1646-87-3	6.3E+01	n	8.2E+02	n					2.0E+01	n	4.0E+00	4.4E-03	n	8.8E-04	
1.7E+01	I	4.9E-03	I	3.0E-05	I			V		1		Aldrin	309-00-2	3.9E-02	c*	1.8E-01	c	5.7E-04	c	2.5E-03	c	9.2E-04	c		1.5E-04	c		
				2.5E-01	I					1	0.1	Allyl	74223-64-6	1.6E+04	n	2.1E+05	nm					4.9E+03	n		1.9E+00	n		
				5.0E-03	I	1.0E-04	X	V		1	1.1E+05	Allyl Alcohol	107-18-6	3.5E+00	n	1.5E+01	n	1.0E-01	n	4.4E-01	n	2.1E-01	n		4.2E-05	n		
2.1E-02	C	6.0E-06	C	1.0E-03	I	V				1	1.4E+03	Allyl Chloride	77-05-1	7.2E-01	c**	3.2E+00	c**	4.7E-01	c**	2.0E+00	c**	7.3E-01	c**		2.3E-04	c**		
				1.0E+00	P	5.0E-03	P			1		Aluminum	7429-90-5	7.7E+04	n	1.1E+06	nm	5.2E+00	n	2.2E+01	n	2.0E+04	n		3.0E+04	n		
				4.0E-04	I					1		Aluminum Phosphide	20889-78-8	3.1E+01	n	4.7E+02	n					8.0E+00	n					
				3.0E-04	I					1	0.1	Amidrin	61485-29-4	1.9E+01	n	2.5E+02	n					5.9E+00	n		2.1E+03	n		
2.1E+01	C	6.0E-03	C	9.0E-03	I					1	0.1	Ametryn	834-12-8	5.7E+02	n	7.4E+03	n					1.5E+02	n		1.6E-01	n		
										1	0.1	Amnophenyl #	92-67-1	2.6E-02	c	1.1E-01	c	4.7E-04	c	2.0E-03	c	3.0E-03	c		1.5E-05	c		
				8.0E-02	P					1	0.1	Aminophenol, m-	591-27-5	5.1E+03	n	6.6E+04	n					1.6E+03	n		6.1E-01	n		
				2.0E-02	P					1	0.1	Aminophenol, p-	123-30-8	1.3E+03	n	1.6E+04	n					4.0E+02	n		1.5E-01	n		
				2.5E-03	I					1	0.1	Amtraz	33089-61-1	1.6E+02	n	2.1E+03	n					8.2E+00	n		4.2E+00	n		
				1.0E-01	I	V				1		Ammonia	7664-41-7					1.0E+02	n	4.4E+02	n							
				2.0E-01	I			X	V	1	1.4E+04	Ammonium Sulfamate	7773-06-0	1.6E+04	n	2.3E+05	nm					4.0E+03	n					
				3.0E-03	X	V				1		Amyl Alcohol, tert-	79-85-4	8.2E+01	n	3.4E+02	n	3.1E+00	n	1.3E+01	n	6.3E+00	n		1.3E-03	n		
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I			1	0.1	Aniline	62-53-3	9.5E+01	c**	4.0E+02	c*	1.0E+00	n	4.4E+00	n	1.3E+01	c*		4.6E-03	c*		
4.0E-02	P			2.0E-03	X					1	0.1	Anthraquinone, 9,10-	84-65-1	1.4E+01	c**	5.7E+01	c*					1.4E+00	c*		1.4E-02	c*		
				4.0E-04	I			0.15				Antimony (metallic)	7440-36-0	3.1E+01	n	4.7E+02	n					7.8E+00	n	6.0E+00	3.5E-01	n	2.7E-01	
				5.0E-04	H			0.15				Antimony Pentoxide	1314-80-9	3.9E+01	n	5.8E+02	n					9.7E+00	n					
				9.0E-04	H			0.15				Antimony Potassium Tartrate	11071-15-1	7.0E+01	n	1.1E+03	n					1.8E+01	n					
				4.0E-04	H			0.15				Antimony Tetroxide	1332-81-6	3.1E+01	n	4.7E+02	n					7.8E+00	n					
						2.0E-04	I		0.15			Antimony Trioxide	1309-64-4	2.8E+05	nm	1.2E+06	nm	2.1E-01	n	8.8E-01	n							
2.5E-02	I	7.1E-06	I	1.3E-02	I					1	0.1	Apollo	74115-24-5	8.2E+02	n	1.1E+04	n					2.3E+02	n		1.4E+01	n		
				5.0E-02	H					1	0.1	Aramite	140-57-8	2.2E+01	c	9.2E+01	c	4.0E-01	c	1.7E+00	c	1.3E+00	c		1.5E-02	c		
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C			1	0.03	Arsenic, Inorganic	7440-38-2	6.8E-01	cR	3.0E+00	cR	6.5E-04	c*	2.9E-03	c*	5.2E-02	c	1.0E+01	1.5E-03	c	2.9E-01	
				3.5E-06	C	5.0E-05	I			1		Arsine	7784-42-1	2.7E-01	n	4.1E+00	n	5.2E-02	n	2.2E-01	n	7.0E-02	n					
				9.0E-03	I					1	0.1	Assure	76578-14-8	5.7E+02	n	7.4E+03	n					1.2E+02	n		1.9E+00	n		
				5.0E-02	I					1	0.1	Asulam	3337-71-1	3.2E+03	n	4.1E+04	n					1.0E+03	n		2.6E-01	n		
2.3E-01	C			3.5E-02	I					1	0.1	Atrazine	1912-24-9	2.4E+00	c	1.0E+01	c					3.0E-01	c	3.0E+00	1.9E-04	c	1.9E-03	
8.8E-01	C	2.5E-04	C							1	0.1	Auramine	492-80-8	6.2E-01	c	2.6E+00	c	1.1E-02	c	4.9E-02	c	6.6E-02	c		6.0E-04	c		
				4.0E-04	I					1	0.1	Avermectin B1	65195-55-3	2.5E+01	n	3.3E+02	n					8.0E+00	n		1.4E+01	n		
1.1E-01	I	3.1E-05	I	1.0E+00	P	7.0E-06	P	V		1	0.1	Azobenzene	103-33-3	5.6E+00	c	2.6E+01	c	9.1E-02	c	4.0E-01	c	1.2E-01	c		9.2E-04	c		
										1	0.1	Azodicarbonamide	123-77-3	8.6E+03	n	4.0E+04	n					3.1E-02	n	2.0E+04	n	6.8E+00	n	

Regional Screening Level (RSL) Summary Table (TR=1E-6, HQ=1) June 2015 (revised)

Toxicity and Chemical-specific Information													Contaminant		Screening Levels								Protection of Ground Water SSLs							
SFO (mg/kg-day) ⁻¹	Key	IUR (ug/m ³ -y) ⁻¹	Key	RfD _o (mg/kg-day)	Key	RfC _i (mg/m ³ -y)	Key	Vo	muta- gen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)		
1.3E+01	I							V				3.2E+02	Benzotrifluoride	98-07-7	5.3E-02	c	2.5E-01	c					2.9E-03	c		6.5E-08	c			
1.7E-01	I	4.9E-05 2.4E-03	C	1.0E-01 2.0E-03	P	1.0E-03 2.0E-05	P	V				1.5E+03	Benzyl Alcohol Benzyl Chloride Beryllium and compounds	100-51-6 100-44-7 7440-41-7	6.3E+00 1.1E+00 1.6E+02	n c* n	8.2E+04 4.8E+00 2.3E+03	n c* n	5.7E-02 4.8E+00 1.2E-03	c* c* c*	2.5E-01 5.1E-03	c*	2.0E+03 8.9E-02 2.5E+01	n c* c*		4.8E-01 9.7E-05 1.9E+01	n c* n		3.2E+00	
8.0E-03	I			5.0E-01	I	4.0E-04	X	V				1.0E+03	Bidrin Bifenox Biphenrin	141-66-2 42576-02-3 82657-04-3	6.3E+00 5.7E+02 9.5E+02	n n n	8.2E+01 7.4E+03 1.2E+04	n n n							2.0E+00 1.0E+02 3.0E+02	n n n		4.7E-04 7.6E-01 1.4E+03	n n n	
7.0E-02	H	1.0E-05	H	4.0E-02	I			V				1.0E+03	Biphenyl, 1,1'- Bis(2-chloro-1-methylethyl) ether Bis(2-chloroethoxy)methane	92-52-4 108-60-1 111-91-1	4.7E+01 4.9E+00 1.9E+02	n c n	2.0E+02 2.2E+01 2.5E+03	n c n	4.2E-01 2.8E-01	n c	1.8E+00 1.2E+00	n	8.3E-01 3.6E-01 5.9E+01	n n n		8.7E-03 1.3E-04 1.3E-02	n c n			
1.1E+00	I	3.3E-04	I					V				5.1E+03	Bis(2-chloroethyl)ether Bis(chloromethyl)ether Bisphenol A	111-44-4 542-88-1 80-05-7	2.3E-01 8.3E-05 3.2E+03	c n n	1.0E+00 3.6E-04 4.1E+04	c c n	8.5E-03 3.6E-04 4.5E-05	c c c	3.7E-02 2.0E-04	c c	1.4E-02 7.2E-05 7.7E+02	c c n		3.6E-06 1.7E-08 5.8E+01	c c n			
2.2E+02	I	6.2E-02	I	5.0E-02	I			V				0.1	Boron And Borates Only Boron Trichloride Boron Trifluoride	7440-42-8 10294-34-5 7637-07-2	1.6E+04 1.6E+05 3.1E+03	n nm n	2.3E+05 2.3E+06 4.7E+04	nm nm n	2.1E+01 2.1E+01 1.4E+01	n n n	8.8E+01 8.8E+01 5.7E+01	n n n	4.0E+03 4.2E+01 2.6E+01	n n n		1.3E+01 n n	n n n			
7.0E-01	I			4.0E-03	I			V				2.4E+03	Bromate Bromo-2-chloroethane, 1- Bromobenzene	15541-45-4 107-04-0 108-86-1	9.9E-01 2.6E-02 2.9E+02	c c ns	4.7E+00 1.1E-01 1.8E+03	c c ns	4.7E+00 4.7E-03 6.3E+01	c c n	2.0E-02 2.0E+02	c c	7.4E-03 6.2E+01	c c	1.0E+01	8.5E-04 2.1E-06 4.2E-02	c c n	7.7E-02		
2.0E+00	X	6.0E-04	X	8.0E-03	I	6.0E-02	I	V				6.8E+02	Bromochloromethane Bromodichloromethane Bromotorm	74-97-5 93-31-2 75-25-2	1.5E+02 2.9E-01 1.9E+01	n c* c*	6.3E+02 1.3E+00 8.6E+01	n c c	4.2E+01 7.6E-02 2.6E+00	n c c	1.8E+02 3.3E-01 1.1E+01	n c c	8.3E+01 1.3E-01 3.3E+00	n c c		2.1E-02 3.6E-05 8.7E-04	n c c	2.2E-02	2.1E-02	
6.2E-02	I	3.7E-05	C	2.0E-02	I			V				4.0E+03	Bromomethane Bromophos Bromoxynil	74-83-9 2104-96-3 1689-84-5	8.8E+00 3.9E+02 1.3E+03	n n n	3.0E+01 5.8E+03 1.6E+04	n n n	5.2E+00 n n	n n n	2.2E+01 n n	n n n	7.5E+00 3.5E+01 3.3E+02	n n n		1.9E-03 1.5E-01 2.8E-01	n n n			
7.9E-03	I	1.1E-06	I	1.4E-03	I	5.0E-03	I	V				3.8E+03	Bromoxynil Octanoate Butadiene, 1,3- Butanol, N-	1689-99-2 106-99-0 71-36-3	1.6E+03 5.8E-02 7.8E+03	n c* ns	2.3E+04 2.6E-01 1.2E+05	n c* nms	9.4E-02 4.1E-01	c* c*	4.1E-01 c*	n	1.4E-02 1.8E-02 2.0E+03	c c c		1.2E+00 9.9E-06 4.1E-01	n c n			
1.9E-03	P			2.0E-01	I			V				2.1E+04	Butyl Benzyl Phthalate Butyl alcohol, sec- Butylate	85-68-7 78-92-2 2008-41-5	2.9E+02 1.3E+05 3.9E+03	c* nms n	1.2E+03 1.5E+06 5.8E+04	c nms n	1.2E+03 3.1E+04 n	c n n	1.3E+05 n	n	1.6E+02 2.4E+04 4.6E+02	c n n		2.3E-01 5.0E+00 4.5E-01	c n n			
2.0E-04	C	5.7E-08	C	3.0E-01	P			V				1.0E+01	Butylated hydroxyanisole Butylated hydroxytoluene Butylbenzene, n	25013-16-5 128-37-0 104-51-8	2.7E+03 1.5E-02 3.9E+03	c c ns	1.1E+04 6.4E+02 5.8E+04	c c ns	4.9E+01 6.4E+02	c c	2.2E+02 n	c	2.4E+02 3.3E+00 1.0E+03	c c c		4.5E-01 9.7E-02 3.2E+00	c c n			
3.6E-03	P			5.0E-02	P			V				1.1E+02	Butylbenzene, sec- Butylbenzene, tert- Cacodylic Acid	135-98-8 98106-6 75-60-5	7.8E+03 7.8E+03 1.3E+03	ns ns n	1.2E+05 1.2E+05 1.6E+04	nms nms n					2.0E+03 6.9E+02 4.0E+02	n n n		5.9E+00 1.6E+00 n	n n n			
5.0E-01	C	1.8E-03 1.8E-03 1.5E-01	I I C	1.0E-03 5.0E-04 2.0E-02	I I C	1.0E-05 1.0E-05 2.0E-04	A A C			0.025 0.05 0.025	0.001 0.001		Cadmium (Diet) Cadmium (Water) Calcium Chromate	7440-43-9 7440-43-9 13765-19-0	7.1E+01 3.0E-01	n c	9.8E+02 6.2E+00	n c			1.6E-03 6.8E-06	c** c	6.8E-03 8.2E-05	c** c	9.2E+00 4.1E-02	n c	5.0E+00	6.9E-01 c	n c	3.8E-01
1.5E-01	C	4.3E-05	C	5.0E-01	I	2.2E-03	C					0.1	Caprolactam Captafol Captan	105-60-2 2425-06-1 133-06-2	3.1E+04 3.6E+00 2.4E+02	n c* c*	4.0E+05 1.5E+01 1.0E+03	nm c* c	2.3E+00 6.5E-02 4.3E+00	n c* c	9.6E+00 2.9E-01 1.9E+01	n c* c*	9.9E+03 4.0E-01 3.1E+01	n c* c*		2.5E+00 7.1E-04 2.2E-02	n c* c*			
2.3E-03	C	6.8E-07	C	1.0E-01	I			V				0.1	Carbaryl Carbofuran Carbon Disulfide	63-25-2 1563-66-2 75-15-0	6.3E+03 3.2E+02 7.7E+02	n n ns	8.2E+04 4.1E+03 3.5E+03	n n ns	7.3E+02 n	n	3.1E+03 n	n	1.8E+03 9.4E+01 8.1E+02	n n n		1.7E+00 3.7E-02 2.4E-01	n n n	1.6E-02		
7.0E-02	I	6.0E-06	I	4.0E-03	I	1.0E-01	I	V				4.6E+02	Carbon Tetrachloride Carbosulfan Carboxin	56-23-5 55285-14-8 5234-68-4	6.5E-01 6.3E-02 6.3E+03	c n n	2.9E+00 8.2E+03 8.2E+04	c n n	4.7E-01 n	c	2.0E+00 n	c	4.5E-01 5.1E+01 1.9E+03	c n n		1.8E-04 1.2E+00 1.0E+00	c n n	1.9E-03		
4.0E-01	H			1.0E-01	I			V				0.1	Ceric oxide Chloral Hydrate Chloramben	1306-38-3 302-17-0 133-90-4	1.3E+06 7.8E+03 9.5E+02	nm nm n	5.4E+06 1.2E+05 1.2E+04	nm nm n	9.4E-01 n	n	3.9E+00 n	n	2.0E+03 2.9E+02	n n		4.0E-01 7.0E-02	n n			
3.5E-01	I	1.0E-04	I	5.0E-04	I	7.0E-04	I	V				0.1	Chloranil Chlordane Chlordecone (Kepone)	118-75-2 12789-03-6 143-50-0	1.3E+00 1.7E+00 5.4E-02	c c* c	5.7E+00 7.5E+00 2.3E-01	c c* c	2.8E-02 c*	1.2E-01 c*	4.5E-02 c*	2.0E+00	1.8E-01 4.5E-02 3.5E-03	c c* c		1.5E-04 3.0E-03 1.2E-04	c c* c	1.4E-01		
1.0E+01	I	4.6E-03	C	3.0E-04	I			V				0.1	Chlorfenvinphos Chlorimuron, Ethyl- Chlorine	470-90-6 90982-32-4 7782-50-5	4.4E+01 1.3E+03 1.8E-01	n n n	5.7E+02 1.6E+04 7.8E-01	n n n	5.7E+02 n	n	6.4E-01 n	n	1.1E+01 3.9E+02 3.0E-01	n n n		3.1E-02 1.3E-01 1.4E-04	n n n			
3.0E-02	I	2.0E-04	I	3.0E-02	I			V				1.2E+03	Chlorine Dioxide Chlorite (Sodium Salt) Chloro-1,1-difluoroethane, 1-	10049-04-4 7758-19-2 75-68-3	2.3E+03 2.3E+03 5.4E+04	n n ns	3.4E+04 3.5E+04 2.3E+05	n n nms	2.1E-01 n	n	8.8E-01 n	n	4.2E-01 6.0E+02 1.0E+05	n n n		5.2E+01 n n	n n n			
4.6E-01	H			3.0E-04	I	2.0E-02	H	2.0E-02	I	V		0.1	Chloro-1,3-butadiene, 2- Chloro-2-methylaniline HCl, 4- Chloro-2-methylaniline, 4-	126-99-8 3165-93-3 95-69-2	1.0E-02 1.2E+00 5.4E+00	c c c*	4.4E-02 5.0E+00 2.3E+01	c c c	9.4E-03 c	c	4.1E-02 c	c	1.9E-02 1.7E-01 6.9E-01	c c c*		9.8E-06 1.5E-04 3.9E-04	c c c*			
2.7E-01	X							V				2.8E+04	Chloroacetaldehyde, 2-	107-20-0	2.6E+00	c	1.2E+01	c					2.9E-01	c		5.8E-05	c			

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Screening Levels								Protection of Ground Water SSLs			
SFO (mg/kg-day) ⁻¹	ke IUR (ug/m ³ -day) ⁻¹	ke RfD _o (mg/kg-day)	ke RfC _i (mg/m ³ -day)	ke V o mutagen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)
		2.0E-03	H					Chloroacetic Acid	79-11-8	1.3E+02	n	1.6E+03	n	3.1E-02	n	1.3E-01	n	4.0E+01	n	6.0E+01	8.1E-03	n	1.2E-02
				3.0E-05	I		0.1	Chloroacetophenone, 2-	532-27-4	4.3E+04	n	1.8E+05	nm	3.1E-02	n	1.3E-01	n						
2.0E-01	P	4.0E-03	I				0.1	Chloroaniline, p-	106-47-8	2.7E+00	c*	1.1E+01	c					3.6E-01	c		1.6E-04	c	
1.1E-01	C	3.1E-05	C	5.0E-02	P	V		Chlorobenzene	108-90-7	2.8E+02	n	1.3E+03	ns	5.2E+01	n	2.2E+02	n	7.8E+01	n	1.0E+02	5.3E-02	n	6.8E-02
		2.0E-02	I				0.1	Chlorobenzilate	510-15-6	4.9E+00	c	2.1E+01	c	9.1E-02	c	4.0E-01	c	3.1E-01	c		1.0E-03	c	
		3.0E-02	X				0.1	Chlorobenzoic Acid, p-	74-11-3	1.9E+03	n	2.5E+04	n					5.1E+02	n		1.3E-01	n	
		3.0E-03	P	3.0E-01	P	V		Chlorobenzotrifluoride, 4-	98-56-6	2.1E+02	ns	2.5E+03	ns	3.1E+02	n	1.3E+03	n	3.5E+01	n		1.2E-01	n	
		4.0E-02	P			V		Chlorobutane, 1-	109-69-3	3.1E+03	ns	4.7E+04	ns					6.4E+02	n		2.6E-01	n	
				5.0E+01	I	V		Chlorodifluoromethane	75-45-6	4.9E+04	ns	2.1E+05	nms	5.2E+04	n	2.2E+05	n	1.0E+05	n		4.3E+01	n	
3.1E-02	C	2.3E-05	I	2.0E-02	P	V		Chloroethanol, 2-	107-07-3	1.6E+03	n	2.3E+04	n					4.0E+02	n		8.1E-02	n	
		1.0E-02	I	9.8E-02	A	V		Chloroform	67-66-3	3.2E-01	c	1.4E+00	c	1.2E-01	c	5.3E-01	c	2.2E-01	c	8.0E+01(F)	6.1E-05	c	2.2E-02
				9.0E-02	I	V		Chloromethane	74-87-3	1.1E+02	n	4.6E+02	n	9.4E+01	n	3.9E+02	n	1.9E+02	n		4.9E-02	n	
2.4E+00	C	6.9E-04	C			V		Chloromethyl Methyl Ether	107-30-2	2.0E-02	c	8.9E-02	c	4.1E-03	c	1.8E-02	c	6.5E-03	c		1.4E-06	c	
3.0E-01	P	3.0E-03	P	1.0E-05	X		0.1	Chloronitrobenzene, o-	88-73-3	1.8E+00	c	7.7E+00	c	1.0E-02	c	4.4E-02	c	2.3E-01	c		2.2E-04	c	
6.3E-03	P	1.0E-03	P	6.0E-04	P			Chloronitrobenzene, p-	100-00-5	6.3E+01	n	3.6E+02	c**	6.3E-01	n	2.6E+00	n	1.1E+01	c**		1.0E-02	c**	
		5.0E-03	I			V		Chlorophenol, 2-	95-57-8	3.9E+02	n	5.8E+03	n					9.1E+01	n		7.4E-02	n	
				4.0E-04	C	V		Chloropicrin	76-06-2	2.0E+00	n	8.2E+00	n	4.2E-01	n	1.8E+00	n	8.3E-01	n		2.5E-04	n	
3.1E-03	C	8.9E-07	C	1.5E-02	I	V		Chlorothaloni	1897-45-6	1.8E+02	c**	7.4E+02	c*	3.2E+00	c	1.4E+01	c	2.2E+01	c*		4.9E-02	c*	
		2.0E-02	I			V		Chlorotoluene, o-	95-49-8	1.6E+03	ns	2.3E+04	ns					2.4E+02	n		2.3E-01	n	
		2.0E-02	X			V		Chlorotoluene, p-	106-43-4	1.6E+03	ns	2.3E+04	ns					2.5E+02	n		2.4E-01	n	
2.4E+02	C	6.9E-02	C				0.1	Chlorzotocin	54749-90-5	2.3E-03	c	9.6E-03	c	4.1E-05	c	1.8E-04	c	3.2E-04	c		7.1E-08	c	
		2.0E-01	I				0.1	Chlorpropham	101-21-3	1.3E+04	n	1.6E+05	nm					2.8E+03	n		2.6E+00	n	
		1.0E-03	A				0.1	Chlorpyrifos	2921-89-2	6.3E+01	n	8.2E+02	n					8.4E+00	n		1.2E-01	n	
		1.0E-02	H				0.1	Chlorpyrifos Methyl	5598-13-0	6.3E+02	n	8.2E+03	n					1.2E+02	n		5.4E-01	n	
		5.0E-02	I				0.1	Chlorsulfuron	64902-72-3	3.2E+03	n	4.1E+04	n					9.9E+02	n		8.3E-01	n	
		8.0E-04	H				0.1	Chlorzoxiprol	60238-56-4	5.1E+01	n	6.6E+02	n					2.8E+00	n		7.3E-02	n	
5.0E-01	J	8.4E-02	S	1.5E+00	I		0.013	Chromium(III) Insoluble Salts	16065-83-1	1.2E+05	nm	1.8E+06	nm					2.2E+04	n		4.0E+07	n	
		3.0E-03	I	1.0E-04	I	M	0.025	Chromium(VI)	13540-29-9	3.0E-01	c	6.3E+00	c	1.2E-05	c	1.5E-04	c	3.5E-02	c	1.0E+02	6.7E-04	c	1.8E+05
							0.013	Chromium, Total	7440-47-3														
9.0E-03	P	3.0E-04	P	6.0E-06	P			Cobalt	7440-48-4	2.3E+01	n	3.5E+02	n	3.1E-04	c*	1.4E-03	c*	6.0E+00	n		2.7E-01	n	
6.2E-04	I	4.0E-02	H			V	M	Coke Oven Emissions	8007-45-2					1.6E-03	c	2.0E-02	c						
								Copper	7440-50-8	3.1E+03	n	4.7E+04	n					8.0E+02	n	1.3E+03	2.8E+01	n	4.6E+01
		5.0E-02	I	6.0E-01	C		0.1	Cresol, m-	108-39-4	3.2E+03	n	4.1E+04	n	6.3E+02	n	2.6E+03	n	9.3E+02	n		7.4E-01	n	
		5.0E-02	I	6.0E-01	C		0.1	Cresol, o-	95-49-7	3.2E+03	n	4.1E+04	n	6.3E+02	n	2.6E+03	n	9.3E+02	n		7.5E-01	n	
		1.0E-01	A	6.0E-01	C		0.1	Cresol, p-	106-44-5	6.3E+03	n	8.2E+04	n	6.3E+02	n	2.6E+03	n	1.9E+03	n		1.5E+00	n	
		1.0E-01	A				0.1	Cresol, p-chloro-m-	59150-7	6.3E+03	n	8.2E+04	n					1.4E+03	n		1.7E+00	n	
1.9E+00	H	1.0E-01	A	6.0E-01	C		0.1	Cresols	1319-77-3	6.3E+03	n	8.2E+04	n	6.3E+02	n	2.6E+03	n	1.9E+03	n		1.5E+00	n	
		1.0E-03	P			V		Crotonaldehyde, trans-	123-73-9	3.7E-01	c	1.7E+00	c					4.0E-02	c		8.2E-06	c	
		1.0E-01	I	4.0E-01	I	V		Cumene	98-82-8	1.9E+03	ns	9.9E+03	ns	4.2E+02	n	1.8E+03	n	4.5E+02	n		7.4E-01	n	
2.2E-01	C	6.3E-05	C				0.1	Cupferron	135-20-6	2.5E+00	c	1.0E+01	c	4.5E-02	c	1.9E-01	c	3.5E-01	c		6.1E-04	c	
8.4E-01	H	2.0E-03	H				0.1	Cyanazine	21125-46-2	6.5E-01	c	2.7E+00	c					8.7E-02	c		4.1E-05	c	
								Cyanides															
		1.0E-03	I					~Calcium Cyanide	592-01-8	7.8E+01	n	1.2E+03	n					2.0E+01	n				
		5.0E-03	I					~Copper Cyanide	544-92-3	3.9E+02	n	5.8E+03	n					1.0E+02	n				
		6.0E-04	I	8.0E-04	S	V		~Cyanide (CN-)	57-12-5	2.7E+00	n	1.2E+01	n	8.3E-01	n	3.5E+00	n	1.5E+00	n	2.0E+02	1.5E-02	n	2.0E+00
		1.0E-03	I			V		~Cyanogen	460-19-5	7.8E+01	n	1.2E+03	n					2.0E+01	n				
		9.0E-02	I			V		~Cyanogen Bromide	506-68-3	7.0E+03	n	1.1E+05	nm					1.8E+03	n				
		5.0E-02	I			V		~Cyanogen Chloride	506-77-4	3.9E+03	n	5.8E+04	n					1.0E+03	n				
		6.0E-04	I	8.0E-04	I	V		~Hydrogen Cyanide	74-90-8	2.3E+01	n	1.5E+02	n	8.3E-01	n	3.5E+00	n	1.5E+00	n		1.5E-02	n	
		2.0E-03	I					~Potassium Cyanide	151-50-8	1.6E+02	n	2.3E+03	n					4.0E+01	n				
		5.0E-03	I				0.04	~Potassium Silver Cyanide	506-61-6	3.9E+02	n	5.8E+03	n					8.2E+01	n				
		1.0E-01	I				0.04	~Silver Cyanide	506-64-9	7.8E+03	n	1.2E+05	nm					1.8E+03	n				
		1.0E-03	I					~Sodium Cyanide	143-33-9	7.8E+01	n	1.2E+03	n					2.0E+01	n	2.0E+02			
		2.0E-04	P					~Thiocyanates	NA	1.6E+01	n	2.3E+02	n					4.0E+00	n				
		2.0E-04	X			V		~Thiocyanic Acid	463-56-9	1.6E+01	n	2.3E+02	n					4.0E+00	n				
		5.0E-02	I					~Zinc Cyanide	557-21-1	3.9E+03	n	5.8E+04	n					1.0E+03	n				
				6.0E+00	I	V		Cyclohexane	110-82-7	6.5E+03	ns	2.7E+04	ns	6.3E+03	n	2.6E+04	n	1.3E+04	n		1.3E+01	n	
2.3E-02	H						0.1	Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	2.4E+01	c	1.0E+02	c					2.4E+00	c		1.4E-02	c	
		5.0E+00	I	7.0E-01	P	V		Cyclohexanone	108														

Regional Screening Level (RSL) Summary Table (TR-1E-6, HQ=1) June 2015 (revised)

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Screening Levels								Protection of Ground Water SSLs							
SFO (mg/kg-day) ⁻¹	ke IUR (ug/m ³) ⁻¹	ke RfD _o	ke RfC ₁ (mg/m ³)	ke V o	muta-gen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)			
3.4E-01	I 9.7E-05	C			V				DDE, p,p'	72-55-9	2.0E+00	c	9.3E+00	c	2.9E-02	c	1.3E-01	c	4.6E-02	c			1.1E-02	c			
3.4E-01	I 9.7E-05	I	5.0E-04	I			1	0.03	DDT	50-29-3	1.9E+00	c*	8.5E+00	c*	2.9E-02	c	1.3E-01	c	2.3E-01	c*			7.7E-02	c*			
			1.0E-02	I			1	0.1	Dacthal	1861-32-1	6.3E+02	n	8.2E+03	n					1.2E+02	n			1.5E-01	n			
			3.0E-02	I			1	0.1	Dalapon	75-99-0	1.9E+03	n	2.5E+04	n					6.0E+02	n	2.0E+02		1.2E-01	n	4.1E-02		
7.0E-04	I		7.0E-03	I			1	0.1	Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5	4.4E+02	n	3.3E+03	c**					1.1E+02	c**			6.2E+01	c**			
			4.0E-05	I			1	0.1	Demeton	8065-48-3	2.5E+00	n	3.3E+01	n					6.7E-01	n				n			
1.2E-03	I		6.0E-01	I			1	0.1	Di(2-ethylhexyl)adipate	103-23-1	4.5E+02	c*	1.9E+03	c					6.5E+01	c	4.0E+02		4.7E+00	c	2.9E+01		
6.1E-02	H						1	0.1	Diallate	2303-16-4	8.9E+00	c	3.8E+01	c					5.2E-01	c			7.8E-04	c			
			7.0E-04	A			1	0.1	Diazinon	333-41-5	4.4E+01	n	5.7E+02	n					1.0E+01	n			6.5E-02	n			
			1.0E-02	X		V	1		Dibenzothiophene	132-65-0	7.8E+02	n	1.2E+04	n					6.5E+01	n			1.2E+00	n			
8.0E-01	P	6.0E-03	P	2.0E-04	P	2.0E-04	I	V	M	1	9.8E+02																
			4.0E-04	X		V	1		Dibromobenzene, 1,3-	96-12-8	5.3E-03	c	6.4E-02	c	1.7E-04	c	2.0E-03	c	3.3E-04	c*	2.0E-01		1.4E-07	c*	8.6E-05		
							1		Dibromobenzene, 1,3-	108-36-1	3.1E+01	n	4.7E+02	ns					5.3E+00	n			5.1E-03	n			
			1.0E-02	I		V	1		Dibromobenzene, 1,4-	106-37-6	7.8E+02	n	1.2E+04	n					1.3E+02	n			1.2E-01	n			
8.4E-02	I	2.7E-05	C	2.0E-02	I		V	1	8.0E+02		7.5E-01	c	3.3E+00	c	1.0E-01	c	4.5E-01	c	1.7E-01	c	8.0E+01(F)		4.5E-05	c	2.1E-02		
2.0E+00	I	6.0E-04	I	9.0E-03	I	9.0E-03	I	V	1	1.3E+03	3.6E-02	c	1.6E-01	c	4.7E-03	c	2.0E-02	c	7.5E-03	c	5.0E-02		2.1E-06	c	1.4E-05		
			1.0E-02	H	4.0E-03	X	V	1	2.8E+03		74-95-3	2.3E+01	n	9.8E+01	n	4.2E+00	n	1.8E+01	n	8.0E+00	n			2.0E-03	n		
			3.0E-04	P				1	0.1	Dibutyltin Compounds	NA	1.9E+01	n	2.5E+02	n					6.0E+00	n				n		
			3.0E-02	I			1	0.1	Dicamba	1918-00-9	1.9E+03	n	2.5E+04	n					5.7E+02	n			1.5E-01	n			
			4.2E-03	P		V	1	5.2E+02	Dichloro-2-butene, 1,4-	764-41-0	8.3E-03	c	3.6E-02	c	6.7E-04	c	2.9E-03	c	1.3E-03	c			6.2E-07	c			
			4.2E-03	P		V	1	5.2E+02	Dichloro-2-butene, cis-1,4-	1476-11-5	7.4E-03	c	3.2E-02	c	6.7E-04	c	2.9E-03	c	1.3E-03	c			6.2E-07	c			
			4.2E-03	P		V	1	7.6E+02	Dichloro-2-butene, trans-1,4-	110-57-6	7.4E-03	c	3.2E-02	c	6.7E-04	c	2.9E-03	c	1.3E-03	c			6.2E-07	c			
5.0E-02	I		4.0E-03	I			1	0.1	Dichloroacetic Acid	79-43-6	1.1E+01	c*	4.6E+01	c*					1.5E+00	c*	6.0E+01		3.1E-04	c*	1.2E-02		
			9.0E-02	I	2.0E-01	H	V	1	3.8E+02		1.8E+03	ns	9.3E+03	ns	2.1E+02	n	8.8E+02	n	3.0E+02	n	6.0E+02		3.0E-01	n	5.8E-01		
5.4E-03	C	1.1E-05	C	7.0E-02	A	8.0E-01	I	V	1	1.0E-06-7	2.6E+00	c	1.1E+01	c	2.6E-01	c	1.1E+00	c	4.8E-01	c	7.5E+01		4.6E-04	c	7.2E-02		
4.5E-01	I	3.4E-04	C					1	0.1	Dichlorobenzidine, 3,3'-	91-94-1	1.2E+00	c	5.1E+00	c	8.3E-03	c	3.6E-02	c	1.2E-01	c			8.1E-04	c		
			9.0E-03	X		V	1	0.1	Dichlorobenzophenone, 4,4'-	90-99-2	5.7E+02	n	7.4E+03	n					7.8E+01	n			4.7E-01	n			
			2.0E-01	I	1.0E-01	X	V	1	8.5E+02	Dichlorodifluoromethane	75-71-8	8.7E+01	n	3.7E+02	n	1.0E+02	n	4.4E+02	n	2.0E+02	n			3.0E-01	n		
5.7E-03	C	1.6E-06	C	2.0E-01	P		V	1	1.7E+03	Dichloroethane, 1,1,2-	75-34-3	3.6E+00	c	1.6E+01	c	1.8E+00	c	7.7E+00	c	2.7E+00	c			7.8E-04	c		
9.1E-02	I	2.6E-05	I	6.0E-03	X	7.0E-03	P	V	1	3.0E+03	4.6E-01	c*	2.0E+00	c*	1.1E-01	c*	4.7E-01	c*	1.7E-01	c*	5.0E+00		4.8E-05	c*	1.4E-03		
			5.0E-02	I	2.0E-01	I	V	1	1.2E+03	Dichloroethane, 1,2-	107-06-2	2.3E+02	n	1.0E+03	n	2.1E+02	n	8.8E+02	n	2.8E+02	n	7.0E+00		1.0E-01	n	2.5E-03	
			2.0E-03	I		V	1	2.4E+03	Dichloroethylene, 1,2-cis-	156-59-2	1.6E+02	n	2.3E+03	n					3.6E+01	n	7.0E+01		1.1E-02	n	2.1E-02		
			2.0E-02	I		V	1	1.9E+03	Dichloroethylene, 1,2-trans-	156-60-5	1.6E+03	n	2.3E+04	ns					3.6E+02	n	1.0E+02		1.1E-01	n	3.1E-02		
			3.0E-03	I		V	1	0.1	Dichlorophenol, 2,4-	120-83-2	1.9E+02	n	2.5E+03	n					4.6E+01	n			5.4E-02	n			
			1.0E-02	I		V	1	0.05	Dichlorophenoxy Acetic Acid, 2,4-	94-75-7	7.0E+02	n	9.6E+03	n					1.7E+02	n	7.0E+01		4.5E-02	n	1.8E-02		
			8.0E-03	I		V	1	0.1	Dichlorophenoxybutyric Acid, 4-(2,4-	94-82-6	5.1E+02	n	6.6E+03	n					1.2E+02	n			1.1E-01	n			
3.6E-02	C	1.0E-05	C	9.0E-02	A	4.0E-03	I	V	1	1.4E+03	Dichloropropane, 1,2-	79-87-5	1.0E+00	c*	4.4E+00	c*	2.8E-01	c*	1.2E+00	c*	4.4E-01	c*	5.0E+00		1.5E-04	c*	1.7E-03
			2.0E-02	P		V	1	1.5E+03	Dichloropropane, 1,3-	142-28-9	1.6E+03	ns	2.3E+04	ns					3.7E+02	n			1.3E-01	n			
			3.0E-03	I		V	1	0.1	Dichloropropanol, 2,3-	616-23-9	1.9E+02	n	2.5E+03	n					5.9E+01	n			1.3E-02	n			
			3.0E-02	I	2.0E-02	I	V	1	1.6E+03	Dichloropropene, 1,3-	542-75-6	1.8E+00	c*	8.2E+00	c*	7.0E-01	c*	3.1E+00	c*	4.7E-01	c*			1.7E-04	c*		
2.9E-01	I	8.3E-05	C	5.0E-04	I	5.0E-04	I		1	0.1	Dichlorvos	62-73-7	1.9E+00	c*	7.9E+00	c*	3.4E-02	c*	1.5E-01	c*	2.6E-01	c*			8.1E-05	c*	
			8.0E-02	P	3.0E-04	X	V	1	77-73-6	Dicycloptadiene	77-73-6	1.3E+00	n	5.4E+00	n	3.1E-01	n	1.3E+00	n	6.3E-01	n			2.2E-03	n		
1.6E+01	I	4.6E-03	I	5.0E-05	I			1	0.1	Dieldrin	60-51-1	3.4E-02	c*	1.4E-01	c	6.1E-04	c	2.7E-03	c	1.7E-03	c			6.9E-05	c		
			3.0E-04	C		5.0E-03	I		1	0.1	NA						9.4E-03	c	4.1E-02	c							
			2.0E-03	P	2.0E-04	P		1	0.1	Diethanolamine	111-42-2	1.3E+02	n	1.6E+03	n	2.1E-01	n	8.8E-01	n	4.0E+01	n			8.1E-03	n		
			3.0E-02	P	1.0E-04	P		1	0.1	Diethylene Glycol Monobutyl Ether	112-34-5	1.9E+03	n	2.4E+04	n	1.0E-01	n	4.4E-01	n	6.0E+02	n			1.3E-01	n		
			6.0E-02	P	3.0E-04	P		1	0.1	Diethylene Glycol Monoethyl Ether	111-90-0	3.8E+03	n	4.8E+04	n	3.1E-01	n	1.3E+00	n	1.2E+03	n			2.4E-01	n		
			1.0E-03	P		V	1	0.1	Diethylformamide	617-84-5	7.8E+01	n	1.2E+03	n					2.0E+01	n			4.1E-03	n			
3.5E+02	C	1.0E-01	C					1	1.1E+05	Diethylstilbestrol	56-53-1	1.6E-03	c	6.6E-03	c	2.8E-05	c	1.2E-04	c	4.9E-05	c			2.7E-05	c		
			8.0E-02	I			1	0.1	Difenzoquat	43222-48-6	5.1E+03	n	6.6E+04	n					1.6E+03	n				n			
			2.0E-02	I			1	0.1	Diflubenzuron	35367-38-5	1.3E+03	n	1.6E+04	n					2.9E+02	n			3.3E-01	n			
			4.0E+01	I	V		1	1.4E+03	Difluoroethane, 1,1-	75-37-6	4.8E+04	ns	2.0E+05	nms	4.2E+04	n	1.8E+05	n	8.3E+04	n							

Regional Screening Level (RSL) Summary Table (TR=1E-6, HQ=1) June 2015 (revised)

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Screening Levels								Protection of Ground Water SSLs					
SFO (mg/kg-day) ⁻¹	ke IUR (ug/m ³) ⁻¹	ke RfD _o (mg/kg-day)	ke RfC ₁ (mg/m ³)	ke V o c	muta-gen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)	
5.5E+02	C	1.6E-01	C					1.9E+05	Dimethylhydrazine, 1,2-	540-73-8	8.8E-04	c	4.1E-03	c	1.8E-05	c	7.7E-05	c	2.8E-05	c			6.5E-09	c	
		2.0E-02	I					0.1	Dimethylphenol, 2,4-	105-67-9	1.3E+03	n	1.6E+04	n					3.6E+02	n			4.2E-01	n	
		6.0E-04	I					0.1	Dimethylphenol, 2,6-	576-26-1	3.8E+01	n	4.9E+02	n					1.1E+01	n			1.3E-02	n	
		1.0E-03	I					0.1	Dimethylphenol, 3,4-	95-65-8	6.3E+01	n	8.2E+02	n					1.8E+01	n			2.1E-02	n	
4.5E-02	C	1.3E-05	C					1.1E+03	Dimethylvinylchloride	513-37-1	2.1E-01	c	9.4E-01	c	2.2E-01	c	9.4E-01	c	3.3E-01	c			2.0E-04	c	
		8.0E-05	X					0.1	Dinitro-o-cresol, 4,6-	534-52-1	5.1E+00	n	6.6E+01	n					1.5E+00	n			2.6E-03	n	
		2.0E-03	I					0.1	Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5	1.3E+02	n	1.6E+03	n					2.3E+01	n			7.7E-01	n	
		1.0E-04	P					0.1	Dinitrobenzene, 1,2-	528-29-0	6.3E+00	n	8.2E+01	n					1.9E+00	n			1.8E-03	n	
		1.0E-04	I					0.1	Dinitrobenzene, 1,3-	99-65-0	6.3E+00	n	8.2E+01	n					2.0E+00	n			1.8E-03	n	
		1.0E-04	P					0.1	Dinitrobenzene, 1,4-	100-25-4	6.3E+00	n	8.2E+01	n					2.0E+00	n			1.8E-03	n	
		2.0E-03	I					0.1	Dinitrophenol, 2,4-	51-28-5	1.3E+02	n	1.6E+03	n					3.9E+01	n			4.4E-02	n	
6.8E-01	I							0.1	Dinitrotoluene Mixture, 2,4/2,6-	NA	8.0E-01	c	3.4E+00	c					1.1E-01	c			1.5E-04	c	
3.1E-01	C	8.9E-05	C					0.102	Dinitrotoluene, 2,4-	121-14-2	1.7E+00	c*	7.4E+00	c	3.2E-02	c	1.4E-01	c	2.4E-01	c			3.2E-04	c	
1.5E+00	P							0.099	Dinitrotoluene, 2,6-	606-20-2	3.6E-01	c*	1.5E+00	c					4.8E-02	c			6.7E-05	c	
		2.0E-03	S					0.006	Dinitrotoluene, 2-Amino-4,6-	35572-78-2	1.5E+02	n	2.3E+03	n					3.9E+01	n			3.0E-02	n	
		2.0E-03	S					0.009	Dinitrotoluene, 4-Amino-2,6-	19406-51-0	1.5E+02	n	2.3E+03	n					3.9E+01	n			3.0E-02	n	
4.5E-01	X							0.1	Dinitrotoluene, Technical grade	25321-14-6	1.2E+00	c*	5.1E+00	c					1.6E-01	c			2.2E-04	c	
		9.0E-04	X					0.1	Dioxseb	88-85-7	6.3E+01	n	8.2E+02	n					1.5E+01	n		7.0E+00	1.3E-01	n	6.2E-02
1.0E-01	I	5.0E-06	I					1.2E+05	Dioxane, 1,4-	123-91-1	5.3E+00	c	2.4E+01	c	5.6E-01	c*	2.5E+00	c*	4.6E-01	c			9.4E-05	c	
6.2E+03	I	1.3E+00	I					0.03	Dioxins	NA	1.0E-04	c	4.7E-04	c	2.2E-06	c	9.4E-06	c	1.3E-05	c			1.7E-05	c	
1.3E+05	C	3.8E+01	C					0.03	- Hexachlorodibenzo-p-dioxin, Mixture ~ I CDD, 2,3,7,8-	1746-01-6	4.8E-06	c*	2.2E-05	c*	7.4E-08	c	3.2E-07	c	1.2E-07	c	3.0E-05		5.9E-08	c	1.5E-05
		3.0E-02	I					0.1	Diphenamid	957-51-7	1.9E+03	n	2.5E+04	n					5.3E+02	n			5.2E+00	n	
		8.0E-04	X					0.1	Diphenyl Sulfone	127-83-9	5.1E+01	n	6.6E+02	n					1.5E+01	n			3.6E-02	n	
		2.5E-02	I					0.1	Diphenylamine	122-39-4	1.6E+03	n	2.1E+04	n					3.1E+02	n			5.8E-01	n	
8.0E-01	I	2.2E-04	I					0.1	Diphenylhydrazine, 1,2-	122-86-7	6.8E-01	c	2.9E+00	c	1.3E-02	c	5.6E-02	c	7.7E-02	c			2.5E-04	c	
		2.2E-03	I					0.1	Diquat	89-00-7	1.4E+02	n	1.8E+03	n					4.4E+01	n		2.0E+01	8.3E-01	n	3.7E-01
7.1E+00	C	1.4E-01	C					0.1	Direct Black 38	1937-37-7	7.6E-02	c	3.2E-01	c	2.0E-05	c	8.8E-05	c	1.1E-02	c			5.3E+00	c	
7.4E+00	C	1.4E-01	C					0.1	Direct Blue 61	2802-43-2	7.3E-02	c	3.1E-01	c	2.0E-05	c	8.8E-05	c	1.1E-02	c			1.7E+01	c	
6.7E+00	C	1.4E-01	C					0.1	Direct Brown 95	16071-86-6	8.1E-02	c	3.4E-01	c	2.0E-05	c	8.8E-05	c	1.2E-02	c				c	
		4.0E-05	I					0.1	Disulfoton	298-04-4	2.5E+00	n	3.3E+01	n					5.0E-01	n			9.4E-04	n	
		1.0E-02	I					0.1	Dithiane, 1,4-	505-29-3	7.8E+02	n	1.2E+04	n					2.0E+02	n			9.7E-02	n	
		2.0E-03	I					0.1	Diuron	330-54-1	1.3E+02	n	1.6E+03	n					3.6E+01	n			1.5E-02	n	
		4.0E-03	I					0.1	Dodine	2439-10-3	2.5E+02	n	3.3E+03	n					8.0E+01	n			4.1E-01	n	
		2.5E-02	I					0.1	EPTO	759-94-4	2.0E+03	n	2.9E+04	n					3.8E+02	n			2.0E-01	n	
		6.0E-03	I					0.1	Endosulfan	116-29-7	4.7E+02	n	7.0E+03	n					1.0E+02	n			1.4E+00	n	
		2.0E-02	I					0.1	Endothal	146-73-3	1.3E+03	n	1.6E+04	n					3.8E+02	n		1.0E+02	9.1E-02	n	2.4E-02
		3.0E-04	I					0.1	Erdrin	72-20-8	1.9E+01	n	2.5E+02	n					2.3E+00	n		2.0E+00	9.2E-02	n	8.1E-02
9.9E-03	I	1.2E-06	I					1.1E+04	Epichlorohydrin	106-89-8	1.9E+01	n	8.2E+01	n	1.0E+00	n	4.4E+00	n	2.0E+00	n			4.5E-04	n	
		6.0E-03	P	1.0E-03	I	V		1.1E+04	Epoxybutane, 1,2-	106-88-7	1.6E+02	n	6.7E+02	n	2.1E+01	n	8.8E+01	n	4.2E+01	n			9.2E-03	n	
		5.0E-03	I					0.1	Ethephon	16672-87-0	3.2E+02	n	4.1E+03	n					1.0E+02	n			2.1E-02	n	
		5.0E-04	I					0.1	Ethion	563-12-2	3.2E+01	n	4.1E+02	n					4.3E+00	n			8.5E-03	n	
		1.0E-01	P	6.0E-02	P	V		3.1E+04	Ethoxyethanol Acetate, 2-	111-15-9	2.6E+03	n	1.4E+04	n	6.3E+01	n	2.6E+02	n	1.2E+02	n			2.5E-02	n	
		9.0E-02	P	2.0E-01	I	V		1.1E+05	Ethoxyethanol, 2-	110-80-5	5.2E+03	n	4.7E+04	n	2.1E+02	n	8.8E+02	n	3.4E+02	n			6.8E-02	n	
		9.0E-01	I	7.0E-02	P	V		1.1E+04	Ethyl Acetate	141-78-6	6.2E+02	n	2.6E+03	n	7.3E+01	n	3.1E+02	n	1.4E+02	n			3.1E-02	n	
4.8E-02	H							2.5E+03	Ethyl Acrylate	140-88-5	1.4E+01	c**	6.8E+01	c**	8.3E+00	n	3.5E+01	n	1.6E+00	c**			3.5E-04	c**	
		1.0E+01	I	V				2.1E+03	Ethyl Chloride (Chloroethane)	75-00-3	1.4E+04	ns	5.7E+04	ns	1.0E+04	n	4.4E+04	n	2.1E+04	n			5.9E+00	n	
		2.0E-01	I	V				1.0E+04	Ethyl Ether	60-29-7	1.6E+04	ns	2.3E+05	nms					3.9E+03	n			8.8E-01	n	
		9.0E-02	H	3.0E-01	P	V		1.1E+03	Ethyl Methacrylate	97-63-2	1.4E+03	ns	7.1E+03	ns	3.1E+02	n	1.3E+03	n	4.6E+02	n			1.1E-01	n	
		1.0E-05	I					0.1	Ethyl-p-nitrophenyl Phosphonate	2104-64-5	6.3E-01	n	8.2E+00	n					8.9E-02	n			2.8E-03	n	
1.1E-02	C	2.5E-06	C					4.8E+02	Ethylbenzene	100-41-4	5.8E+00	c	2.5E+01	c	1.1E+00	c	4.9E+00	c	1.5E+00	c			1.7E-03	c	7.8E-01
		1.0E-01	I	1.0E+00	I	V		1.1E+03	Ethylene Cyanohydrin	109-78-4	4.4E+03	n	5.7E+04	n					1.4E+03	n			2.8E-01	n	
		9.0E-02	P					1.9E+05	Ethylene Diamine	107-15-3	7.0E+03	n	1.1E+05	nm					1.8E+03	n			4.1E-01	n	
		2.0E+00	I	4.0E-01	C			1.0E+01	Ethylene Glycol	107-21-1	1.3E+05	nm	1.6E+06	nm	4.2E+02	n	1.8E+03	n	4.0E+04	n			8.1E+00	n	
		1.0E-01	I	1.6E+00	I			1.0E+01	Ethylene Glycol Monobutyl Ether	111-76-2	6.3E+03	n	8.2E+04	n	1.7E+03	n	7.0E+03	n	2.0E+03	n			4.1E-01	n	
3.1E																									

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information											Contaminant		Screening Levels								Protection of Ground Water SSLs							
SFO (mg/kg-day) ⁻¹	ke (y)	IUR (ug/m ³) ⁻¹	ke (y)	RfD _o (mg/kg-day)	ke (y)	RfC ₁ (mg/m ³) ⁻¹	ke (y)	muta-gen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)	
				8.0E-02	I					1	0.1	Fluridone	59756-60-4	5.1E+03	n	6.6E+04	n					1.4E+03	n		1.6E+02	n		
				2.0E-02	I					1	0.1	Flurprimidol	56425-91-3	1.3E+03	n	1.6E+04	n					3.4E+02	n		1.6E+00	n		
				6.0E-02	I					1	0.1	Flutolanil	66332-96-5	3.8E+03	n	4.9E+04	n					9.5E+02	n		5.0E+00	n		
				1.0E-02	I					1	0.1	Fluvalinate	69409-94-5	6.3E+02	n	8.2E+03	n					2.0E+02	n		2.9E+02	n		
3.5E-03	I			1.0E-01	I					1	0.1	Folpet	133-07-3	1.6E+02	c*	6.6E+02	c					2.0E+01	c*		4.7E-03	c*		
1.9E-01	I									1	0.1	Fomesafen	72178-02-0	2.9E+00	c	1.2E+01	c					3.9E-01	c		1.3E-03	c		
				2.0E-03	I					1	0.1	Fonofos	944-22-9	1.3E+02	n	1.6E+03	n					2.4E+01	n		4.7E-02	n		
1.3E-05	I			2.0E-01	I	9.8E-03	A	V		1		Formaldehyde	50-00-0	1.7E+01	c*	7.3E+01	c*	2.2E-01	c*	9.4E-01	c*	4.3E-01	c*		8.7E-05	c*		
				9.0E-01	P	3.0E-04	X	V		1	1.1E+05	Formic Acid	64-18-6	2.9E+01	n	1.2E+02	n	3.1E-01	n	1.3E+00	n	6.3E-01	n		1.3E-04	n		
				3.0E+00	I					1	0.1	Fosetyl-AL	39148-24-8	1.9E+05	nm	2.5E+06	nm					6.0E+04	n			n		
												Furans																
				1.0E-03	X			V		1	0.03	-Dibenzofuran	132-64-9	7.3E+01	n	1.0E+03	n					7.9E+00	n		1.5E-01	n		
				1.0E-03	I			V		1	0.03	-Furan	110-00-9	7.3E+01	n	1.0E+03	n					1.9E+01	n		7.3E-03	n		
				9.0E-01	I	2.0E+00	I	V		1	0.03	-Tetrahydrofuran	109-99-9	1.8E+04	n	9.6E+04	n	2.1E+03	n	8.8E+03	n	3.4E+03	n		7.5E-01	n		
3.8E+00	H									1	0.1	Furazolidone	67-45-8	1.4E-01	c	6.0E-01	c					2.0E-02	c		3.9E-05	c		
				3.0E-03	I	5.0E-02	H	V		1	1.0E+04	Furfural	98-01-1	2.1E+02	n	2.6E+03	n	5.2E+01	n	2.2E+02	n	3.8E+01	n		8.1E-03	n		
1.5E+00	C	4.3E-04	C							1	0.1	Furium	531-82-8	3.6E-01	c	1.5E+00	c	6.5E-03	c	2.9E-02	c	5.0E-02	c		8.8E-05	c		
3.0E-02	I	8.6E-06	C							1	0.1	Furmecyclo	60568-05-0	1.8E+01	c	7.7E+01	c	3.3E-01	c	1.4E+00	c	1.1E+00	c		1.2E-03	c		
				4.0E-04	I					1	0.1	Glufosinate, Ammonium	77182-82-2	2.5E+01	n	3.3E+02	n					8.0E+00	n		1.8E-03	n		
						8.0E-05	C			1	0.1	Glutaraldehyde	111-30-8	1.1E+05	nm	4.8E+05	nm	8.3E-02	n	3.5E-01	n					n		
				4.0E-04	I	1.0E-03	H	V		1	1.1E+05	Glycidyl	785-34-4	2.2E+01	n	1.9E+02	n	1.0E+00	n	4.4E+00	n	1.7E+00	n		3.3E-04	n		
				1.0E-01	I					1	0.1	Glycosate	1071-83-6	6.3E+03	n	8.2E+04	n					2.0E+03	n	7.0E+02	8.8E+00	n	3.1E+00	
				3.0E-03	I					1	0.1	Goal	428/14-U3-3	1.9E+02	n	2.5E+03	n					3.2E+01	n		2.5E+00	n		
				1.0E-02	X			V		1		Guandine	113-00-8	7.8E+02	n	1.2E+04	n					2.0E+02	n		4.5E-02	n		
				2.0E-02	P					1	0.1	Guandine Chloride	50-01-1	1.3E+03	n	1.6E+04	n					4.0E+02	n			n		
				3.0E-03	A	1.0E-02	A			1	0.1	Guthion	88-50-0	1.9E+02	n	2.5E+03	n	1.0E+01	n	4.4E+01	n	5.6E+01	n		1.7E-02	n		
				5.0E-05	I					1	0.1	Haloxypol, Methyl	66806-40-2	3.2E+00	n	4.1E+01	n					7.6E-01	n		8.4E-03	n		
				1.3E-02	I					1	0.1	Harmony	78277-27-3	8.2E+02	n	1.1E+04	n					2.6E+02	n		7.8E-02	n		
4.5E+00	I	1.3E-03	I	5.0E-04	I			V		1		Heptachlor	76-41-5	1.3E-01	c	6.3E-01	c	2.2E-03	c	9.4E-03	c	1.4E-03	c	4.0E-01	1.1E-04	c	3.3E-02	
9.1E+00	I	2.6E-03	I	1.3E-05	I			V		1		Heptachlor Epoxide	1024-57-3	7.0E-02	c*	3.3E-01	c*	1.1E-03	c	4.7E-03	c	1.4E-03	c*	2.0E-01	2.8E-05	c*	4.1E-03	
				2.0E-03	I			V		1		Hexabromobenzene	87-82-1	1.6E+02	n	2.3E+03	n					4.0E+01	n		2.3E-01	n		
				2.0E-04	I			V		1	0.1	Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BDE-153)	68631-49-2	1.3E+01	n	1.6E+02	n					4.0E+00	n			n		
1.6E+00	I	4.6E-04	I	8.0E-04	I			V		1		Hexachlorobenzene	148-74-1	2.1E-01	c	9.6E-01	c	6.1E-03	c	2.7E-02	c	9.8E-03	c	1.0E+00	1.2E-04	c	1.3E-02	
7.8E-02	I	2.2E-05	I	1.0E-03	P			V		1	1.7E+01	Hexachlorobutadiene	87-68-3	1.2E+00	c*	5.3E+00	c	1.3E-01	c	5.6E-01	c	1.4E-01	c*		2.6E-04	c*		
6.3E+00	I	1.8E-03	I	8.0E-03	A					1	0.1	Hexachlorocyclohexane, Alpha	319-84-6	8.6E-02	c	3.6E-01	c	1.6E-03	c	6.8E-03	c	7.1E-03	c		4.1E-05	c		
1.8E+00	I	5.3E-04	I							1	0.1	Hexachlorocyclohexane, Beta	319-85-7	3.0E-01	c	1.3E+00	c	5.3E-03	c	2.3E-02	c	2.5E-02	c		1.4E-04	c		
1.1E+00	C	3.1E-04	C	3.0E-04	I					1	0.4	Hexachlorocyclohexane, Gamma (Lindane)	58-89-9	5.7E-01	c*	2.5E+00	c	9.1E-03	c	4.0E-02	c	4.1E-02	c*	2.0E-01	2.4E-04	c*	1.2E-03	
1.8E+00	I	5.1E-04	I							1	0.1	Hexachlorocyclohexane, Technical	608-73-1	3.0E-01	c	1.3E+00	c	5.5E-03	c	2.4E-02	c	2.5E-02	c		1.4E-04	c		
4.0E-02	I	1.1E-05	C	6.0E-03	I	2.0E-04	I	V		1	1.6E+01	Hexachlorocyclopentadiene	77-47-4	1.8E+00	n	7.5E+00	n	2.1E-01	n	8.8E-01	n	4.1E-01	n	5.0E+01	1.3E-03	n	1.6E-01	
				7.0E-04	I	3.0E-02	I	V		1		Hexachloroethane	67-72-1	1.8E+00	c*	8.0E+00	c*	2.6E-01	c	1.1E+00	c	3.3E-01	c*		2.0E-04	c*		
				3.0E-04	I					1	0.1	Hexachlorophene	70-30-4	1.9E+01	n	2.5E+02	n					6.0E+00	n		8.0E+00	n		
1.1E-01	I			3.0E-03	I					1	0.015	Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	6.1E+00	c*	2.8E+01	c					7.0E-01	c*		2.7E-04	c*		
				4.0E-04	P	1.0E-05	I	V		1	5.2E+03	Hexamethylene Diisocyanate, 1,6-Hexamethylphosphoramide	822-06-0 680-31-9	3.1E+00 2.5E+01	n n	1.3E+01 3.3E+02	n n	1.0E-02	n	4.4E-02	n	2.1E-02	n	8.0E+00	n		2.1E-04	n
				6.0E-02	H	7.0E-01	I	V		1	1.4E+02	Hexane, N-	110-54-3	5.4E+02	ns	2.5E+03	ns	7.3E+02	n	3.1E+03	n	3.2E+02	n		2.3E+00	n		
				2.0E+00	P					1	0.1	Hexanedioic Acid	124-04-9	1.3E+05	nm	1.6E+06	nm					4.0E+04	n		9.9E+00	n		
				5.0E-03	I	3.0E-02	I	V		1	3.3E+03	Hexanone, 2-	591-78-6	2.0E+02	n	1.3E+03	n	3.1E+01	n	1.3E+02	n	3.8E+01	n		8.8E-03	n		
				3.3E-02	I					1	0.1	Hexazinone	51235-04-2	2.1E+03	n	2.7E+04	n					6.4E+02	n		3.0E-01	n		
3.0E+00	I	4.9E-03	I			3.0E-05	P	V		1		Hydrazine	302-01-2	2.3E-01	c	1.1E+00	c	5.7E-04	c*	2.5E-03	c*	1.1E-03	c*			c*		
3.0E+00	I	4.9E-03	I							1		Hydrazine Sulfate	10034-93-2	2.3E-01	c	1.1E+00	c	5.7E-04	c	2.5E-03	c	2.6E-02	c			c		
				2.0E-02	I			V		1		Hydrogen Chloride	7647-01-0	2.8E+07	nm	1.2E+08	nm	2.1E+01	n	8.8E+01	n	4.2E+01	n			n		
				4.0E-02	C	1.4E-02	C	V		1		Hydrogen Fluoride	7664-39-3	3.1E+03	n	4.7E+04	n											

Regional Screening Level (RSL) Summary Table (TR=1E-6, HQ=1) June 2015 (revised)

Toxicity and Chemical-specific Information											Contaminant		Screening Levels								Protection of Ground Water SSLs				
SFO (mg/kg-day) ⁻¹	Key IUR (ug/m ³ -y) ⁻¹	Key RfD _o (mg/kg-day)	Key RfC ₁ (mg/m ³ -y)	Key V _o	Key mutagen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)	
			3.0E-01	A	V		1		JP-7	NA	4.3E+08	nm	1.8E+09	nm	3.1E+02	n	1.3E+03	n	6.3E+02	n			n		
		7.5E-02	I				1	0.1	Kerb	23950-58-5	4.7E+03	n	6.2E+04	n					1.2E+03	n		1.2E+00	n		
		2.0E-03	I				1	0.1	Lactofen	77501-63-4	1.3E+02	n	1.6E+03	n					2.5E+01	n		1.2E+00	n		
									Lead Compounds																
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M	0.025	1	0.1														
8.5E-03	C	1.2E-05	C						1																
2.8E-01	C	8.0E-05	C						1	0.1															
8.5E-03	C	1.2E-05	C						1	0.1															
		1.0E-07	I		V		1	2.4E+00																	
		2.0E-03	I				1	0.1	Linuron	330-55-2	1.3E+02	n	1.6E+03	n					3.3E+01	n		2.9E-02	n		
		2.0E-03	P				1		Lithium	7439-93-2	1.6E+02	n	2.3E+03	n					4.0E+01	n		1.2E+01	n		
		2.0E-01	I				1	0.1	Londax	83055-99-8	1.3E+04	n	1.6E+05	nm					3.9E+03	n		1.0E+00	n		
		5.0E-04	I				1	0.1	MCPA	94-74-6	3.2E+01	n	4.1E+02	n					7.5E+00	n		2.0E-03	n		
		1.0E-02	I				1	0.1	MCPB	94-81-5	6.3E+02	n	8.2E+03	n					1.5E+02	n		5.8E-02	n		
		1.0E-03	I				1	0.1	MCPP	93-65-2	6.3E+01	n	8.2E+02	n					1.6E+01	n		4.6E-03	n		
		2.0E-02	I				1	0.1	Malathion	121-75-5	1.3E+03	n	1.6E+04	n					3.9E+02	n		1.0E-01	n		
		1.0E-01	I	7.0E-04	C		1	0.1	Maleic Anhydride	108-31-6	6.3E+03	n	8.0E+04	n	7.3E-01	n	3.1E+00	n	1.9E+03	n		3.8E-01	n		
		5.0E-01	I				1	0.1	Maleic Hydrazide	123-33-1	3.2E+04	n	4.1E+05	nm					1.0E+04	n		2.1E+00	n		
		1.0E-04	P				1	0.1	Malononitrile	109-77-3	6.3E+00	n	8.2E+01	n					2.0E+00	n		4.1E-04	n		
		3.0E-02	H				1	0.1	Mancozeb	8018-01-7	1.9E+03	n	2.5E+04	n					5.4E+02	n		1.0E-01	n		
		5.0E-03	I				1	0.1	Maneb	12427-38-2	3.2E+02	n	4.1E+03	n					9.8E+01	n		1.4E-01	n		
		1.4E-01	I	5.0E-05	I		1		Manganese (Diet)	7439-96-5															
		2.4E-02	S	5.0E-05	I		0.04		Manganese (Non-diet)	7439-96-5	1.8E+03	n	2.6E+04	n	5.2E-02	n	2.2E-01	n	4.3E+02	n		2.8E+01	n		
		9.0E-05	H				1	0.1	Mephosolan	950-10-7	5.7E+00	n	7.4E+01	n					1.8E+00	n		2.6E-03	n		
		3.0E-02	I				1	0.1	Meprosolone	24307-28-4	1.9E+03	n	2.5E+04	n					6.0E+02	n		2.0E-01	n		
									Mercury Compounds																
		3.0E-04	I	3.0E-04	S		0.07		Mercuric Chloride (and other Mercury salts)	7487-94-7	2.3E+01	n	3.5E+02	n	3.1E-01	n	1.3E+00	n	5.7E+00	n	2.0E+00		n		
				3.0E-04	I	V		3.1E+00	Mercury (elemental)	7439-97-6	9.4E+00	ns	4.0E+01	ns	3.1E-01	n	1.3E+00	n	6.3E-01	n	2.0E+00		n		
		1.0E-04	I				1		Methyl Mercury	22967-92-6	7.8E+00	n	1.2E+02	n					2.0E+00	n		3.3E-02	n	1.0E-01	
		8.0E-05	I				1	0.1	Phenylmercuric Acetate	62-38-4	5.1E+00	n	6.6E+01	n					1.6E+00	n		5.0E-04	n		
		3.0E-05	I		V		1		Merphos	150-50-5	2.3E+00	n	3.5E+01	n					6.0E-01	n		5.9E-02	n		
		3.0E-05	I				1	0.1	Merphos Oxide	78-48-8	1.9E+00	n	2.5E+01	n					8.5E-02	n		4.2E-04	n		
		6.0E-02	I				1	0.1	Metalaxyl	57837-19-1	3.8E+02	n	4.9E+04	n					1.2E+03	n		3.3E-01	n		
		1.0E-04	I	3.0E-02	P	V		4.6E+03	Methacrylonitrile	126-98-7	7.5E+00	n	1.0E+02	n	3.1E+01	n	1.3E+02	n	1.9E+00	n		4.3E-04	n		
		5.0E-05	I				1	0.1	Methamidophos	10265-92-6	3.2E+00	n	4.1E+01	n					1.0E+00	n		2.1E-04	n		
		2.0E+00	I	2.0E+01	I	V		1.1E+05	Methanol	67-56-1	1.2E+06	nms	1.2E+06	nms	2.1E+04	n	8.8E+04	n	2.0E+04	n		4.1E+00	n		
		1.0E-03	I				1	0.1	Methidathion	950-37-8	6.3E+01	n	8.2E+02	n					1.9E+01	n		4.7E-03	n		
		2.5E-02	I				1	0.1	Methomyl	16752-11-5	1.6E+03	n	2.1E+04	n					5.0E+02	n		1.1E-01	n		
4.9E-02	C	1.4E-05	C				1	0.1	Methoxy-5-nitroaniline, 2-	99-59-2	1.1E+01	c	4.7E+01	c	2.0E-01	c	8.8E-01	c	1.5E+00	c	4.0E+01	5.3E-04	c		
		5.0E-03	I				1	0.1	Methoxychlor	72-43-5	3.2E+02	n	4.1E+03	n					3.7E+01	n		2.0E+00	n	2.2E+00	
		8.0E-03	P	1.0E-03	P	V		1.2E+05	Methoxyethanol Acetate, 2-	110-49-6	1.1E+02	n	5.1E+02	n	1.0E+00	n	4.4E+00	n	2.1E+00	n		4.2E-04	n		
		5.0E-03	P	2.0E-02	I	V		1.1E+05	Methoxyethanol, 2-	109-86-4	3.3E+02	n	3.5E+03	n	2.1E+01	n	8.8E+01	n	2.9E+01	n		5.9E-03	n		
		1.0E+00	X		V		1	2.9E+04	Methyl Acetate	79-20-9	7.8E+00	ns	1.2E+06	nms					2.0E+04	n		4.1E+00	n		
		3.0E-02	H	2.0E-02	P	V		6.8E+03	Methyl Acrylate	96-33-3	1.4E+02	n	6.0E+02	n	2.1E+01	n	8.8E+01	n	3.9E+01	n		8.3E-03	n		
		6.0E-01	I	5.0E+00	I	V		2.8E+04	Methyl Ethyl Ketone (2-Butanone)	78-93-3	2.7E+04	n	1.9E+05	nms	5.2E+03	n	2.2E+04	n	5.6E+03	n		1.2E+00	n		
		1.0E-03	X				1	1.8E+05	Methyl Hydrazine	60-34-4	4.4E-01	c**	1.9E+00	c**	2.8E-03	c**	1.2E-02	c**	5.6E-03	c**		1.3E-06	c**		
		8.0E-02	H	3.0E+00	I	V		3.4E+03	Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1	5.3E+03	ns	5.6E+04	ns	3.1E+03	n	1.3E+04	n	1.2E+03	n		2.8E-01	n		
				1.0E-03	C	V		1.7E+04	Methyl Isocyanate	624-83-9	4.6E+00	n	1.9E+01	n	1.0E+00	n	4.4E+00	n	2.1E+00	n		5.9E-04	n		
		1.4E+00	I	7.0E-01	I	V		2.4E+03	Methyl Methacrylate	80-62-6	4.4E+03	ns	1.9E+04	ns	7.3E+02	n	3.1E+03	n	1.4E+03	n		3.0E-01	n		
		2.5E-04	I				1	0.1	Methyl Parathion	298-00-0	1.6E+01	n	2.1E+02	n					4.5E+00	n		7.4E-03	n		
		6.0E-02	X				1	0.1	Methyl Phosphonic Acid	993-13-5	3.8E+03	n	4.9E+04	n					1.2E+03	n		2.4E-01	n		
9.9E-02	C	2.8E-05	C				1	0.1	Methyl Styrene (Mixed Isomers)	25013-15-4	2.4E+02	n	1.6E+03	ns	4.2E+01	n	1.8E+02	n	3.8E+01	n		6.2E-02	n		
							1	0.1	Methyl methanesulfonate	66-27-3	5.5E+00	c	2.3E+01	c	1.0E-01	c	4.4E-01	c	7.9E-01	c		1.6E-04	c		
1.8E-03	C	2.6E-07	C			3.0E+00	I	V	1	8.9E+03	Methyl tert-Butyl Ether (MTBE)	1634-04-4	4.7E+01	c	2.1E+02	c	1.1E+01	c	4.7E+01	c	1.4E+01	c	3.2E-03	c	
		3.0E-04	X				1	0.1	Methyl-1,4-benzenediamine dithyodichloride, 2-	615-45-2	1.9E+01	n	2.5E+02	n					6.0E+00	n		3.6E-03	n		
9.0E-03	P	2.0E-02	X				1	0.1	Methyl-5-Nitroaniline, 2-	99-55-8	6.0E+01	c*	2.6E+02	c*					8.1E+00	c*		4.5E-03	c*		
8.3E+00	C	2.4E-03	C				1	0.1	Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	6.5E-02	c	2.8E-01	c	1.2E-03	c	5.1E-03	c	9.4E-03	c		3.2E-06	c		
1.3E-01	C	3.7E-05	C				1	0.1	Methylamliane Hydrochloride, 2-	636-21-5	4.2E+00	c	1.8E+01	c	7.6E-02	c	3.3E-01	c	6.0E-01	c		2.6E-04	c		
				1.0E-02	A		1	0.1	Methylarsonic acid	124-58-3	6.3E+02	n	8.2E+03												

Regional Screening Level (RSL) Summary Table (TR=1E-6, HQ=1) June 2015 (revised)

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Screening Levels								Protection of Ground Water SSLs						
SFO (mg/kg-day) ⁻¹	ke IUR (ug/m ³) ⁻¹	ke RfD _o (mg/kg-day)	ke RfC _i (mg/m ³)	ke RfC _v (mg/m ³)	mutagen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)		
4.6E-02	I	1.3E-05	C						Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-81-1	1.2E+01	c	5.0E+01	c	2.2E-01	c	9.4E-01	c	4.6E-01	c			2.6E-03	c		
1.6E+00	C	4.6E-04	C	2.0E-02	C	1	0.1		Methylenebisbenzenamine, 4,4'-	101-77-9	3.4E-01	c	1.4E+00	c	6.1E-03	c	2.7E-02	c	4.7E-02	c			2.1E-04	c		
				6.0E-04	I	1	0.1		Methylenediphenyl Diisocyanate	101-68-8	8.5E+05	nm	3.6E+06	nm	6.3E-01	n	2.6E+00	n								
		7.0E-02	H		V	1		5.0E+02	Methylstyrene, Alpha-	98-83-9	5.5E+03	ns	8.2E+04	ns					7.8E+02	n			1.2E+00	n		
1.5E-01	I								Metolachlor	51218-45-2	9.5E+03	n	1.2E+05	nm									2.7E+03	n	3.2E+00	
2.5E-02	I								Metribuzin	21087-64-9	1.6E+03	n	2.1E+04	n									4.9E+02	n	1.5E-01	
3.0E+00	P				V	1		3.4E-01	Mineral oils	8012-95-1	2.3E+05	nms	3.5E+06	nms									6.0E+04	n	2.4E+03	
1.8E+01	C	5.1E-03	C	2.0E-04	I				Mirex	2385-85-5	3.6E-02	c	1.7E-01	c	5.5E-04	c	2.4E-03	c	8.8E-04	c			6.3E-04	c		
		2.0E-03	I					0.1	Molinate	2212-67-1	1.3E+02	n	1.6E+03	n									3.0E+01	n	1.7E-02	
		5.0E-03	I						Molybdenum	7439-98-7	3.9E+02	n	5.8E+03	n									1.0E+02	n	2.0E+00	
1.0E-01	I								Monochloramine	10599-90-3	7.8E+03	n	1.2E+05	nm								4.0E+03		2.0E+03	n	
2.0E-03	P				V	1			Monomethylaniline	100-61-8	1.3E+02	n	1.6E+03	n									3.8E+01	n	1.4E-02	
3.0E-04	X								N,N-Diphenyl-1,4-benzenediamine	74-31-7	1.9E+01	n	2.5E+02	n									3.6E+00	n	3.7E-01	
2.0E-03	X				V	1			Naled	300-76-5	1.6E+02	n	2.3E+03	n									4.0E+01	n	1.8E-02	
3.0E-02	X	1.0E-01	P	V		1			Naphtha, High Flash Aromatic (HFAN)	64742-95-6	2.3E+03	n	3.5E+04	n	1.0E+02	n	4.4E+02	n	1.5E+02	n			1.0E+01	n		
1.8E+00	C	0.0E+00	C						Naphthylamine, 2-	91-59-8	3.0E-01	c	1.3E+00	c									3.9E-02	c	2.0E-04	
		1.0E-01	I						Napropamide	15299-99-7	6.3E+03	n	8.2E+04	n									1.6E+03	n	1.1E+01	
2.6E-04	C	1.1E-02	C	1.4E-05	C	1	0.1		Nickel Acetate	373-02-4	6.7E+02	n	8.1E+03	n	1.1E-02	c**	4.7E-02	c**	2.2E+02	n						
2.6E-04	C	1.1E-02	C	1.4E-05	C	1	0.1		Nickel Carbonate	3333-67-3	6.7E+02	n	8.1E+03	n	1.1E-02	c**	4.7E-02	c**	2.2E+02	n						
2.6E-04	C	1.1E-02	C	1.4E-05	C	V	1		Nickel Carbonyl	13463-39-3	8.2E+02	n	1.1E+04	n	1.1E-02	c**	4.7E-02	c**	2.2E+02	c**					c**	
2.6E-04	C	1.1E-02	C	1.4E-05	C		0.04		Nickel Hydroxide	12054-48-7	8.2E+02	n	1.1E+04	n	1.1E-02	c**	4.7E-02	c**	2.0E+02	n					n	
2.6E-04	C	1.1E-02	C	2.0E-05	C		0.04		Nickel Oxide	1313-99-1	8.4E+02	n	1.2E+04	n	1.1E-02	c**	4.7E-02	c**	2.0E+02	n					n	
2.4E-04	I	1.1E-02	C	1.4E-05	C		0.04		Nickel Refinery Dust	NA	8.2E+02	n	1.1E+04	n	1.2E-02	c**	5.1E-02	c**	2.2E+02	n					3.2E+01	
2.6E-04	C	2.0E-02	I	9.0E-05	A		0.04		Nickel Soluble Salts	7440-02-0	1.5E+03	n	2.2E+04	n	1.1E-02	c**	4.7E-02	c**	3.9E+02	n					2.6E+01	
1.7E+00	C	4.8E-04	I	1.1E-02	C	1.4E-05	C	0.04	Nickel Sulfide	12035-72-2	4.1E-01	c	1.9E+00	c	5.8E-03	c**	2.6E-02	c**	4.5E-02	c					c	
2.6E-04	C	1.1E-02	C	1.4E-05	C	1	0.1		Nickelocene	1271-28-9	6.7E+02	n	8.1E+03	n	1.1E-02	c**	4.7E-02	c**	2.2E+02	n					n	
		1.6E+00	I						Nitrate	14787-55-8	1.3E+05	nm	1.9E+06	nm									3.2E+04	n	1.0E+04	
									Nitrate+Nitrite (as N)	NA													1.0E+04	n	1.0E+04	
1.0E-01	I								Nitrite	14787-86-0	7.8E+03	n	1.2E+05	nm									2.0E+03	n		
1.0E-02	X	5.0E-05	X			1	0.1		Nitroanthine, 2-	88-74-4	6.3E+02	n	8.0E+03	n	5.2E-02	n	2.2E-01	n	1.9E+02	n					8.0E-02	
4.0E-03	P	6.0E-03	P			1	0.1		Nitroaniline, 4-	100-01-6	2.7E+01	c**	1.1E+02	c*	6.3E+00	n	2.6E+01	n	3.8E+00	c*					1.6E-03	
4.0E-05	I	2.0E-03	I	9.0E-03	I	V	1	3.1E+03	Nitrobenzene	98-95-3	5.1E+00	c*	2.2E+01	c*	7.0E-02	c	3.1E-01	c	1.4E-01	c*					9.2E-05	
		3.0E+03	P					0.1	Nitrocellulose	9004-70-0	1.9E+06	nm	2.5E+09	nm									6.0E+07	n	1.3E+04	
		7.0E-02	H					0.1	Nitrofurantoin	67-20-9	4.4E+03	n	5.7E+04	n									1.4E+03	n	6.1E-01	
1.3E+00	C	3.7E-04	C						Nitrofurazone	59-87-0	4.2E-01	c	1.8E+00	c	7.6E-03	c	3.3E-02	c	6.0E-02	c					5.4E-05	
1.7E-02	P	1.0E-04	P						Nitroglycerin	55-63-0	6.3E+00	n	8.2E+01	n									2.0E+00	n	8.5E-04	
		1.0E-01	I						Nitroguanidine	596-88-7	6.3E+03	n	8.2E+04	n									2.0E+03	n	4.8E-01	
8.8E-06	P			5.0E-03	P	V	1	1.8E+04	Nitromethane	75-52-5	5.4E+00	c*	2.4E+01	c*	3.2E-01	c*	1.4E+00	c*	6.4E-01	c*					1.4E-04	
2.7E-03	H			2.0E-02	I	V	1	4.9E+03	Nitropropane, 2-	79-46-9	1.4E-02	c	6.0E-02	c	1.0E-03	c	4.5E-03	c	2.1E-03	c					5.4E-07	
7.0E+00	C	7.7E-03	C						Nitroso-N-ethylurea, N-	759-73-9	4.5E-03	c	8.5E-02	c	1.3E-04	c	1.6E-03	c	9.2E-04	c					2.2E-07	
1.2E+02	C	3.4E-02	C					0.1	Nitroso-N-methylurea, N-	684-93-5	1.0E-03	c	1.9E-02	c	3.0E-05	c	3.6E-04	c	2.1E-04	c					4.6E-08	
5.4E+00	I	1.6E-03	I						Nitroso-di-N-butylamine, N-	924-16-3	9.9E-02	c	4.6E-01	c	1.8E-03	c	7.7E-03	c	2.7E-03	c					5.5E-06	
7.0E+00	I	2.0E-03	C						Nitroso-di-N-propylamine, N-	621-64-7	7.8E-02	c	3.3E-01	c	1.4E-03	c	6.1E-03	c	1.1E-02	c					8.1E-06	
2.8E+00	I	8.0E-04	C						Nitrosodiethanolamine, N-	1116-54-7	1.9E-01	c	8.2E-01	c	3.5E-03	c	1.5E-02	c	2.8E-02	c					5.6E-06	
1.5E+02	I	4.3E-02	I						Nitrosodithylamine, N-	55-19-5	2.1E-04	c	1.5E-02	c	2.4E-05	c	2.9E-04	c	1.7E-04	c					6.0E-08	
5.1E+01	I	1.4E-02	I	8.0E-06	P	4.0E-05	X	V	M	62-75-9	2.0E-03	c	3.4E-02	c	7.2E-05	c	8.8E-04	c	1.1E-04	c					2.8E-08	
4.9E-03	I	2.6E-06	C						Nitrosodiphenylamine, N-	86-30-6	1.1E+02	c	4.7E+02	c	1.1E+00	c	4.7E+00	c	1.2E+01	c					6.6E-02	
2.2E+01	I	6.3E-03	C						Nitrosomethylthylamine, N-	10595-95-6	2.0E-02	c	9.1E-02	c	4.5E-04	c	1.9E-03	c	7.1E-04	c					2.0E-07	
6.7E+00	C	1.9E-03	C						Nitrosomorpholine [N-]	59-89-2	8.1E-02	c	3.4E-01	c	1.5E-03	c	6.5E-03	c	1.2E-02	c					2.8E-06	
9.4E+00	C	2.7E-03	C						Nitrosopiperidine [N-]	100-75-4	5.8E-02	c	2.4E-01	c	1.0E-03	c	4.5E-03	c	8.2E-03	c					4.4E-06	
2.1E+00	I	6.1E-04	I						Nitrosopyrrolidine, N-	930-55-2	2.6E-01	c	1.1E+00	c	4.6E-03	c	2.0E-02	c	3.7E-02	c					1.4E-05	
		1.0E-04	X						Nitrotoluene, m-	99-08-1	6.3E+00	n	8.2E+01	n									1.7E+00	n	1.6E-03	
2.2E-01	P	9.0E-04	P						Nitrotoluene, o-	88-72-2	3.2E+00	c*	1.5E+01	c*									3.1E-01	c*	2.9E-04	
1.6E-02	P	4.0E-03	P						Nitrotoluene, p-	99-99-0	3.4E+01	c**	1.4E+02	c*												

Toxicity and Chemical-specific Information													Contaminant		Screening Levels								Protection of Ground Water SSLs					
SFO (mg/kg-day) ⁻¹	ke	IUR (ug/m ³) ⁻¹	ke	RfD _o (mg/kg-day)	ke	RfC _i (mg/m ³) ⁻¹	ke	Vo	muta-gen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)
		6.0E-03	H							1	0.1		Parathion	56-38-2	3.8E+02	n	4.9E+03	n					8.6E+01	n		4.3E-01	n	
		5.0E-02	H						V	1			Pebulate	1114-71-2	3.9E+03	n	5.8E+04	n					5.6E+02	n		4.6E-01	n	
		4.0E-02	I							1	0.1		Pendimethalin	40487-42-1	2.5E+03	n	3.3E+04	n					1.8E+02	n		2.1E+00	n	
		2.0E-03	I							1	0.1		Pentabromodiphenyl Ether	32534-81-9	1.3E+02	n	1.6E+03	n					4.0E+01	n		1.7E+00	n	
		1.0E-04	I							1	0.1		Pentabromodiphenyl ether, 2,2',4,4',5-(BDE-99)	60348-60-9	6.3E+00	n	8.2E+01	n					2.0E+00	n		8.7E-02	n	
		8.0E-04	I						V	1			Pentachlorobenzene	608-93-5	6.3E+01	n	9.3E+02	n					3.2E+00	n		2.4E-02	n	
9.0E-02	P											4.5E+02	Pentachloroethane	76-01-7	7.7E+00	c	3.6E+01	c					6.4E-01	c		3.1E-04	c	
2.6E-01	H			3.0E-03	I				V	1			Pentachloronitrobenzene	82-68-8	2.7E+00	c*	1.3E+01	c					1.2E-01	c		1.4E-03	c	
4.0E-01	I	5.1E-06	C	5.0E-03	I					1	0.25		Pentachlorophenol	87-86-5	1.0E+00	c	4.0E+00	c	5.5E-01	c	2.4E+00	c	4.0E-02	c	1.0E+00	4.0E-04	c	1.0E-02
4.0E-03	X			2.0E-03	P					1	0.1		Pentaerythritol tetranitrate (PETN)	78-11-5	1.3E+02	n	5.7E+02	c**					1.9E+01	c**		2.8E-02	c**	
						1.0E+00	P	V		1		3.9E+02	Pentane, n-	109-66-0	8.1E+02	ns	3.4E+03	ns	1.0E+03	n	4.4E+03	n	2.1E+03	n		1.0E+01	n	
													Perchlorates															
		7.0E-04	I							1			-Ammonium Perchlorate	7790-98-9	5.5E+01	n	8.2E+02	n					1.4E+01	n			n	
		7.0E-04	I							1			-Lithium Perchlorate	7791-03-9	5.5E+01	n	8.2E+02	n					1.4E+01	n			n	
		7.0E-04	I							1			-Perchlorate and Perchlorate Salts	14797-73-0	5.5E+01	n	8.2E+02	n					1.4E+01	n	1.5E+01(F)		n	
		7.0E-04	I							1			-Potassium Perchlorate	7778-74-7	5.5E+01	n	8.2E+02	n					1.4E+01	n			n	
		7.0E-04	I							1			-Sodium Perchlorate	7601-89-0	5.5E+01	n	8.2E+02	n					1.4E+01	n			n	
		2.0E-02	P						V	1			Perfluorobutane Sulfonate	375-73-5	1.6E+03	n	2.3E+04	n					3.8E+02	n		2.1E-01	n	
2.2E-03	C	6.3E-07	C	5.0E-02	I					1	0.1		Permethrin	52645-53-1	3.2E+03	n	4.1E+04	n					1.0E+03	n		2.4E+02	n	
										1	0.1		Phenacetin	62-44-2	2.5E+02	c	1.0E+03	c	4.5E+00	c	1.9E+01	c	3.4E+01	c		9.7E-03	c	
		2.5E-01	I							1	0.1		Phenmedipham	13684-63-4	1.6E+04	n	2.1E+05	nm					4.0E+03	n		2.1E+01	n	
		3.0E-01	I	2.0E-01	C					1	0.1		Phenol	108-95-2	1.9E+04	n	2.5E+05	nm	2.1E+02	n	8.8E+02	n	5.8E+03	n		3.3E+00	n	
		5.0E-04	X							1	0.1		Phenothiazine	92-84-2	3.2E+01	n	4.1E+02	n					4.3E+00	n		1.4E-02	n	
4.7E-02	H			6.0E-03	I					1	0.1		Phenylenediamine, m-	108-45-2	3.8E+02	n	4.9E+03	n					1.2E+02	n		3.2E-02	n	
										1	0.1		Phenylenediamine, o-	95-54-5	1.2E+01	c	4.9E+01	c					1.6E+00	c		4.4E-04	c	
		1.9E-01	H							1	0.1		Phenylenediamine, p-	108-50-3	1.2E+04	n	1.6E+05	nm					3.8E+03	n		1.0E+00	n	
1.9E-03	H									1	0.1		Phenylphenol, 2-	90-43-7	2.8E+02	c	1.2E+03	c					3.0E+01	c		4.0E-01	c	
		2.0E-04	H							1	0.1	1.6E+03	Phorate	298-02-2	1.3E+01	n	1.6E+02	n					3.0E+00	n		3.4E-03	n	
						3.0E-04	I	V		1			Phosgene	75-44-5	3.1E-01	n	1.3E+00	n	3.1E-01	n	1.3E+00	n						
		2.0E-02	I							1	0.1		Phosmet	732-11-8	1.3E+03	n	1.6E+04	n					3.7E+02	n		8.2E-02	n	
		4.9E+01	P							1			Phosphates, inorganic															
										1			-Aluminum metaphosphate	13776-88-0	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Ammonium polyphosphate	68333-79-9	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Calcium pyrophosphate	7790-76-3	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Diammonium phosphate	7783-28-0	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Dicalcium phosphate	7757-93-9	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Dimagnesium phosphate	7782-75-4	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Dipotassium phosphate	7758-11-4	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Disodium phosphate	7658-79-4	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Monoaluminum phosphate	13530-50-2	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Monoammonium phosphate	1122-16-1	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Monocalcium phosphate	7758-23-8	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Monomagnesium phosphate	7757-86-0	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Monopotassium phosphate	1118-11-0	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Monosodium phosphate	7558-80-7	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Polyphosphoric acid	8017-16-1	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Potassium tripolyphosphate	13845-36-8	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Sodium acid pyrophosphate	7758-16-9	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Sodium aluminum phosphate (acidic)	7785-88-8	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Sodium aluminum phosphate (anhydrous)	10279-59-1	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Sodium aluminum phosphate (tetrahydrate)	10305-76-7	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Sodium hexametaphosphate	10124-56-8	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Sodium polyphosphate	68915-31-1	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Sodium trimetaphosphate	7785-84-4	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P							1			-Sodium tripolyphosphate	7758-29-4	3.8E+06	nm	5.7E+07	nm					9.7E+05	n			n	
		4.9E+01	P																									

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information											Contaminant		Screening Levels							Protection of Ground Water SSLs						
SFO (mg/kg-day) ⁻¹	ke IUR (ug/m ³) ⁻¹	ke RfD _o (mg/kg-day)	ke RfC _i (mg/m ³)	ke RfC _v (mg/m ³)	mutagen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)		
1.4E-02	I	2.4E-06	C	2.0E-02	I			1	0.1	Phthalates																
				1.0E+00	I			1	0.1	-Bis(2-ethylhexyl)phthalate	117-81-7	3.9E+01	c*	1.6E+02	c	1.2E+00	c	5.1E+00	c	5.6E+00	c*	6.0E+00	1.3E+00	c*	1.4E+00	
					I			1	0.1	-Butylphthalyl Butylglycolate	85-70-1	6.3E+04	n	8.2E+05	nm					1.3E+04	n		3.0E+02	n		
				1.0E-01	I			1	0.1	-Dibutyl Phthalate	84-74-2	6.3E+03	n	8.2E+04	n					9.0E+02	n		2.3E+00	n		
				8.0E-01	I			1	0.1	-Diethyl Phthalate	84-66-2	5.1E+04	n	6.6E+05	nm					1.5E+04	n		6.1E+00	n		
				1.0E-01	I		V	1	0.1	-Dimethylterephthalate	120-61-6	7.8E+03	n	1.2E+05	nm					1.9E+03	n		4.9E-01	n		
				1.0E-02	P			1	0.1	-Octyl Phthalate, di-N-	117-84-0	6.3E+02	n	8.2E+03	n					2.0E+02	n		5.7E+01	n		
				1.0E+00	H			1	0.1	-Phthalic Acid, P-	100-21-0	6.3E+04	n	8.2E+05	nm					1.9E+04	n		6.8E+00	n		
				2.0E+00	I	2.0E-02	C	1	0.1	-Phthalic Anhydride	85-44-9	1.3E+05	nm	1.6E+06	nm	2.1E+01	n	8.8E+01	n	3.9E+04	n		8.5E+00	n		
				7.0E-02	I			1	0.1	Picloram	1918-02-1	4.4E+03	n	5.7E+04	n					1.4E+03	n	5.0E+02	3.8E-01	n	1.4E-01	
				1.0E-04	X			1	0.1	Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3	6.3E+00	n	8.2E+01	n					2.0E+00	n		1.3E-03	n		
				1.0E-02	I			1	0.1	Pyrimphos, Methyl	29232-93-7	6.3E+02	n	8.2E+03	n					1.2E+02	n		1.2E-01	n		
3.0E+01	C	8.6E-03	C	7.0E-06	H			1	0.1	Polybrominated Biphenyls	59536-65-1	1.8E-02	c*	7.7E-02	c*	3.3E-04	c	1.4E-03	c	2.6E-03	c*			c*		
				7.0E-02	S	2.0E-05	S			Polychlorinated Biphenyls (PCBs)																
					I		V	1	0.14	-Aroclor 1016	12674-11-2	4.1E+00	n	2.7E+01	c**	1.4E-01	c	6.1E-01	c	2.2E-01	c**		2.1E-02	c**		
2.0E+00	S	5.7E-04	S							-Aroclor 1221	11104-28-2	1.7E-01	c	7.2E-01	c	4.9E-03	c	2.1E-02	c	4.6E-03	c		7.9E-05	c		
2.0E+00	S	5.7E-04	S							-Aroclor 1232	11411-16-5	1.7E-01	c	7.2E-01	c	4.9E-03	c	2.1E-02	c	4.6E-03	c		7.9E-05	c		
2.0E+00	S	5.7E-04	S							-Aroclor 1242	53469-21-9	2.3E-01	c	9.7E-01	c	4.9E-03	c	2.1E-02	c	7.8E-03	c		1.2E-03	c		
2.0E+00	S	5.7E-04	S							-Aroclor 1248	12672-29-6	2.3E-01	c	9.4E-01	c	4.9E-03	c	2.1E-02	c	7.8E-03	c		1.2E-03	c		
2.0E+00	S	5.7E-04	S	2.0E-05	I			1	0.14	-Aroclor 1254	11097-69-1	2.4E-01	c**	9.7E-01	c*	4.9E-03	c	2.1E-02	c	7.8E-03	c*		2.0E-03	c*		
2.0E+00	S	5.7E-04	S							-Aroclor 1260	11096-82-5	2.4E-01	c	9.9E-01	c	4.9E-03	c	2.1E-02	c	7.8E-03	c		5.5E-03	c		
				6.0E-04	X			1	0.14	-Aroclor 5460	11126-42-4	3.5E+01	n	4.4E+02	n					1.2E+01	n		2.0E+00	n		
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1	0.14	-Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	1.2E-01	c*	5.1E-01	c*	2.5E-03	c	1.1E-02	c	4.0E-03	c		2.8E-03	c	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1	0.14	-Hexachlorobiphenyl, 2,3,4,4',5,5'- (PCB 167)	52663-72-6	1.2E-01	c*	5.1E-01	c*	2.5E-03	c	1.1E-02	c	4.0E-03	c		1.7E-03	c	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1	0.14	-Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	1.2E-01	c*	5.1E-01	c*	2.5E-03	c	1.1E-02	c	4.0E-03	c		1.7E-03	c	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1	0.14	-Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 156)	38380-08-4	1.2E-01	c*	5.1E-01	c*	2.5E-03	c	1.1E-02	c	4.0E-03	c		1.7E-03	c	
3.9E+03	E	1.1E+00	E	2.3E-08	E	1.3E-06	E	V	1	0.14	-Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	1.2E-04	c*	5.1E-04	c*	2.5E-06	c	1.1E-05	c	4.0E-06	c		1.7E-06	c	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1	0.14	-Pentachlorobiphenyl, 2,3,4,4',5'- (PCB 123)	65610-44-3	1.2E-01	c*	5.0E-01	c*	2.5E-03	c	1.1E-02	c	4.0E-03	c		1.0E-03	c	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1	0.14	-Pentachlorobiphenyl, 2,3',4,4',5'- (PCB 118)	31588-00-6	1.2E-01	c*	5.0E-01	c*	2.5E-03	c	1.1E-02	c	4.0E-03	c		1.0E-03	c	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1	0.14	-Pentachlorobiphenyl, 2,3,3',4,4',5'- (PCB 105)	32598-14-4	1.2E-01	c*	5.0E-01	c*	2.5E-03	c	1.1E-02	c	4.0E-03	c		1.0E-03	c	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1	0.14	-Pentachlorobiphenyl, 2,3,4,4',5'- (PCB 114)	74472-37-0	1.2E-01	c*	5.0E-01	c*	2.5E-03	c	1.1E-02	c	4.0E-03	c		1.0E-03	c	
1.3E+04	E	3.8E+00	E	7.0E-09	E	4.0E-07	E	V	1	0.14	-Pentachlorobiphenyl, 3,3',4,4',5'- (PCB 126)	57465-28-8	3.7E-05	c*	1.5E-04	c*	7.4E-07	c	3.2E-06	c	1.2E-06	c		3.0E-07	c	
2.0E+00	I	5.7E-04	I					1	0.14	-Polychlorinated Biphenyls (high risk)	1336-36-3	2.3E-01	c	9.7E-01	c	4.9E-03	c	2.1E-02	c							
4.0E-01	I	1.0E-04	I					1	0.14	-Polychlorinated Biphenyls (low risk)	1336-36-3					2.8E-02	c	1.2E-01	c	4.4E-02	c	5.0E-01	6.8E-03	c	7.8E-02	
7.0E-02	I	2.0E-05	I					1	0.14	-Polychlorinated Biphenyls (lowest risk)	1336-36-3					1.4E-01	c	6.1E-01	c							
1.3E+01	E	3.8E-03	E	7.0E-06	E	4.0E-04	E	V	1	0.14	-Tetrachlorobiphenyl, 3,3',4,4' (PCB 77)	32598-13-3	3.8E-02	c*	1.6E-01	c*	7.4E-04	c	3.2E-03	c	6.0E-03	c*		9.4E-04	c*	
3.9E+01	E	1.1E-02	E	2.3E-06	E	1.3E-04	E	V	1	0.14	-Tetrachlorobiphenyl, 3,4,4',5'- (PCB 81)	70362-50-4	1.2E-02	c*	4.9E-02	c*	2.5E-04	c	1.1E-03	c	4.0E-04	c		6.2E-05	c	
				6.0E-04	I			1	0.1	Polymeric Methylene Diethylene Disocyanate (PMDI)	9016-87-9	8.5E+05	nm	3.6E+06	nm	6.3E-01	n	2.6E+00	n							
				6.0E-02	I		V	1	0.13	Polymeric Aromatic Hydrocarbons (PAHs)																
				3.0E-01	I		V	1	0.13	-Acenaphthene	83-32-9	3.6E+03	n	4.5E+04	n					5.3E+02	n		5.5E+00	n		
7.3E-01	E	1.1E-04	C					1	0.13	-Anthracene	120-12-7	1.8E+04	n	2.3E+05	nm					1.8E+03	n		5.8E+01	n		
								1	0.13	-Benz[a]anthracene	56-55-3	1.6E-01	c	2.9E+00	c	9.2E-03	c	1.1E-01	c	1.2E-02	c		4.3E-03	c		
1.2E+00	C	1.1E-04	C					1	0.13	-Benzo[j]fluoranthene	205-85-2	4.2E-01	c	1.8E+00	c	2.6E-02	c	1.1E-01	c	6.5E-02	c		7.8E-02	c		
7.3E+00	I	1.1E-03	C				M	1	0.13	-Benzo[a]pyrene	50-32-8	1.6E-02	c	2.9E-01	c	9.2E-04	c	1.1E-02	c	3.4E-03	c	2.0E-01	4.0E-03	c	2.4E-01	
7.3E-01	E	1.1E-04	C				M	1	0.13	-Benzo[b]fluoranthene	205-99-2	1.6E-01	c	2.9E+00	c	9.2E-03	c	1.1E-01	c	3.4E-02	c		4.1E-02	c		
7.3E-02	E	1.1E-04	C				M	1	0.13	-Benzo[k]fluoranthene	207-08-9	1.6E+00	c	2.9E+01	c	9.2E-03	c	1.1E-01	c	3.4E-01	c		4.0E-01	c		
7.3E-03	E	1.1E-05	C	8.0E-02	I		V	1	0.13	-Chloronaphthalene, Beta-	91-58-7	4.8E+03	n	6.0E+04	n					7.5E+02	n		3.8E+00	n		
							M	1	0.13	-Chrysene	2181-01-9	1.6E+01	c	2.9E+02	c	9.2E-02	c	1.1E+00	c	3.4E+00	c		1.2E+00	c		
7.3E+00	E	1.2E-03	C				M	1	0.13	-Dibenzo[a,h]anthracene	53-70-3	1.6E-02	c	2.9E-01	c	8.4E-04	c	1.0E-02	c	3.4E-03	c		1.3E-02	c		
1.2E+01	C	1.1E-03	C				M	1	0.13	-Dibenzo[a,e]pyrene	192-65-4	4.2E-02	c	1.8E-01	c	2.6E-03	c	1.1E-02	c	6.5E-03	c		8.4E-02	c		
2.5E+02	C	7.1E-02	C				M	1	0.13	-Dimethylbenz(a)anthracene																

Regional Screening Level (RSL) Summary Table (TR-1E-6, HQ=1) June 2015 (revised)

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Screening Levels								Protection of Ground Water SSLs					
SFO (mg/kg-day) ⁻¹	ke IUR (ug/m ³ -y) ⁻¹	ke RfD _o (mg/kg-day)	ke RfC _i (mg/m ³ -y)	ke RfC _v (mg/m ³ -y)	o mutagen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)	
		5.0E-03	I					0.1	Propanil	709-98-8	3.2E+02	n	4.1E+03	n					8.2E+01	n		4.5E-02	n		
		2.0E-02	I					0.1	Propargite	2312-35-8	1.3E+03	n	1.6E+04	n					1.6E+02	n		1.2E+01	n		
		2.0E-03	I		V			1.1E+05	Propargyl Alcohol	107-19-7	1.6E+02	n	2.3E+03	n					4.0E+01	n		8.1E-03	n		
		2.0E-02	I					0.1	Propazine	139-40-2	1.3E+03	n	1.6E+04	n					3.4E+02	n		3.0E-01	n		
		2.0E-02	I					0.1	Propam	122-42-9	1.3E+03	n	1.6E+04	n					3.5E+02	n		2.2E-01	n		
		1.3E-02	I					0.1	Propiconazole	60207-90-1	8.2E+02	n	1.1E+04	n					2.1E+02	n		6.9E-01	n		
				8.0E-03	I	V			3.3E+04	Propionaldehyde	123-38-6	7.5E+01	n	3.1E+02	n	8.3E+00	n	3.5E+01	n	1.7E+01	n		3.4E-03	n	
		1.0E-01	X	1.0E+00	X	V			2.6E+02	Propyl benzene	103-65-1	3.8E+03	ns	2.4E+04	ns	1.0E+03	n	4.4E+03	n	6.6E+02	n		1.2E+00	n	
				3.0E+00	C	V			3.5E+02	Propylene	115-07-1	2.2E+03	ns	9.3E+03	ns	3.1E+03	n	1.3E+04	n	6.3E+03	n		6.0E+00	n	
		2.0E+01	P					0.1	Propylene Glycol	57-55-6	1.3E+06	nm	1.6E+07	nm					4.0E+05	n		8.1E+01	n		
				2.7E-04	A			0.1	Propylene Glycol Dinitrate	6423-43-4	3.9E+05	nm	1.6E+06	nm	2.8E-01	n	1.2E+00	n							
		7.0E-01	H		V				8.5E+04	Propylene Glycol Monoethyl Ether	1569-02-4	5.5E+04	n	8.2E+05	nms					1.4E+04	n		2.8E+00	n	
		7.0E-01	H	2.0E+00	I	V			1.1E+05	Propylene Glycol Monomethyl Ether	107-98-2	4.1E+04	n	3.7E+05	nms	2.1E+03	n	8.8E+03	n	3.2E+03	n		6.5E-01	n	
2.4E-01	I	3.7E-06	I			3.0E-02	I	V		7.8E+04	Propylene Oxide	75-56-9	2.1E+00	c	9.7E+00	c	7.6E-01	c*	3.3E+00	c*	2.7E-01	c		5.6E-05	c
		2.5E-01	I					0.1	Pursuit	81335-77-5	1.6E+04	n	2.1E+05	nm					4.7E+03	n		4.1E+00	n		
		2.5E-02	I					0.1	Pyridin	51630-58-1	1.6E+03	n	2.1E+04	n					5.0E+02	n		3.2E-02	n		
		1.0E-03	I			V			5.3E+05	Pyridine	110-86-1	7.8E+01	n	1.2E+03	n					2.0E+01	n		6.8E-03	n	
		5.0E-04	I					0.1	Quinalphos	13593-03-8	3.2E+01	n	4.1E+02	n					5.1E+00	n		4.3E-02	n		
3.0E+00	I							0.1	Quinoline	91-22-5	1.8E-01	c	7.7E-01	c					2.4E-02	c		7.8E-05	c		
				3.0E-02	A			1	Refractory Ceramic Fibers	NA	4.3E+07	nm	1.8E+08	nm	3.1E+01	n	1.3E+02	n							
		3.0E-02	I					0.1	Resmethrin	10453-86-8	1.9F+03	n	2.5E+04	n					6.7E+01	n		4.2E+01	n		
		5.0E-02	H		V			1	Ronnel	299-94-3	3.9E+03	n	5.8E+04	n					4.1E+02	n		3.7E+00	n		
2.2E-01	C	6.3E-05	C					0.1	Rotenone	83-79-4	2.5E+02	n	3.3E+03	n					6.1E+01	n		3.2E+01	n		
						M		0.1	Safrole	94-59-7	5.5E-01	c	1.0E+01	c	1.6E-02	c	1.9E-01	c	9.5E-02	c		5.9E-05	c		
		2.5E-02	I					0.1	Savay	78587-05-0	1.6E+03	n	2.1E+04	n					1.1E+02	n		5.0E-01	n		
		5.0E-03	I					1	Selenious Acid	7783-60-8	3.9E+02	n	5.8E+03	n					1.0E+02	n		5.2E-01	n	2.6E-01	
		5.0E-03	I	2.0E-02	C			1	Selenium	7782-49-2	3.9E+02	n	5.8E+03	n	2.1E+01	n	8.8E+01	n	1.0E+02	n	5.0E+01				
		5.0E-03	C	2.0E-02	C			1	Selenium Sulfide	7446-34-6	3.9E+02	n	5.8E+03	n	2.1E+01	n	8.8E+01	n	1.0E+02	n					
		9.0E-02	I					0.1	Sethoxydim	74051-80-2	5.7E+03	n	7.4E+04	n					1.0E+03	n		9.3E+00	n		
				3.0E-03	C			1	Silica (crystalline, respirable)	7631-86-9	4.3E+06	nm	1.8E+07	nm	3.1E+00	n	1.3E+01	n							
		5.0E-03	I			0.04			Silver	7440-22-4	3.9E+02	n	5.8E+03	n					9.4E+01	n		8.0E-01	n		
1.2E-01	H	5.0E-03	I					0.1	Simazine	122-34-9	4.5E+00	c*	1.9E+01	c					6.1E-01	c	4.0E+00	3.0E-04	c	2.0E-03	
		1.3E-02	I					0.1	Sodium Acifluorfen	62476-59-9	8.2E+02	n	1.1E+04	n					2.6E+02	n		2.1E+00	n		
		4.0E-03	I					1	Sodium Azide	26628-22-8	3.1E+02	n	4.7E+03	n					8.0E+01	n					
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M	0.025	Sodium Dichromate	10588-01-9	3.0E-01	c	6.2E+00	c	6.8E-06	c	8.2E-05	c	4.1E-02	c				
2.7E-01	H	3.0E-02	I					0.1	Sodium Diethylidithiocarbamate	148-18-5	2.0E+00	c	8.5E+00	c					2.9E-01	c					
		5.0E-02	A	1.3E-02	C			1	Sodium Fluoride	7681-49-4	3.9E+03	n	5.8E+04	n	1.4E+01	n	5.7E+01	n	1.0E+03	n					
		2.0E-05	I					0.1	Sodium Fluoroacetate	82-74-8	1.3E+00	n	1.6E+01	n					4.0E-01	n		8.1E-05	n		
		1.0E-03	H					1	Sodium Metavanadate	13718-26-8	7.8E+01	n	1.2E+03	n					2.0E+01	n					
2.4E-02	H	3.0E-02	I					0.1	Stirofos (Tetrachlorovinphos)	961-11-5	2.3E+01	c*	9.6E+01	c					2.8E+00	c		8.1E-03	c		
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M	0.025	Strontium Chromate	7789-06-2	3.0E-01	c	6.2E+00	c	6.8E-06	c	8.2E-05	c	4.1E-02	c				
		6.0E-01	I					1	Strontium, Stable	7440-24-6	4.7E+04	n	7.0E+05	nm					1.2E+04	n		4.2E+02	n		
		3.0E-04	I					0.1	Strychnine	57-24-9	1.9E+01	n	2.5E+02	n					5.9E+00	n		6.5E-02	n		
		2.0E-01	I	1.0E+00	I	V			8.7E+02	Styrene	100-42-5	6.0E+03	ns	3.5E+04	ns	1.0E+03	n	4.4E+03	n	1.2E+03	n	1.0E+02	1.3E+00	n	1.1E-01
		3.0E-03	P					0.1	Styrene-Acrylonitrile (SAN) Trimer	NA	1.9E+02	n	2.5E+03	n					4.8E+01	n					
		1.0E-03	P	2.0E-03	X			0.1	Sulfolane	126-33-0	6.3E+01	n	8.2E+02	n	2.1E+00	n	8.8E+00	n	2.0E+01	n		4.4E-03	n		
		8.0E-04	P					0.1	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9	5.1E+01	n	6.6E+02	n					1.1E+01	n		6.5E-02	n		
				1.0E-03	C	V		1	Sulfur Trioxide	7446-11-9	1.4E+06	nm	6.0E+06	nm	1.0E+00	n	4.4E+00	n	2.1E+00	n					
				1.0E-03	C			1	Sulfuric Acid	7664-93-9	1.4E+06	nm	6.0E+06	nm	1.0E+00	n	4.4E+00	n							
		2.5E-02	I					0.1	Systhane	88671-89-0	1.6E+03	n	2.1E+04	n					4.5E+02	n		5.6E+00	n		
		3.0E-02	H					1	TCMTB	21564-17-0	1.9E+03	n	2.5E+04	n					4.8E+02	n		3.3E+00	n		
		7.0E-02	I					0.1	Tebuthiuron	34014-18-1	4.4E+03	n	5.7E+04	n					1.4E+03	n		3.9E-01	n		
		2.0E-02	H					0.1	Temephos	3383-96-8	1.3E+03	n	1.6E+04	n					4.0E+02	n		7.6E+01	n		
		1.3E-02	I					0.1	Terbacil	5902-51-2	8.2E+02	n	1.1E+04	n					2.5E+02	n		7.5E-02	n		
		2.5E-05	H		V			3.1E+01	Terbufos	13071-79-9	2.0E+00	n	2.9E+01	n					2.4E-01	n		5.2E-04	n		
		1.0E-03	I					0.1	Terbutryn	886-50-0	6.3E+01	n	8.2E+02	n					1.3E+01	n		1.9E-02	n		
		1.0E-04	I					0.1	Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1	6.3E+00	n	8.2E+01	n					2.0E+00	n		5.3E-02	n		
		3.0E-04	I		V			1	Tetrachlorobenzene, 1,2,																

Regional Screening Level (RSL) Summary Table (TR-1E-6, HQ=1) June 2015 (revised)

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Screening Levels								Protection of Ground Water SSLs					
SFO (mg/kg-day) ⁻¹	ke IUR (ug/m ³ -y) ⁻¹	ke RfD _o (mg/kg-day)	ke RfC _i (mg/m ³ -y)	ke Vo	muta-gen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)	
		2.0E-03	P			1	0.0007		Tetryl (Trinitrophenylmethyl nitramine)	479-45-8	1.6E+02	n	2.3E+03	n						3.9E+01	n		3.7E-01	n	
		7.0E-06	X			1			Thallium (I) Nitrate	10102-45-1	5.5E-01	n	8.2E+00	n						1.4E-01	n			n	
		1.0E-05	X			1			Thallium (Soluble Salts)	7440-28-0	7.8E-01	n	1.2E+01	n						2.0E-01	n	2.0E+00	1.4E-02	n	1.4E-01
		6.0E-06	X		V	1	0.1		Thallium Acetate	563-68-8	3.8E-01	n	4.9E+00	n						1.2E-01	n			n	
		2.0E-05	X			1	0.1		Thallium Carbonate	6533-73-9	1.3E+00	n	1.6E+01	n						4.0E-01	n			n	
		6.0E-06	X			1			Thallium Chloride	7791-12-0	4.7E-01	n	7.0E+00	n						1.2E-01	n			n	
		2.0E-05	X			1			Thallium Sulfate	7446-18-6	1.6E+00	n	2.3E+01	n						4.0E-01	n			n	
		1.0E-02	I			1	0.1		Thiobencarb	28249-77-6	6.3E+02	n	8.2E+03	n						1.6E+02	n		5.5E-01	n	
		7.0E-02	X			1	0.0075		Thiodiglycol	111-48-8	5.4E+03	n	7.9E+04	n						1.4E+03	n		2.8E-01	n	
		3.0E-04	H			1	0.1		Thiofanox	39196-18-4	1.9E+01	n	2.5E+02	n						5.3E+00	n		1.8E-03	n	
		8.0E-02	I			1	0.1		Thiophanate, Methyl	23564-05-8	5.1E+03	n	6.6E+04	n						1.6E+03	n		1.4E+00	n	
		5.0E-03	I			1	0.1		Thiram	137-26-8	3.2E+02	n	4.1E+03	n						9.8E+01	n		1.4E-01	n	
		6.0E-01	H			1			Tin	7440-31-5	4.7E+04	n	7.0E+05	nm	1.0E-01	n	4.4E-01	n		1.2E+04	n		3.0E+03	n	
				1.0E-04	A	V	1		Titanium Tetrachloride	7750-45-0	1.4E+05	nm	6.0E+05	nm	1.0E-01	n	4.4E-01	n		2.1E-01	n			n	
		8.0E-02	I	5.0E+00	I	V	1	8.2E+02	Toluene	108-88-3	4.9E+03	ns	4.7E+04	ns	5.2E+03	n	2.2E+04	n		1.1E+03	n	1.0E+03	7.6E-01	n	6.9E-01
1.8E-01	X	2.0E-04	X			1	0.1		Toluene-2,5-diamine	95-70-5	3.0E+00	c**	1.3E+01	c*						4.3E-01	c**		1.3E-04	c**	
3.0E-02	P	4.0E-03	X			1	0.1		Toluidine, p-	106-49-0	1.8E+01	c*	7.7E+01	c*						2.5E+00	c*		1.1E-03	c*	
		3.0E+00	P		V	1		3.4E-01	Total Petroleum Hydrocarbons (Aliphatic High)	NA	2.3E+05	nms	3.5E+06	nms						6.0E+04	n		2.4E+03	n	
				6.0E-01	P	V	1	1.4E+02	Total Petroleum Hydrocarbons (Aliphatic Low)	NA	5.2E+02	ns	2.2E+03	ns	6.3E+02	n	2.6E+03	n		1.3E+03	n		8.8E+00	n	
		1.0E-02	X	1.0E-01	P	V	1	6.9E+00	Total Petroleum Hydrocarbons (Aliphatic Medium)	NA	9.6E+01	ns	4.4E+02	ns	1.0E+02	n	4.4E+02	n		1.0E+02	n		1.5E+00	n	
		4.0E-02	P			1	0.1		Total Petroleum Hydrocarbons (Aromatic High)	NA	2.5E+03	n	3.3E+04	n						8.0E+02	n		8.9E+01	n	
		4.0E-03	P	3.0E-02	P	V	1	1.8E+03	Total Petroleum Hydrocarbons (Aromatic Low)	NA	8.2E+01	n	4.2E+02	n	3.1E+01	n	1.3E+02	n		3.3E+01	n		1.7E-02	n	
		4.0E-03	P	3.0E-03	P	V	1		Total Petroleum Hydrocarbons (Aromatic Medium)	NA	1.1E+02	n	6.0E+02	n	3.1E+00	n	1.3E+01	n		5.5E+00	n		2.3E-02	n	
1.1E+00	I	3.2E-04	I			1	0.1		Toxaphene	8001-35-2	4.9E-01	c	2.1E+00	c	8.8E-03	c	3.8E-02	c		1.5E-02	c	3.0E+00	2.4E-03	c	4.6E-01
		7.5E-03	I			1	0.1		Tralometrin	66841-25-6	4.7E+02	n	6.2E+03	n						1.5E+02	n		5.8E+01	n	
		3.0E-04	A		V	1			Tri-n-butyltin	688-73-3	2.3E+01	n	3.5E+02	n						3.7E+00	n		8.2E-02	n	
		8.0E+01	X			1	0.1		Triacetin	102-76-1	5.1E+06	nm	6.6E+07	nm						1.6E+06	n		4.5E+02	n	
		1.3E-02	I		V	1			Triallate	2303-17-5	1.0E+03	n	1.5E+04	n						1.2E+02	n		2.6E-01	n	
		1.0E-02	I			1	0.1		Triasulfuron	82097-50-5	6.3E+02	n	8.2E+03	n						2.0E+02	n		2.1E-01	n	
		5.0E-03	I		V	1			Tribromobenzene, 1,2,4-	615-59-3	3.9E+02	n	5.8E+03	n						4.5E+01	n		6.4E-02	n	
9.0E-03	P	1.0E-02	P			1	0.1		Tributyl Phosphate	126-73-8	6.0E+01	c*	2.6E+02	c*						5.1E+00	c*		2.5E-02	c*	
		3.0E-04	P			1	0.1		Tributyltin Compounds	NA	1.9E+01	n	2.5E+02	n						6.0E+00	n			n	
		3.0E-04	I			1	0.1		Tributyltin Oxide	56-35-9	1.9E+01	n	2.5E+02	n						5.7E+00	n		2.9E+02	n	
7.0E-02	I	3.0E+01	H	V		1		9.1E+02	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1	4.0E+04	ns	1.7E+05	nms	3.1E+04	n	1.3E+05	n		5.5E+04	n		1.4E+02	n	
		2.0E-02	I			1	0.1		Trichloroacetic Acid	76-03-9	7.8E+00	c	3.3E+01	c						1.1E+00	c	6.0E+01	2.2E-04	c	1.2E-02
2.9E-02	H					1	0.1		Trichloroaniline-HCl, 2,4,6-	33683-50-2	1.9E+01	c	7.9E+01	c						2.7E+00	c		7.4E-03	c	
7.0E-03	X	3.0E-05	X			1	0.1		Trichloroaniline, 2,4,6-	634-93-5	1.9E+00	n	2.5E+01	n						4.0E-01	n		3.6E-03	n	
		8.0E-04	X		V	1			Trichlorobenzene, 1,2,3-	87-61-6	6.3E+01	n	9.3E+02	n						7.0E+00	n		2.1E-02	n	
2.9E-02	P	1.0E-02	I	2.0E-03	P	V	1	4.0E+02	Trichlorobenzene, 1,2,4-	120-82-1	2.4E+01	c**	1.1E+02	c**	2.1E+00	n	8.8E+00	n		1.1E+00	c**	7.0E+01	3.3E-03	c**	2.0E-01
		2.0E+00	I	5.0E+00	I	V	1	6.4E+02	Trichloroethane, 1,1,1-	71-55-6	8.1E+03	ns	3.6E+04	ns	5.2E+03	n	2.2E+04	n		8.0E+03	n	2.0E+00	2.8E+00	n	7.0E-02
5.7E-02	I	1.6E-05	I			1		2.2E+03	Trichloroethane, 1,1,2-	79-00-5	1.1E+00	c**	5.0E+00	c**	1.8E-01	c**	7.7E-01	c**		2.8E-01	c**	5.0E+00	8.9E-05	c**	1.6E-03
4.6E-02	I	4.1E-06	I	5.0E-04	I	2.0E-03	I	V	M	1	6.9E+02														
		3.0E-01	I	7.0E-01	H	V	1	1.2E+03	Trichloroethylene	79-01-6	9.4E-01	c**	6.0E+00	c**	4.8E-01	c**	3.0E+00	c**		4.9E-01	c**	5.0E+00	1.8E-04	c**	1.8E-03
		1.0E-01	I			1	0.1		Trichlorofluoromethane	75-69-4	7.3E+02	n	3.1E+03	ns	7.3E+02	n	3.1E+03	n		1.1E+03	n		7.3E-01	n	
		1.1E-02	I	3.1E-06	I				Trichlorophenol, 2,4,6-	95-95-4	6.3E+03	n	8.2E+04	n						1.2E+03	n		4.4E+00	n	
		1.0E-02	I			1	0.1		Trichlorophenol, 2,4,6-	88-06-2	4.9E+01	c**	2.1E+02	c**	9.1E-01	c	4.0E+00	c		4.0E+00	c**		1.5E-02	c**	
		1.0E-02	I			1	0.1		Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5	6.3E+02	n	8.2E+03	n						1.6E+02	n		6.7E-02	n	
		8.0E-03	I			1	0.1		Trichlorophenoxypropionic Acid, -2,4,5	93-72-1	5.1E+02	n	6.6E+03	n						1.1E+02	n	5.0E+01	6.1E-02	n	2.8E-02
3.0E+01	I	5.0E-03	I		V	1		1.3E+03	Trichloropropane, 1,1,2-	598-77-6	3.9E+02	n	5.8E+03	ns						8.8E+01	n		3.5E-02	n	
		4.0E-03	I	3.0E-04	I	V	M	1	1.4E+03	Trichloropropane, 1,2,3-	96-18-4	5.1E+03	c	1.1E-01	c	3.1E-01	n	1.3E+00	n		7.5E-04	c		3.2E-07	c
		3.0E-03	X	3.0E-04	P	V	1	4.5E+02	Trichloropropene, 1,2,3-	96-19-5	7.3E-01	n	3.1E+00	n	3.1E-01	n	1.3E+00	n		6.2E-01	n		3.1E-04	n	
		2.0E-02	A			1	0.1		Tricresyl Phosphate (TCP)	1330-78-5	1.3E+03	n	1.6E+04	n						1.6E+02	n		1.5E+01	n	
		3.0E-03	I			1	0.1		Tridiphane	58138-08-2	1.9E+02	n	2.5E+03	n						1.8E+01	n		1.3E-01	n	
				7.0E-03	I	V	1	2.8E+04	Triethylamine	121-44-8	1.2E+02	n	4.8E+02	n	7.3E+00	n	3.1E+01	n		1.5E+01	n		4.		

Regional Screening Level (RSL) Summary Table (TR=1E-6, HQ=1) June 2015 (revised)

Toxicity and Chemical-specific Information											Contaminant		Screening Levels										Protection of Ground Water SSLs				
SFO (mg/kg-day) ⁻¹	Key	IUR (ug/m ³ -y) ⁻¹	Key	RfD _o (mg/kg-day)	Key	RfC _i (mg/m ³ -y)	Key	muta- gen	GIABS	ABS	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)	Risk-based SSL (mg/kg)	key	MCL-based SSL (mg/kg)
2.0E-02	P			7.0E-03	P				1	0.1		Tris(2-chloroethyl)phosphate	115-96-8	2.7E+01	c*	1.1E+02	c*					3.8E+00	c*		3.8E-03	c*	
3.2E-03	P			1.0E-01	P				1	0.1		Tris(2-ethylhexyl)phosphate	78-42-2	1.7E+02	c*	7.2E+02	c					2.4E+01	c*		1.2E+02	c*	
				3.0E-03	I	4.0E-05	A		1			Uranium (Soluble Salts)	NA	2.3E+02	n	3.5E+03	n	4.2E-02	n	1.8E-01	n	6.0E+01	n	3.0E+01	2.7E+01	n	1.4E+01
1.0E+00	C	2.9E-04	C					M	1	0.1		Urethane	51-79-6	1.2E-01	c	2.3E+00	c	3.5E-03	c	4.2E-02	c	2.5E-02	c		5.6E-06	c	
		8.3E-03	P	9.0E-03	I	7.0E-06	P		0.026			Vanadium Peroxide	1314-62-1	4.6E+02	c**	2.0E+03	c**	3.4E-04	c*	1.5E-03	c*	1.5E+02	n			n	
				5.0E-03	S	1.0E-04	A		0.026			Vanadium and Compounds	7440-62-2	3.9E+02	n	5.8E+03	n	1.0E-01	n	4.4E-01	n	8.6E+01	n		8.6E+01	n	
				1.0E-03	I		V		1			Vermolate	1929-77-7	7.8E+01	n	1.2E+03	n					1.1E+01	n		8.9E-03	n	
				2.5E-02	I				1	0.1		Vinclozolin	50471-44-8	1.6E+03	n	2.1E+04	n					4.4E+02	n		3.4E-01	n	
				1.0E+00	H	2.0E-01	I	V	1		2.8E+03	Vinyl Acetate	108-05-4	9.1E+02	n	3.8E+03	ns	2.1E+02	n	8.8E+02	n	4.1E+02	n		8.7E-02	n	
		3.2E-05	H			3.0E-03	I	V	1		3.4E+03	Vinyl Bromide	593-60-2	1.2E-01	c*	5.2E-01	c*	8.8E-02	c*	3.8E-01	c*	1.8E-01	c*		5.1E-05	c*	
7.2E-01	I	4.4E-06	I	3.0E-03	I	1.0E-01	I	V	M	1	3.9E+03	Vinyl Chloride	75-01-4	5.9E-02	c	1.7E+00	c	1.7E-01	c	2.8E+00	c	1.9E-02	c	2.0E+00	6.5E-06	c	6.9E-04
				3.0E-04	I				1	0.1		Wartann	81-81-2	1.9E+01	n	2.5E+02	n					5.6E+00	n		5.9E-03	n	
				2.0E-01	S	1.0E-01	S	V	1		3.9E+02	Xylene, p	106-42-3	5.6E+02	ns	2.4E+03	ns	1.0E+02	n	4.4E+02	n	1.9E+02	n		1.9E-01	n	
				2.0E-01	S	1.0E-01	S	V	1		3.9E+02	Xylene, m	108-38-3	5.5E+02	ns	2.4E+03	ns	1.0E+02	n	4.4E+02	n	1.9E+02	n		1.9E-01	n	
				2.0E-01	S	1.0E-01	S	V	1		4.3E+02	Xylene, o	95-47-6	6.5E+02	ns	2.8E+03	ns	1.0E+02	n	4.4E+02	n	1.9E+02	n		1.9E-01	n	
				2.0E-01	I	1.0E-01	I	V	1		2.6E+02	Xylenes	1320-20-7	6.5E+02	ns	2.8E+03	ns	1.0E+02	n	4.4E+02	n	1.9E+02	n	1.0E+04	1.9E-01	n	9.8E+00
				3.0E-04	I				1			Zinc Phosphide	1314-84-7	2.3E+01	n	3.5E+02	n					6.0E+00	n			n	
				3.0E-01	I				1			Zinc and Compounds	7440-66-6	2.3E+04	n	3.5E+05	nm					6.0E+03	n		3.7E+02	n	
				5.0E-02	I				1	0.1		Zinc	72122-67-7	3.2E+03	n	4.1E+04	n					9.9E+02	n		2.9E+00	n	
				8.0E-05	X				1			Zirconium	7440-67-7	6.3E+00	n	9.3E+01	n					1.6E+00	n		4.8E+00	n	

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k e IUR (ug/m ³) ⁻¹	k e RfD _o (mg/kg-day)	k e RfC _i (mg/m ³)	k e v o c muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)			
1.8E-02	C	5.1E-06	C	1.5E-01	I	1.0E+00	1.0E-01	1.4E+09		ALAR	1596-84-5	3.9E+01	1.4E+02	7.5E+05	3.0E+01	1.2E+04	4.9E+04		9.5E+03			
8.7E-03	I	4.0E-03	I	4.0E-03	I	1.0E+00	1.0E-01	1.4E+09		Acephate	30560-19-1	8.0E+01	2.8E+02		6.2E+01	3.1E+02	1.3E+03		2.5E+02			
		2.2E-06	I		I	1.0E+00		1.1E+05	8.7E+03	Acetaldehyde	75-07-0			1.1E+01			8.2E+01		8.2E+01			
			I	2.0E-02	I	1.0E+00	1.0E-01	1.4E+09		Acetochlor	34256-82-1					1.6E+03	6.6E+03		1.3E+03			
		9.0E-01	I	3.1E+01	A	1.0E+00	1.1E+05	1.4E+09	1.4E+04	Acetone	67-64-1					7.0E+04		4.4E+05	6.1E+04			
			I	2.0E-03	X	1.0E+00	1.1E+05	1.4E+09	2.4E+04	Acetone Cyanohydrin	75-86-5							5.0E+01	5.0E+01			
			I	6.0E-02	I	1.0E+00	1.3E+05	1.4E+09	1.3E+04	Acetonitrile	75-05-8							8.1E+02	8.1E+02			
3.8E+00	C	1.3E-03	C	1.0E-01	I	1.0E+00	2.5E+03	1.4E+09	6.0E+04	Acetophenone	98-86-2					7.8E+03			7.8E+03			
			I	1.0E-01	I	1.0E+00	1.0E-01	1.4E+09		Acetylaminofluorene, 2-	93-96-3	1.8E-01	6.5E-01	2.9E+03	1.4E-01							
5.0E-01	I	1.0E-04	I	5.0E-04	I	1.0E+00	2.3E+04	1.4E+09	6.9E+03	Acrolein	107-02-8	3.1E-01	1.2E+00	1.4E+04	2.4E-01	3.9E+01	6.6E+02	1.4E-01	1.4E-01			
			I	6.0E-03	I	1.0E+00	1.4E+09	1.4E+09		Acrylamide	79-06-1					1.6E+02	6.6E+02	8.5E+06	1.3E+02			
			I	1.0E-03	I	1.0E+00	1.1E+05	1.4E+09	9.5E+04	Acrylic Acid	79-10-7					3.9E+04	9.9E+01	9.9E+01	9.9E+01			
5.4E-01	I	6.8E-05	I	4.0E-02	A	1.0E+00	1.1E+04	1.4E+09	7.7E+03	Acrylonitrile	107-13-1	1.3E+00		3.2E-01	2.5E-01	3.1E+03		1.6E+01	1.6E+01			
			I	6.0E-03	P	1.0E+00	1.0E+00	1.4E+09		Adiponitrile	111-69-3							8.5E+06	8.5E+06			
5.6E-02	C	1.0E-02	I	1.0E-02	I	1.0E+00	1.0E-01	1.4E+09		Aldrichlor	15972-60-8	1.2E+01	4.4E+01		9.7E+00	7.8E+02	3.3E+03		6.3E+02			
			I	1.0E-03	I	1.0E+00	1.0E-01	1.4E+09		Aldicarb	116-06-3					7.8E+01	3.3E+02		6.3E+01			
			I	1.0E-03	I	1.0E+00	1.0E-01	1.4E+09		Aldicarb Sulfone	1646-88-4					7.8E+01	3.3E+02		6.3E+01			
			I	1.0E-03	I	1.0E+00	1.0E-01	1.4E+09		Aldicarb sulfide	1646-87-3											
1.7E+01	I	4.9E-03	I	3.0E-05	I	1.0E+00	1.4E+09	1.7E+06		Aldrin	309-00-2	4.1E-02		9.8E-01	3.9E-02	2.3E+00			2.3E+00			
			I	2.5E-01	I	1.0E+00	1.0E-01	1.4E+09		Allyl	74223-64-6					2.0E+04	8.2E+04		1.6E+04			
			I	5.0E-03	I	1.0E+00	1.1E+05	1.4E+09	3.4E+04	Allyl Alcohol	107-18-6					3.9E+02		3.6E+00	3.5E+00			
2.1E-02	C	6.0E-06	C	1.0E-03	I	1.0E+00	1.4E+03	1.4E+09	1.6E+03	Allyl Chloride	107-05-1	3.3E+01		7.4E-01	7.2E-01	7.8E+04		1.7E+00	1.7E+00			
			I	5.0E-03	P	1.0E+00	1.4E+09	1.4E+09		Aluminum	7429-90-5					7.8E+04			7.7E+04			
			I	4.0E-04	I	1.0E+00	1.4E+09	1.4E+09		Aluminum Phosphide	20859-73-8					3.1E+01		7.1E+06	3.1E+01			
			I	3.0E-04	I	1.0E+00	1.0E-01	1.4E+09		Amdriol	87486-29-4					2.3E+01	9.9E+01		1.9E+01			
2.1E+01	C	6.0E-03	C	9.0E-03	I	1.0E+00	1.0E-01	1.4E+09		Ametyln	834-12-8	3.3E-02	1.2E-01	6.4E+02	2.6E-02	7.0E+02	3.0E+03		5.7E+02			
			I	8.0E-02	P	1.0E+00	1.0E-01	1.4E+09		Amidobiphenyl, 4-	92-87-1											
			I	2.0E-02	P	1.0E+00	1.0E-01	1.4E+09		Aminophenol, m-	591-27-5					6.3E+03	2.6E+04		5.1E+03			
			I	2.5E-03	I	1.0E+00	1.0E-01	1.4E+09		Aminophenol, p-	123-30-8					1.6E+03	6.6E+03		1.3E+03			
			I	1.0E-01	I	1.0E+00	1.0E-01	1.4E+09		Amtraz	33089-61-1					2.0E+02	8.2E+02		1.6E+02			
			I	2.0E-01	I	1.0E+00	1.4E+09	1.4E+09		Ammonia	7664-41-7					1.6E+04		8.2E+01	1.6E+04			
			I	3.0E-03	X	1.0E+00	1.4E+09	2.6E+04		Ammonium Sulfamate	7773-06-0								8.2E+01			
			I	1.0E-03	I	1.0E+00	1.4E+09	1.4E+09		Amyl Alcohol, tert	75-85-4								1.6E+04			
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E+00	1.0E-01	1.4E+09		Aniline	62-53-3	1.2E+02	4.3E+02	2.4E+06	9.5E+01	5.5E+02	2.3E+03	1.4E+06	4.4E+02			
4.0E-02	P	2.0E-03	X	4.0E-04	I	1.0E+00	1.0E-01	1.4E+09		Anthraquinone, 9,10-	84-65-1	1.7E+01	6.2E+01		1.4E+01	1.6E+02	6.6E+02		1.3E+02			
			I	1.5E-01	I	1.0E+00	1.4E+09	1.4E+09		Antimony (metallic)	7440-36-0					3.1E+01			3.1E+01			
			I	5.0E-04	H	1.5E-01	1.4E+09	1.4E+09		Antimony Pentoxide	1314-60-9					3.9E+01			3.9E+01			
			I	9.0E-04	H	1.5E-01	1.4E+09	1.4E+09		Antimony Potassium Tartrate	11071-15-1					7.0E+01			7.0E+01			
			I	4.0E-04	H	1.5E-01	1.4E+09	1.4E+09		Antimony Tetroxide	1332-81-6					3.1E+01			3.1E+01			
			I	2.0E-04	I	1.5E-01	1.4E+09	1.4E+09		Antimony Trioxide	1309-64-4							2.8E+05	2.8E+05			
2.5E-02	I	7.1E-06	I	1.3E-02	I	1.0E+00	1.0E-01	1.4E+09		Apollo	74115-24-5	2.8E+01	9.9E+01	5.4E+05	2.2E+01	1.0E+03	4.3E+03		8.2E+02			
			I	5.0E-02	H	1.0E+00	1.0E-01	1.4E+09		Aramite	140-57-8					3.9E+03	1.6E+04		3.2E+03			
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.0E+00	3.0E-02	1.4E+09		Arsenic, Inorganic	7440-38-2	7.7E-01	5.5E+00	8.9E+02	6.8E-01	3.9E+01	3.3E+02	2.1E+04	3.5E+01			
			I	3.5E-06	C	1.0E+00	1.4E+09	1.4E+09		Arsine	7784-42-1					2.7E-01		7.1E+04	2.7E-01			
			I	9.0E-03	I	1.0E+00	1.0E-01	1.4E+09		Assure	76578-14-8					7.0E+02	3.0E+03		5.7E+02			
			I	5.0E-02	I	1.0E+00	1.0E-01	1.4E+09		Asulam	3337-71-1					3.9E+03	1.6E+04		3.2E+03			
2.3E-01	C	3.5E-02	I	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09		Atrazine	1912-24-9	3.0E+00	1.1E+01		2.4E+00	2.7E+03	1.2E+04		2.2E+03			
8.8E-01	C	2.5E-04	C	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09		Auramine	492-80-8	7.9E-01	2.8E+00	1.5E+04	6.2E-01							
1.1E-01	I	3.1E-05	I	4.0E-04	I	1.0E+00	1.0E-01	1.4E+09		Avermectin B1	65195-55-3	6.3E+00		4.7E+01	5.6E+00	3.1E+01	1.3E+02		2.5E+01			
			I	1.0E+00	P	7.0E-06	P	1.0E+00	5.2E+05	Azobenzene	103-33-3					7.8E+04	3.3E+05	9.9E+03	8.6E+03			
			I	5.0E-04	H	7.0E-02	1.4E+09	1.4E+09		Azodicarbonamide	123-77-3											
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	2.5E-02	1.4E+09	Barium	7440-39-3	3.1E-01		9.2E+00	3.0E-01	1.6E+04		7.1E+05	1.5E+04			
			I	4.0E-03	I	1.0E+00	1.0E-01	1.4E+09		Barium Chromate	10294-40-3					1.6E+03		2.8E+05	1.6E+03			
			I	3.0E-02	I	1.0E+00	1.0E-01	1.4E+09		Baygon	114-26-1					3.1E+02	1.3E+03		2.5E+02			
			I	2.5E-02	I	1.0E+00	1.0E-01	1.4E+09		Bayleton	43121-43-3					2.3E+03	9.9E+03		1.9E+03			
			I	3.0E-01	I	1.0E+00	1.4E+09	3.1E+05		Baythroid	68359-37-5					2.0E+03	8.2E+03		1.6E+03			
			I	1.0E+00	V	1.0E+00	1.4E+09	1.4E+09		Benefin	1861-40-1					2.3E+04			2.3E+04			
			I	5.0E-02	I	1.0E+00	1.0E-01	1.4E+09		Benomyl	17804-35-2					3.9E+03	1.6E+04		3.2E+03			
			I	3.0E-02	I	1.0E+00	1.0E-01	1.4E+09		Bentazon	25057-89-0					2.3E+03	9.9E+03		1.9E+03			
			I	1.0E-01	I	1.0E+00	1.2E+03	1.4E+09	2.3E+04	Benzaldehyde	100-52-7					7.8E+03			7.8E+03			
5.5E-02	I	7.8E-06	I	4.0E-03	I	1.0E+00	1.8E+03	1.4E+09	3.5E+03	Benzene	71-43-2	1.3E+01	2.5E+01</									

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _i (mg/m ³)	k _e y	v	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)	
1.3E+01	I			4.0E+00	I					1.0E+00	1.0E-01				Benzoic Acid	65-85-0					3.1E+05	1.3E+06		2.5E+05	
										1.0E+00	1.0E-01	3.2E+02		1.4E+09	Benzotrithloride	98-07-7	5.3E-02			5.3E-02					
1.7E-01	I	4.9E-05	C	1.0E-01	P	1.0E-03	P	V		1.0E+00	1.0E-01	1.5E+03		1.4E+09	Benzyl Alcohol	100-51-6	4.1E+00		1.5E+00	1.1E+00	7.8E+03	3.3E+04		6.3E+03	
		2.4E-03	I	2.0E-03	I	2.0E-05	I			1.0E+00	1.0E-01	7.0E-03		1.4E+09	Benzyl Chloride	100-44-7			1.6E+03	1.6E+03	1.6E+02		2.7E+01	2.3E+01	
				1.0E-04	I	1.0E-04	X	V		1.0E+00	1.0E-01			1.4E+09	Beryllium and compounds	7440-41-7			1.6E+03	1.6E+03	1.6E+02		2.8E+04	1.6E+02	
				9.0E-03	P					1.0E+00	1.0E-01			1.4E+09	Bidrin	141-66-2					7.8E+00	3.3E+01		6.3E+00	
				1.5E-02	I					1.0E+00	1.0E-01			1.4E+09	Bifenox	42576-02-3					7.0E+02	3.0E+03		5.7E+00	
8.0E-03	I	5.0E-01	I	4.0E-04	X	V				1.0E+00	1.0E-01			1.4E+09	Biphenthrin	82657-04-3					1.2E+03	4.9E+03		9.5E+02	
7.0E-02	H	1.0E-05	H	4.0E-02	I					1.0E+00	1.0E-01	1.0E+03		3.5E+04	Biphenyl, 1,1'-	92-52-4	8.7E+01		8.7E+01	8.7E+01	3.9E+04		4.8E+01	4.7E+01	
				3.0E-03	P					1.0E+00	1.0E-01			1.4E+09	Bis(2-chloro-1-methylethyl) ether	108-60-1	9.9E+00		9.8E+00	4.9E+00	3.1E+03			3.1E+03	
										1.0E+00	1.0E-01			1.4E+09	Bis(2-chloroethoxy)methane	111-91-1					2.3E+02	9.9E+02		1.9E+02	
1.1E+00	I	3.3E-04	I							1.0E+00	1.0E-01	5.1E+03		4.3E+04	Bis(2-chloroethyl)ether	111-44-4	6.3E-01		3.6E-01	2.3E-01					
2.2E+02	I	6.2E-02	I							1.0E+00	1.0E-01	4.2E+03		1.9E+03	Bis(chloromethyl)ether	542-88-1	3.2E-03		8.5E-05	8.3E-05					
				5.0E-02	I					1.0E+00	1.0E-01			1.4E+09	Bisphenol A	80-05-7					3.9E+03	1.6E+04		3.2E+03	
				2.0E-01	I	2.0E-02	H			1.0E+00	1.0E-01			1.4E+09	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		1.0E+00	1.0E-01			1.4E+09	Boron Trichloride	10294-34-5					1.6E+05		2.8E+07	1.6E+05	
				4.0E-02	C	1.3E-02	C	V		1.0E+00	1.0E-01			1.4E+09	Boron Trifluoride	7637-07-2					3.1E+03		1.8E+07	3.1E+03	
7.0E-01	I			4.0E-03	I					1.0E+00	1.0E-01			1.4E+09	Bromate	15541-46-4	9.9E-01		9.9E-01	9.9E-01	3.1E+02			3.1E+02	
2.0E+00	X	6.0E-04	X							1.0E+00	1.0E-01	2.4E+03		5.9E+03	Bromo-2-chloroethane, 1-	107-04-0	3.5E-01		2.8E-02	2.6E-02					
				8.0E-03	I	6.0E-02	I	V		1.0E+00	1.0E-01	6.8E+02		8.4E+03	Bromobenzene	108-86-1					6.3E+02		5.2E+02	2.9E+02	
				4.0E-02	X	V				1.0E+00	1.0E-01	4.0E+03		3.5E+03	Bromochloromethane	74-97-5					1.6E+03		1.5E+02	1.5E+02	
6.2E-02	I	3.7E-05	C	2.0E-02	I					1.0E+00	1.0E-01	9.3E+02		4.0E+03	Bromodichloromethane	75-27-4	1.1E+01		3.0E-01	2.9E-01	1.6E+03			1.6E+03	
7.9E-03	I	1.1E-06	I	2.0E-02	I					1.0E+00	1.0E-01	9.2E+02		9.7E+03	Bromotorm	75-25-2	8.8E+01		2.5E+01	1.9E+01	1.6E+03			1.6E+03	
				1.4E-03	I	5.0E-03	I	V		1.0E+00	1.0E-01	3.6E+03		1.4E+09	Bromomethane	74-83-9					1.1E+02		7.3E+00	6.8E+00	
				5.0E-03	H					1.0E+00	1.0E-01			1.2E+05	Bromophos	2104-96-3					3.9E+02			3.9E+02	
				2.0E-02	I					1.0E+00	1.0E-01			1.4E+09	Bromoxynil	1589-84-5					1.6E+03	6.6E+03		1.3E+03	
				2.0E-02	I					1.0E+00	1.0E-01			4.7E+05	Bromoxynil Octanoate	1889-99-2					1.6E+03			1.6E+03	
3.4E+00	C	3.0E-05	I	2.0E-03	I	V				1.0E+00	1.0E-01	6.7E+02		8.7E+02	Butadiene, 1,3-	106-99-0	2.0E-01		8.1E-02	5.8E-02	1.6E+03		1.8E+00	1.8E+00	
				1.0E-01	I					1.0E+00	1.0E-01	7.6E+03		3.0E+04	Butanol, n-	11-36-3					7.8E+03			7.8E+03	
1.9E-03	P			2.0E-01	I					1.0E+00	1.0E-01			1.4E+09	Butyl Benzyl Phthalate	85-68-7	3.7E+02	1.3E+03		2.9E+02	1.6E+04	6.6E+04		1.3E+04	
				2.0E+00	P	3.0E+01	P	V		1.0E+00	1.0E-01	2.1E+04		2.9E+04	Butyl alcohol, sec-	78-92-2					1.6E+06		9.1E+05	1.3E+06	
				5.0E-02	I					1.0E+00	1.0E-01			8.6E+04	Butylate	2008-41-5					3.9E+03			3.9E+03	
2.0E-04	C	5.7E-08	C							1.0E+00	1.0E-01			1.4E+09	Butylated hydroxyanisole	25013-16-5	3.5E+03	1.2E+04	6.7E+07	2.7E+03	3.9E+03			3.9E+03	
3.6E-03	P			3.0E-01	P					1.0E+00	1.0E-01			1.4E+09	Butylated hydroxytoluene	128-37-0	1.9E+02	6.9E+02		1.5E+02	2.3E+04	9.9E+04		1.9E+04	
				5.0E-02	P					1.0E+00	1.0E-01	1.1E+02		8.1E+03	Butylbenzene, n-	104-51-8					3.9E+03			3.9E+03	
				1.0E-01	X					1.0E+00	1.0E-01	1.5E+02		7.4E+03	Butylbenzene, sec-	135-98-8					7.8E+03			7.8E+03	
				1.0E-01	X					1.0E+00	1.0E-01	1.8E+02		7.4E+03	Butylbenzene, tert-	98-06-6					7.8E+03			7.8E+03	
				2.0E-02	A					1.0E+00	1.0E-01			1.4E+09	Cacodylic Acid	75-60-5					1.6E+03	6.6E+03		1.3E+03	
				1.8E-03	I	1.0E-03	I	1.0E-05	A	2.5E-02	1.0E-03			1.4E+09	Cadmium (Diet)	7440-43-9			2.1E+03	2.1E+03	7.8E+01		8.2E+02	1.4E+04	7.1E+01
				1.8E-03	I	5.0E-04	I	1.0E-05	A	5.0E-02	1.0E-03			1.4E+09	Cadmium (Water)	7440-43-9					7.8E+01				
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M		2.5E-02				1.4E+09	Calcium Chromate	13765-19-0	3.1E-01		9.2E+00	3.0E-01	1.6E+03		2.8E+05	1.6E+03	
				5.0E-01	I	2.2E-03	C			1.0E+00	1.0E-01			1.4E+09	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					1.0E+00	1.0E-01			1.4E+09	Captafol	2425-06-1	4.6E+00	1.6E+01	8.9E+04	3.6E+00	1.6E+02	6.6E+02		1.3E+02	
2.3E-03	C	6.6E-07	C	1.3E-01	I					1.0E+00	1.0E-01			1.4E+09	Captan	133-06-2	3.0E+02	1.1E+03	5.8E+06	2.4E+02	1.0E+04	4.3E+04		8.2E+03	
				1.0E-01	I					1.0E+00	1.0E-01			1.4E+09	Carbaryl	63-25-2					7.8E+03	3.3E+04		6.3E+03	
				5.0E-03	I					1.0E+00	1.0E-01			1.4E+09	Carbofuran	1563-66-2					3.9E+02	1.6E+03		3.2E+02	
				1.0E-01	I	7.0E-01	I	V		1.0E+00	1.0E-01	7.4E+02		1.2E+03	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		1.0E+00	1.0E-01	4.6E+02		1.5E+03	Carbon Tetrachloride	56-23-5	9.9E+00		7.0E-01	6.5E-01	3.1E+02		1.6E+02	1.0E+02	
				1.0E-02	I					1.0E+00	1.0E-01			1.4E+09	Carbosulfan	55285-14-8					7.8E+02	3.3E+03		6.3E+02	
				1.0E-01	I					1.0E+00	1.0E-01			1.4E+09	Carboxin	5234-68-4					7.8E+03	3.3E+04		6.3E+03	
				9.0E-04	I					1.0E+00	1.0E-01			1.4E+09	Ceric oxide	1306-38-3							1.3E+06	1.3E+06	
				1.0E-01	I					1.0E+00	1.0E-01			1.5E+05	Chloral Hydrate	302-17-0					7.8E+03			7.8E+03	
				1.5E-02	I					1.0E+00															

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Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -y) ⁻¹	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _i (mg/m ³)	k _e (y)	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)	
2.4E-01	I	6.9E-05	C	1.0E-02	I	7.5E-03	I		1.0E+00	1.0E-01		1.4E+09		Cypermethrin	52315-07-8					7.8E+02	3.3E+03		6.3E+02	
														Cyromazine	66215-27-8					5.9E+02	2.5E+03		4.7E+02	
3.4E-01	I	9.7E-05	C	1.0E-02	I	1.0E-02	I	V	1.0E+00	1.0E-01		1.4E+09	2.1E+06	DDE, p,p'	72-55-9	2.9E+00	1.0E+01	5.5E+04	2.3E+00					
3.4E-01	I	9.7E-05	I	5.0E-04	I	1.0E-02	I		1.0E+00	3.0E-02		1.4E+09		DDT	50-29-3	2.0E+00	2.4E+01	3.9E+04	1.9E+00	3.9E+01	5.5E+02		3.7E+01	
												1.4E+09		Dacthal	1861-32-1					7.8E+02	2.5E+03		6.3E+02	
7.0E-04	I			3.0E-02	I	4.0E-05	I		1.0E+00	1.0E-01		1.4E+09		Dalapon	75-99-0					2.3E+03	9.9E+03		1.9E+03	
												1.4E+09		Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5	9.9E+02	3.5E+03		7.8E+02	5.5E+02	2.3E+03		4.4E+02	
												1.4E+09		Demeton	8065-48-3					3.1E+00	1.3E+01		2.5E+00	
1.2E-03	I			6.0E-01	I				1.0E+00	1.0E-01		1.4E+09		Di(2-ethylhexyl)adipate	103-23-1	5.8E+02	2.1E+03		4.5E+02	4.7E+04	2.0E+05		3.8E+04	
6.1E-02	H			7.0E-04	A				1.0E+00	1.0E-01		1.4E+09		Diallate	2303-16-4	1.1E+01	4.1E+01		8.9E+00	5.5E+01	2.3E+02		4.4E+01	
												1.4E+09		Diazinon	333-41-5					5.5E+01	2.3E+02		4.4E+01	
8.0E-01	P	6.0E-03	P	1.0E-02	X	2.0E-04	I	V	1.0E+00		9.8E+02	1.4E+09	3.2E+04	Dibenzothiophene	132-65-0	1.9E-01		5.4E-03	5.3E-03	7.8E+02			7.8E+02	
											1.6E+02	1.4E+09	1.9E+04	Dibromo-3-chloropropane, 1,2-	96-12-8					1.6E+01		6.7E+00	4.7E+00	
												1.4E+09		Dibromobenzene, 1,3-	108-36-1					3.1E+01			3.1E+01	
8.4E-02	I	2.7E-05	C	1.0E-02	I				1.0E+00		8.0E+02	1.4E+09	8.0E+03	Dibromobenzene, 1,4-	106-37-6					7.8E+02			7.8E+02	
2.0E+00	I	6.0E-04	I	9.0E-03	I	9.0E-03	I	V	1.0E+00		1.3E+03	1.4E+09	8.6E+03	Dibromochloromethane	124-48-1	8.3E+00		8.3E-01	7.5E-01	1.6E+03			1.6E+03	
												1.4E+09		Dibromoethane, 1,2-	106-93-4	3.5E-01		4.0E-02	3.6E-02	7.0E+02		8.1E+01	7.3E+01	
												1.4E+09		Dibromomethane (Methylene Bromide)	74-95-3					7.8E+02		2.4E+01	2.3E+01	
												1.4E+09		Dibutyltin Compounds	NA					2.3E+01	9.9E+01		1.9E+01	
												1.4E+09		Dicamba	1918-00-9					2.3E+03	9.9E+03		1.9E+03	
4.2E-03	P								1.0E+00		5.2E+02	1.4E+09	1.2E+04	Dichloro-2-butene, 1,4-	764-41-0			8.3E-03	8.3E-03					
4.2E-03	P								1.0E+00		5.2E+02	1.4E+09	1.1E+04	Dichloro-2-butene, cis-1,4-	1476-11-5			7.4E-03	7.4E-03					
4.2E-03	P								1.0E+00		7.6E+02	1.4E+09	1.1E+04	Dichloro-2-butene, trans-1,4-	110-57-6			7.4E-03	7.4E-03					
5.0E-02	I	4.0E-03	I	9.0E-02	I	2.0E-01	H	V	1.0E+00	1.0E-01		1.4E+09		Dichloroacetic Acid	79-43-6	1.4E+01	4.9E+01		1.1E+01	3.1E+02	1.3E+03		2.5E+02	
5.4E-03	C	1.1E-05	C	7.0E-02	A	8.0E-01	I	V	1.0E+00		3.8E+02	1.4E+09	1.2E+04	Dichlorobenzene, 1,2-	95-50-1					7.0E+03		2.4E+03	1.8E+03	
												1.4E+09		Dichlorobenzene, 1,4-	106-36-7	1.3E+02		2.7E+00	2.6E+00	5.5E+03		8.7E+03	3.4E+03	
4.5E-01	I	3.4E-04	C	1.0E+00	I	1.0E-01	I		1.0E+00	1.0E-01		1.4E+09		Dichlorobenzidine, 3,3'	91-94-1	1.5E+00	5.5E+00	1.1E+04	1.2E+00					
											8.5E+02	1.4E+09	8.4E+02	Dichlorobenzophenone, 4,4'	90-98-2					7.0E+02	3.0E+03		5.7E+02	
5.7E-03	C	1.6E-06	C	2.0E-01	P				1.0E+00		1.7E+03	1.4E+09	2.1E+03	Dichlorodifluoromethane	75-71-8					1.6E+04		8.8E+01	8.7E+01	
9.1E-02	I	2.6E-05	I	6.0E-03	X	7.0E-03	P	V	1.0E+00		3.0E+03	1.4E+09	4.6E+03	Dichloroethane, 1,1-	75-34-3	1.2E+02		3.7E+00	3.6E+00	1.6E+04			1.6E+04	
											1.2E+03	1.4E+09	1.2E+03	Dichloroethane, 1,2-	107-06-2	7.6E+00		4.9E-01	4.6E-01	4.7E+02		3.3E+01	3.1E+01	
												1.4E+09		Dichloroethylene, 1,1-	75-35-4					3.9E+03		2.4E+02	2.3E+02	
											2.4E+03	1.4E+09	2.5E+03	Dichloroethylene, 1,2-cis-	156-59-2					1.6E+02			1.6E+02	
											1.9E+03	1.4E+09	1.7E+03	Dichloroethylene, 1,2-trans-	156-60-5					1.6E+03			1.6E+03	
											1.0E+00	1.0E-01		Dichlorophenol, 2,4-	120-83-2					2.3E+02	9.9E+02		1.9E+02	
											1.0E+00	5.0E-02		Dichlorophenoxy Acetic Acid, 2,4-	94-75-7					7.8E+02	6.6E+03		7.0E+02	
											1.0E+00	1.0E-01		Dichlorophenoxy butyric Acid, 4-(2,4-	94-82-6					6.3E+02	2.6E+03		5.1E+02	
3.6E-02	C	1.0E-05	C	9.0E-02	A	4.0E-03	I	V	1.0E+00		1.4E+03	1.4E+09	3.8E+03	Dichloropropane, 1,2-	78-87-5	1.9E+01		1.1E+00	1.0E+00	7.0E+03		1.6E+01	1.6E+01	
											1.5E+03	1.4E+09	6.8E+03	Dichloropropane, 1,3-	142-28-9					1.6E+03			1.6E+03	
1.0E-01	I	4.0E-06	I	3.0E-02	I	2.0E-02	I	V	1.0E+00	1.0E-01		1.4E+09		Dichloropropanol, 2,3-	616-23-9	7.0E+00		2.5E+00	1.8E+00	2.3E+02	9.9E+02		1.9E+02	
											1.6E+03	1.4E+09	3.6E+03	Dichloropropene, 1,3-	542-75-6					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		1.0E+00	1.0E-01		1.4E+09		Dichlorvos	62-73-7	2.4E+00	8.5E+00	4.6E+04	1.9E+00	3.9E+01	1.6E+02		7.1E+05	3.2E+01
1.6E+01	I	4.6E-03	I	8.0E-02	P	3.0E-04	X	V	1.0E+00		1.4E+09	4.1E+03		Dicyclopentadiene	77-73-6					6.3E+03		1.3E+00	1.3E+00	
											1.0E+00	1.0E-01		Weldm	80-57-1	4.3E-02	1.5E-01	8.3E+02	3.4E-02	3.9E+00	1.6E+01		3.2E+00	
												1.0E+00	1.0E-01	Diesel Engine Exhaust	NA					1.6E+02	6.6E+02	2.8E+05	1.3E+02	
												1.4E+09		Diethanolamine	111-42-2					2.3E+03	9.9E+03	1.4E+05	1.9E+03	
												1.4E+09		Diethylene Glycol Monoethyl Ether	112-34-5									
												1.4E+09		Diethylene Glycol Monoethyl Ether	111-90-0					4.7E+03	2.0E+04	4.3E+05	3.8E+03	
3.5E+02	C	1.0E-01	C	1.0E-03	P				1.0E+00		1.1E+05	1.4E+09	1.4E+05	Diethylformamide	617-84-5	2.0E-03	7.1E-03	3.8E+01	1.6E-03	7.8E+01			7.8E+01	
											1.0E+00	1.0E-01		Diethylstilbestrol	56-53-1									
												1.4E+09		Difenzoquat	43222-48-6					6.3E+03	2.6E+04		5.1E+03	
												1.4E+09		Diflubenzuron	35367-38-5					1.6E+03	6.6E+03		1.3E+03	
											1.4E+03	1.4E+09	1.2E+03	Difluoroethane, 1,1-	75-37-6							4.8E+04	4.8E+04	
4.4E-02	C	1.3E-05	C						1.0E+00		1.4E+09	1.5E+03		Dihydrosofrole	94-58-6	1.6E+01		3.3E-01	3.2E-01					
											2.3E+03	1.4E+09	3.1E+03	Diisopropyl Ether	108-20-3							2.2E+03	2.2E+03	
											5.3E+02	1.4E+09	3.8E+04	Diisopropyl Methylphosphonate	1445-75-6					6.3E+03			6.3E+03	
1.6E+00	P		</																					

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1				
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)
1.1E+01	P			2.0E-03	I		V		1.0E+00	1.0E+00	8.3E+02	1.4E+09	3.1E+04	Dimethylaniline, N,N-Dimethylbenzidine, 3,3'-	121-69-7 119-93-7	6.3E-02	2.2E-01		4.9E-02	1.6E+02			1.6E+02
				1.0E-01	P	3.0E-02	I	V	1.0E+00		1.1E+05	1.4E+09	1.3E+05	Dimethylformamide	68-12-2					7.8E+03		4.0E+03	2.6E+03
5.5E+02	C	1.6E-01	C	1.0E-04	X	2.0E-06	X	V	1.0E+00		1.7E+05	1.4E+09	1.6E+05	Dimethylhydrazine, 1,1-Dimethylhydrazine, 1,2-	57-14-7 540-73-8	1.3E-03		2.9E-03	8.8E-04	7.8E+00	7.8E+00	3.4E-01	3.2E-01
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Dimethylphenol, 2,4-Dimethylphenol, 2,6-Dimethylphenol, 3,4-	105-67-9 576-26-1 95-65-8					1.6E+03	6.6E+03		1.3E+03
4.5E-02	C	1.3E-05	C	1.0E-03	I		V		1.0E+00	1.0E-01	1.1E+03	1.4E+09	1.0E+03	Dimethylvinylchloride	513-37-1	1.5E+01		2.2E-01	2.1E-01	4.7E+01	2.0E+02	3.3E+02	3.8E+01
				8.0E-05	X				1.0E+00	1.0E-01		1.4E+09		Dinitro-o-cresol, 4,6-Dinitro-o-cyclohexyl Phenol, 4,6-	534-52-1 131-89-5					6.3E+00	2.6E+01		5.1E+00
				2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Dinitrobenzene, 1,2-Dinitrobenzene, 1,3-Dinitrobenzene, 1,4-	528-29-0 99-65-0 100-25-4					7.8E+00	3.3E+01		6.3E+00
				1.0E-04	P				1.0E+00	1.0E-01		1.4E+09		Dinitrophenol, 2,4-Dinitrotoluene Mixture, 2,4/2,6-Dinitrotoluene, 2,4-	51-28-5 NA 121-14-2	1.0E+00	3.6E+00		8.0E-01	1.6E+02	6.6E+02		1.3E+02
6.8E-01	I			3.1E-01	C	8.9E-05	C		1.0E+00	1.0E-01		1.4E+09		Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 4-Amino-2,6-	606-20-2 35572-78-2 19406-51-0	4.6E-01	1.7E+00		3.6E-01	2.3E+01	1.0E+02		1.9E+01
1.5E+00	P			2.0E-03	X				1.0E+00	9.9E-02		1.4E+09		Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 4-Amino-2,6-	606-20-2 35572-78-2 19406-51-0	4.6E-01	1.7E+00		3.6E-01	1.6E+02	6.5E+02		1.3E+02
4.5E-01	X			9.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Dinitrotoluene, Technical grade	25321-14-6	1.5E+00	5.5E+00		1.2E+00	7.0E+01	3.3E+02		5.7E+01
1.0E-01	I	5.0E-06	I	3.0E-02	I	3.0E-02	I	V	1.0E+00	1.0E-01	1.2E+05	1.4E+09	4.0E+04	Dioxane, 1,4-Dioxins	88-85-7 123-91-1	7.0E+00		2.2E+01	5.3E+00	7.8E+01	3.3E+02	1.2E+03	6.3E+01
6.2E+03	I	1.3E+00	I	1.3E+05	C	3.8E+01	C		1.0E+00	3.0E-02		1.4E+09		-Hexachlorodibenzo-p-dioxin, Mixture	NA	1.1E-04	1.3E-03	2.9E+00	1.0E-04	5.5E-05	7.7E-04	8.2E-02	5.1E-05
				3.0E-02	I				1.0E+00	1.0E-01		1.4E+09		-TCDD, 2,3,7,8-	1746-01-6	5.3E-06	6.3E-05	1.4E-04	4.8E-06	2.3E+03	9.9E+03		1.9E+03
				8.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Diphenyl Sulfone	957-51-7					6.3E+01	2.6E+02		5.1E+01
				2.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Diphenylamine	122-39-4					2.0E+03	8.2E+03		1.6E+03
8.0E-01	I	2.2E-04	I	2.2E-03	I				1.0E+00	1.0E-01		1.4E+09		Diphenylhydrazine, 1,2-Diquat	122-66-7 85-00-7	8.7E-01	3.1E+00	1.7E+04	6.8E-01	1.7E+02	7.3E+02		1.4E+02
7.1E+00	C	1.4E-01	C						1.0E+00	1.0E-01		1.4E+09		Direct Black 38	1937-37-7	9.8E-02	3.5E-01	2.7E+01	7.6E-02				
7.4E+00	C	1.4E-01	C						1.0E+00	1.0E-01		1.4E+09		Direct Blue 6	2602-46-2	9.4E-02	3.3E-01	2.7E+01	7.3E-02				
6.7E+00	C	1.4E-01	C	4.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Direct Brown 95	16071-86-6	1.0E-01	3.7E-01	2.7E+01	8.1E-02	3.1E+00	1.3E+01		2.5E+00
				1.0E-02	I		V		1.0E+00		1.4E+09	4.5E+04		Dithane, 1,4-Diuron	505-29-3 330-54-1					7.8E+02			7.8E+02
				4.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Doding	2439-10-3					1.6E+02	6.6E+02		1.3E+02
				2.5E-02	I		V		1.0E+00		1.4E+09	1.2E+05		EPTC	759-94-4					2.0E+03			2.0E+03
				6.0E-03	I		V		1.0E+00		1.4E+09	4.1E+05		Endosulfan	115-29-7					4.7E+02			4.7E+02
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Endothal	145-73-3					1.6E+03	6.6E+03		1.3E+03
9.9E-03	I	1.2E-06	I	3.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Endrin	72-20-8	7.0E+01		4.4E+01	2.7E+01	2.3E+01	9.9E+01		1.9E+01
				6.0E-03	P	1.0E-03	I	V	1.0E+00		1.1E+04	1.4E+09	1.9E+04	Epichlorohydrin	106-89-8					4.7E+02		2.0E+01	1.9E+01
				2.0E-02	I				1.0E+00		1.5E+04	1.4E+09	7.7E+03	Epoxybutane, 1,2-Ethephon	16672-87-0					1.6E+02		1.6E+02	1.6E+02
				5.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Ethion	563-12-2					3.9E+02	1.6E+03		3.2E+02
				1.0E-01	P	6.0E-02	P	V	1.0E+00		3.1E+04	1.4E+09	6.2E+04	Ethoxyethanol Acetate, 2-Ethoxyethanol, 2-	111-15-9					7.8E+03	1.6E+02	3.8E+03	2.6E+03
4.8E-02	H			9.0E-02	P	2.0E-01	I	V	1.0E+00		1.1E+05	1.4E+09	9.8E+04	Ethyl Acetate	110-80-5	1.4E+01			1.4E+01	7.0E+03		2.1E+04	5.2E+03
				9.0E-01	I	7.0E-02	P	V	1.0E+00		1.1E+04	1.4E+09	8.6E+03	Ethyl Acrylate	141-78-6					7.0E+04		6.3E+02	6.2E+02
				5.0E-03	P	8.0E-03	P	V	1.0E+00		2.5E+03	1.4E+09	6.3E+03	Ethyl Chloride (Chloroethane)	140-88-5					3.9E+02		5.3E+01	4.7E+01
				1.0E+01	I		V		1.0E+00		2.1E+03	1.4E+09	1.3E+03	Ethyl Ether	75-00-3					7.8E+01		1.4E+04	1.4E+04
				2.0E-01	I		V		1.0E+00		1.0E+04	1.4E+09	3.1E+03	Ethyl Methacrylate	60-29-7					1.6E+04			1.6E+04
				9.0E-02	H	3.0E-01	P	V	1.0E+00		1.1E+03	1.4E+09	5.8E+03	Ethyl-p-nitrophenyl Phosphonate	97-63-2					7.0E+03		1.8E+03	1.4E+03
1.1E-02	C	2.5E-06	C	1.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Ethylbenzene	2104-64-5	6.3E+01		6.4E+00	5.8E+00	7.8E+01	3.3E+00		6.3E-01
				7.0E-02	P				1.0E+00	1.0E-01	4.8E+02	1.4E+09	5.7E+03	Ethylene Cyanohydrin	100-41-4					7.8E+03		5.9E+03	3.4E+03
				9.0E-02	P		V		1.0E+00		1.9E+05	1.4E+09	1.8E+05	Ethylene Diamine	109-78-4					5.5E+03	2.3E+04		4.4E+03
				2.0E+00	I	4.0E-01	C		1.0E+00	1.0E-01		1.4E+09		Ethylene Glycol	107-15-3					7.0E+03			7.0E+03
				1.0E-01	I	1.6E+00	I		1.0E+00	1.0E-01		1.4E+09		Ethylene Glycol Monobutyl Ether	107-21-1					1.6E+05	6.6E+05	5.7E+08	1.3E+05
3.1E-01	C	8.8E-05	C	3.0E-02	C	V			1.0E+00		1.2E+05	1.4E+09	6.1E+03	Ethylene Oxide	111-76-2	2.2E+00		1.9E-01	1.8E-01	7.8E+03	3.3E+04	2.3E+09	6.3E+03
4.5E-02	C	1.3E-05	C	8.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Ethylene Thiourea	75-21-8					1.6E+05	6.6E+05		1.9E+02
6.5E+01	C	1.9E-02	C				V		1.0E+00		1.5E+05	1.4E+09	2.4E+04	Ethyleneimine	96-45-7	1.5E+01	5.5E+01	2.9E+05	1.2E+01	6.3E+00	2.6E+01		5.1E+00
				3.0E+00	I				1.0E+00	1.0E-01		1.4E+09		Ethylphthalyl Ethyl Glycolate	84-72-0					2.3E+05	9.9E+05		1.9E+05
				8.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Express	101200-48-0					6.3E+02	2.6E+03		5.1E+02

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1				
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³ -y) ⁻¹	k e y	RfD _o (mg/kg-day)	k e y	RfC _i (mg/m ³)	k e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)
2.5E-04	I			1.0E+00	1.0E-01									Penamiphos	22224-92-6					2.0E+01	8.2E+01		1.8E+01
2.5E-02	I			1.0E+00	1.0E-01									Fenprothrin	39515-41-8					2.0E+03	8.2E+03		1.6E+03
1.3E-02	I			1.0E+00	1.0E-01									Fluometuron	2164-17-2					1.0E+03	4.3E+03		8.2E+02
4.0E-02	C	1.3E-02	C	1.0E+00										Fluoride	16984-48-8					3.1E+03		1.8E+07	3.1E+03
6.0E-02	I	1.3E-02	C	1.0E+00										Fluorine (Soluble Fluoride)	7782-41-4					4.7E+03		1.8E+07	4.7E+03
8.0E-02	I			1.0E+00	1.0E-01									Fluridone	59756-60-4					6.3E+03	2.6E+04		5.1E+03
2.0E-02	I			1.0E+00	1.0E-01									Flurprimidol	56425-91-3					1.6E+03	6.6E+03		1.3E+03
6.0E-02	I			1.0E+00	1.0E-01									Flutolanil	66332-96-5					4.7E+03	2.0E+04		3.8E+03
1.0E-02	I			1.0E+00	1.0E-01									Fluvalinate	69409-94-5					7.8E+02	3.3E+03		6.3E+02
3.5E-03	I			1.0E-01	1.0E-01									Folpet	133-07-3	2.0E+02	7.1E+02		1.6E+02	7.8E+03	3.3E+04		6.3E+03
1.9E-01	I			1.0E+00	1.0E-01									Fomesafen	72178-02-0	3.7E+00	1.3E+01		2.9E+00				
		2.0E-03	I	1.0E+00	1.0E-01									Fonofos	944-22-9					1.6E+02	6.6E+02		1.3E+02
		1.3E-05	I	2.0E-01	9.8E-03	A	V		1.0E+00	4.2E+04	1.4E+09	7.8E+04		Formaldehyde	50-00-0			1.7E+01	1.7E+01	1.6E+04	8.0E+02		7.6E+02
9.0E-01	P	3.0E-04	X	V	1.0E+00	1.1E+05	1.4E+09	9.3E+04	1.0E+00	1.0E-01	1.4E+09			Formic Acid	64-18-6					7.0E+04	2.9E+01		2.9E+01
3.0E+00	I			1.0E+00	1.0E-01									Fosetyl-AL	39148-24-8					2.3E+05	9.9E+05		1.9E+05
														Furans									
1.0E-03	X		V	1.0E+00	3.0E-02				1.4E+09	2.0E+05				Dibenzofuran	132-64-9					7.8E+01	1.1E+03		7.3E+01
1.0E-03	I		V	1.0E+00	3.0E-02	6.2E+03	1.4E+09	2.6E+03	1.0E+00	3.0E-02	6.2E+03	1.4E+09		-Furan	110-00-9					7.8E+01	1.1E+03		7.3E+01
9.0E-01	I	2.0E+00	I	V	1.0E+00	3.0E-02	1.7E+05	1.4E+09	1.2E+04					-Tetrahydrofuran	109-99-9					7.0E+04	9.9E+05	2.5E+04	1.8E+04
3.8E+00	H			1.0E+00	1.0E-01									Furazolidone	67-45-8	1.8E-01	6.5E-01		1.4E-01				
1.5E+00	C	4.3E-04	C	3.0E-03	I	5.0E-02	H	V	1.0E+00	1.0E-01	1.0E+04	4.9E+04		Furfural	98-01-1					2.3E+02		2.5E+03	2.1E+02
3.0E-02	I	8.6E-06	C	1.0E+00	1.0E-01									Furium	531-82-8	4.6E-01	1.6E+00	8.9E+03	3.6E-01				
				1.0E+00	1.0E-01									Furmecycloz	60588-05-0	2.3E+01	8.2E+01	4.4E+05	1.8E+01				
		4.0E-04	I	1.0E+00	1.0E-01									Glufosinate, Ammonium	77182-82-2					3.1E+01	1.3E+02		2.5E+01
				8.0E-05	C				1.0E+00	1.0E-01	1.4E+09			Glutaraldehyde	111-30-8							1.1E+05	1.1E+05
4.0E-04	I	1.0E-03	H	V	1.0E+00	1.1E+05	1.4E+09	7.3E+04	1.0E+00	1.0E-01	1.4E+09			Glycidyl	785-34-4					3.1E+01		7.7E+01	2.2E+01
1.0E-01	I			1.0E+00	1.0E-01									Glyphosate, Goal	1071-83-6					7.8E+03	3.3E+04		6.3E+03
3.0E-03	I			1.0E+00	1.0E-01									Guanidine	42874-03-3					2.3E+02	9.9E+02		1.9E+02
1.0E-02	X		V	1.0E+00					1.4E+09	1.5E+05				Guanidine Chloride	133-00-8					7.8E+02			7.8E+02
2.0E-02	P			1.0E+00	1.0E-01				1.4E+09					Guthrie	50-01-1					1.6E+03	6.6E+03		1.3E+03
3.0E-03	A	1.0E-02	A	1.0E+00	1.0E-01				1.4E+09					Harmony	98-50-0					2.3E+02	9.9E+02	1.4E+07	1.9E+02
5.0E-05	I			1.0E+00	1.0E-01				1.4E+09					Haloxyp, Methyl	69806-40-2					3.9E+00	1.6E+01		3.2E+00
1.3E-02	I			1.0E+00	1.0E-01				1.4E+09					Heptachlor	79277-27-3					1.0E+03	4.3E+03		8.2E+02
4.5E+00	I	1.3E-03	I	5.0E-04	I		V		1.0E+00	4.8E+05				Heptachlor Epoxide	76-44-8	1.5E-01	1.0E+00	1.3E-01		3.9E+01			3.9E+01
9.1E+00	I	2.6E-03	I	1.3E-05	I	V			1.0E+00	8.4E+05				Hexachlor Epoxide	1024-57-3	7.6E-02	9.1E-01	7.0E-02		1.0E+00			1.0E+00
1.6E+00	I	4.6E-04	I	8.0E-04	I	V			1.0E+00	3.8E+05				Hexachlorobenzene	87-82-1					1.6E+02			1.6E+02
7.8E-02	I	2.2E-05	I	1.0E-03	P		V		1.0E+00	1.7E+01	1.4E+09	1.1E+04		Hexachlorodiphenyl ether, 2,2',4,4',5,5'-(BDE-163)	69831-49-2	4.3E-01	4.1E-01	2.1E-01		1.6E+01	6.6E+01		1.3E+01
6.3E+00	I	1.8E-03	I	8.0E-03	A				1.0E+00	1.0E-01	1.4E+09			Hexachlorobenzene	118-74-1	1.1E-01	3.9E-01	2.1E+03	8.6E-02	6.3E+01			6.3E+01
1.8E+00	I	5.3E-04	I	1.0E+00	1.0E-01				1.4E+09					Hexachlorobutadiene	87-68-3	8.9E+00	1.4E+00	1.2E+00		7.8E+01			7.8E+01
1.1E+00	C	3.1E-04	C	3.0E-04	I				1.4E+09					Hexachlorocyclohexane, Alpha-	319-84-6	1.1E-01	3.9E-01	2.1E+03	8.6E-02	6.3E+02	2.6E+03		5.1E+02
1.8E+00	I	5.1E-04	I	1.0E+00	1.0E-01				1.4E+09					Hexachlorocyclohexane, Beta-	319-85-7	3.9E-01	1.4E+00	7.2E+03	3.0E-01				
1.1E+00	C	3.1E-04	C	3.0E-04	I				1.0E+00	4.0E-02	1.4E+09			Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	6.3E-01	5.6E+00	1.2E+04	5.7E-01	2.3E+01	2.5E+02		2.1E+01
1.8E+00	I	5.1E-04	I	1.0E+00	1.0E-01				1.4E+09					Hexachlorocyclohexane, Technical	608-73-1	3.9E-01	1.4E+00	7.5E+03	3.0E-01				
4.0E-02	I	1.1E-05	C	6.0E-03	I	2.0E-04	I	V	1.0E+00	1.6E+01	1.4E+09	8.5E+03		Hexachlorocyclopentadiene	77-47-4					4.7E+02		1.8E+00	1.8E+00
7.0E-04	I	3.0E-02	I	V	1.0E+00				1.4E+09	8.0E+03				Hexachloroethane	67-72-1	1.7E+01	2.0E+00	1.8E+00		5.5E+01		2.5E+02	4.5E+01
3.0E-04	I			1.0E+00	1.0E-01				1.4E+09					Hexachlorophene	70-30-4					2.3E+01	9.9E+01		1.9E+01
1.1E-01	I			3.0E-03	I	1.0E+00	1.5E-02		1.4E+09					Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	6.3E+00	1.5E+02		6.1E+00	2.3E+02	6.6E+03		2.3E+02
				1.0E+00					5.2E+03	1.4E+09	3.0E+05			Hexamethylene Diisocyanate, 1,6-Hexamethylphosphoramide	822-06-0					3.1E+01	1.3E+02		3.1E+00
4.0E-04	P			1.0E+00	1.0E-01				1.4E+09					Hexane, N-	680-31-9								2.5E+01
6.0E-02	H	7.0E-01	I	V	1.0E+00	1.4E+02	1.4E+09	8.3E+02	1.0E+00	1.0E-01	1.4E+09			Hexanedioic Acid	110-54-3					4.7E+03		6.1E+02	5.4E+02
2.0E+00	P			1.0E+00	1.0E-01				1.4E+09					Hexanone, 2-	124-04-9					1.6E+05	6.6E+05		1.3E+05
5.0E-03	I	3.0E-02	I	V	1.0E+00	3.3E+03	1.4E+09	1.3E+04	1.0E+00	1.0E-01	1.4E+09			Hexazinone	591-78-6					3.9E+02	4.2E+02		2.0E+02
3.0E+00	I	4.9E-03	I	3.3E-02	I	1.0E+00	1.0E-01		1.4E+09					Hydrazine	51235-04-2	2.3E-01			2.3E-01	2.6E+03	1.1E+04		2.1E+03
3.0E+00	I	4.9E-03	I	1.0E+00					1.4E+09					Hydrazine Sulfate	302-01-2	2.3E-01	7.8E+02	2.3E-01				4.3E+04	4.3E+04
				2.0E-02	I	V			1.0E+00					Hydrogen Chloride	7647-01-0							2.8	

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1				
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg-day)	k e y	RfC _i (mg/m ³)	k e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)
9.5E-04	I			3.0E-01	I		V		1.0E+00		1.0E+04	1.4E+09	2.8E+04	Isobutyl Alcohol	78-83-1				5.7E+02	2.3E+04			2.3E+04
				2.0E-01	I	2.0E+00	C		1.0E+00	1.0E-01		1.4E+09		Isophorone	78-59-1	7.3E+02	2.6E+03			1.6E+04	6.6E+04	2.8E+09	1.3E+04
				1.5E-02	I		V		1.0E+00		1.4E+09	4.2E+05		Isopropalin	33820-53-0					1.2E+03			1.2E+03
				2.0E+00	P	2.0E-01	P V		1.0E+00		1.1E+05	1.4E+09	2.8E+04	Isopropanol	67-63-0					1.6E+05		5.8E+03	5.6E+03
				1.0E-01	I				1.0E+00	1.0E-01		1.4E+09		Isopropyl Methyl Phosphonic Acid	1832-54-8					7.8E+03	3.3E+04		6.3E+03
				5.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Isoxaben	82558-50-7					3.9E+03	1.6E+04		3.2E+03
						3.0E-01	A V		1.0E+00		1.4E+09			JP-7	NA							4.3E+08	4.3E+08
				7.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Kerb	23950-58-5					5.9E+03	2.5E+04		4.7E+03
				2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Lactofen	77501-63-4					1.6E+02	6.6E+02		1.3E+02
														Lead Compounds									
														--Lead Chromate	7758-97-6	3.1E-01		9.2E+00	3.0E-01	1.6E+03		2.8E+05	1.6E+03
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M	2.5E-02		1.4E+09			--Lead Phosphate	7446-27-7	8.2E+01		3.2E+05	8.2E+01				
8.5E-03	C	1.2E-05	C						1.0E+00		1.4E+09			--Lead acetate	301-04-2	2.5E+00	8.8E+00	4.8E+04	1.9E+00				
									1.0E+00	1.0E-01		1.4E+09		--Lead and Compounds	7439-92-1								4.0E+02
									1.0E+00	1.0E-01		1.4E+09		--Lead subacetate	1335-32-6	8.2E+01	2.9E+02	3.2E+05	6.4E+01				
				1.0E-07	I		V		1.0E+00		2.4E+00	1.4E+09	1.9E+03	--Tetraethyl Lead	78-00-2					7.8E-03			7.8E-03
				2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Linuron	330-55-2					1.6E+02	6.6E+02		1.3E+02
				2.0E-03	P				1.0E+00		1.4E+09			Lithium	7439-93-2					1.6E+02			1.6E+02
				2.0E-01	I				1.0E+00	1.0E-01		1.4E+09		Londax	83055-99-6					1.6E+04	6.6E+04		1.3E+04
				5.0E-04	I				1.0E+00	1.0E-01		1.4E+09		MCPA	94-74-6					3.9E+01	1.6E+02		3.2E+01
				1.0E-02	I				1.0E+00	1.0E-01		1.4E+09		MCPB	94-81-5					7.8E+02	3.3E+03		6.3E+02
				1.0E-03	I				1.0E+00	1.0E-01		1.4E+09		MCPD	93-65-2					7.8E+01	3.3E+02		6.3E+01
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Malathion	121-75-5					1.6E+03	6.6E+03		1.3E+03
				1.0E-01	I	7.0E-04	C		1.0E+00	1.0E-01		1.4E+09		Maleic Anhydride	108-31-6					7.8E+03	3.3E+04	9.9E+05	6.3E+03
				5.0E-01	I				1.0E+00	1.0E-01		1.4E+09		Maleic Hydrzade	123-33-1					3.9E+04	1.6E+05		3.2E+04
				1.0E-04	P				1.0E+00	1.0E-01		1.4E+09		Malononitrile	109-77-3					7.8E+00	3.3E+01		6.3E+00
				3.0E-02	H				1.0E+00	1.0E-01		1.4E+09		Mancozeb	8018-01-7					2.3E+03	9.9E+03		1.9E+03
				5.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Manganese (Diet)	12427-38-2					3.9E+02	1.6E+03		3.2E+02
				1.4E-01	I	5.0E-05	I		1.0E+00			1.4E+09		Manganese (Non-diet)	7439-96-5					1.9E+03		7.1E+04	1.8E+03
				2.4E-02	S	5.0E-05	I		4.0E-02			1.4E+09		Manganese (Non-diet)	7439-96-5								
				9.0E-05	H				1.0E+00	1.0E-01		1.4E+09		Mephosalan	850-107-7					7.0E+00	3.0E+01		5.7E+00
				3.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Mepiquat Chloride	24307-26-4					2.3E+03	9.9E+03		1.9E+03
														Mercury Compounds									
				3.0E-04	I	3.0E-04	S		7.0E-02		1.4E+09			--Mercuric Chloride (and other Mercury salts)	7487-94-7					2.3E+01		4.3E+05	2.3E+01
						3.0E-04	I V		1.0E+00	3.1E+00	1.4E+09	3.0E+04		--Mercury (elemental)	7439-97-6							9.4E+00	9.4E+00
				1.0E-04	I				1.0E+00		1.4E+09			--Methyl Mercury	22967-92-6					7.8E+00			7.8E+00
				8.0E-05	I				1.0E+00	1.0E-01		1.4E+09		--Phenylmercuric Acetate	62-38-4					6.3E+00	2.6E+01		5.1E+00
				3.0E-05	I		V		1.0E+00		1.4E+09	1.9E+06		Merphos	150-50-5					2.3E+00			2.3E+00
				3.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Merphos Oxide	88-48-8					2.3E+00	9.9E+00		1.9E+00
				6.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Metakyl	57837-19-1					4.7E+03	2.0E+04		3.8E+03
				1.0E-04	I	3.0E-02	P V		1.0E+00	4.6E+03	1.4E+09	6.8E+03		Methacrylonitrile	126-98-7					7.8E+00		2.1E+02	7.5E+00
				5.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Methamidophos	10265-92-6					3.9E+00	1.6E+01		3.2E+00
				2.0E+00	I	2.0E+01	I V		1.0E+00		1.1E+05	1.4E+09	2.9E+04	Methanol	67-56-1					1.6E+05		6.1E+05	1.2E+05
				1.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Methidathion	950-37-8					7.8E+01	3.3E+02		6.3E+01
				2.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Methomyl	16152-77-5					2.0E+03	8.2E+03		1.6E+03
4.9E-02	C	1.4E-05	C						1.0E+00	1.0E-01		1.4E+09		Methoxy-5-nitroaniline, 2-	99-50-2	1.4E+01	5.0E+01	2.7E+05	1.1E+01	3.9E+02	1.6E+03		3.2E+02
				5.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Methoxychlor	72-43-5					6.3E+02		1.3E+02	1.1E+02
				8.0E-03	P	1.0E-03	P V		1.0E+00		1.2E+05	1.4E+09	1.2E+05	Methoxyethanol Acetate, 2-	110-49-6								
				5.0E-03	P	2.0E-02	I V		1.0E+00		1.1E+05	1.4E+09	1.0E+05	Methoxyethanol, 2-	109-86-4					3.9E+02		2.1E+03	3.3E+02
				1.0E+00	X		V		1.0E+00		2.9E+04	1.4E+09	8.1E+03	Methyl Acetate	79-20-9					7.8E+04			7.8E+04
				3.0E-02	H	2.0E-02	I V		1.0E+00		6.8E+03	1.4E+09	7.0E+03	Methyl Acrylate	96-33-3					2.3E+03		1.5E+02	1.4E+02
				6.0E-01	I	5.0E+00	I V		1.0E+00		2.8E+04	1.4E+09	1.2E+04	Methyl Ethyl Ketone (2-Butanone)	78-93-3					4.7E+04		6.4E+04	2.7E+04
				1.0E-03	X				1.0E+00		1.8E+05	1.4E+09	1.6E+05	Methyl Hydrazine	60-34-4			4.4E-01	4.4E-01	7.8E+01		3.3E+00	3.1E+00
				8.0E-02	H	3.0E+00	I V		1.0E+00		3.4E+03	1.4E+09	1.1E+04	Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1					6.3E+03		3.3E+04	5.3E+03
						1.0E-03	C V		1.0E+00		1.7E+04	1.4E+09	4.4E+03	Methyl Isocyanate	624-83-9								4.6E+00
				1.4E+00	I	7.0E-01	I V		1.0E+00		2.4E+03	1.4E+09	6.3E+03	Methyl Methacrylate	80-62-6					1.1E+05		4.6E+03	4.4E+03
				2.5E-04	I				1.0E+00	1.0E-01		1.4E+09		Methyl Parathion	298-00-0					2.0E+01	8.2E+01		1.6E+01
				6.0E-02	X				1.0E+00	1.0E-01		1.4E+09		Methyl Phosphonic Acid	993-13-5					4.7E+03	2.0E+04		

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg-day)	k e y	RfC _i (mg/m ³)	k e y	v o l a t i l e	m u t a g e n	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)	
1.3E-01	C	3.7E-05	C	1.0E-02	A					1.0E+00	1.0E-01				Methylaniline Hydrochloride, 2- Methylarsonic acid	636-21-5 124-58-3	5.3E+00	1.9E+01	1.0E+05	4.2E+00	7.8E+02	3.3E+03		6.3E+02	
1.0E-01	X	3.0E-04	X	2.0E-04	X					1.0E+00	1.0E-01				Methylbenzene, 1,4-diamine monohydrochloride, 2- Methylbenzene-1,4-diamine sulfate, 2- Methylcholanthrene, 3-	74612-12-7 615-50-9 56-49-5	7.0E+00	2.5E+01		5.4E+00	1.6E+01	6.6E+01		1.3E+01	
2.2E+01	C	6.3E-03	C	3.0E-04	X				M	1.0E+00	1.0E-01						7.0E-03	2.7E-02	2.2E+02	5.5E-03	2.3E+01	9.9E+01		1.9E+01	
2.0E-03	I	1.0E-08	I	6.0E-03	I	6.0E-01	I	V	M	1.0E+00		3.3E+03	1.4E+09	2.2E+03	Methylene Chloride Methylene-bis(2-chloroaniline), 4,4'- Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	75-09-2 101-14-4 101-61-1	7.7E+01	6.0E+00	2.2E+02	5.7E+01	4.7E+02	6.6E+02	1.4E+03	3.5E+02	
1.0E-01	P	4.3E-04	C	2.0E-03	P				M	1.0E+00	1.0E-01						1.5E+00	3.2E+03	3.2E+03	1.2E+00	1.6E+02	6.6E+02		1.3E+02	
4.6E-02	I	1.3E-05	C			2.0E-02	C			1.0E+00	1.0E-01				Methylenediphenyl Diisocyanate Methylstyrene, Alpha-	101-77-9 98-83-9	1.5E+01	5.4E+01	2.9E+05	1.2E+01			2.8E+07 8.5E+05	2.8E+07 8.5E+05	
1.6E+00	C	4.6E-04	C	7.0E-02	H				V	1.0E+00		5.0E+02	1.4E+09	1.3E+04					8.3E+03	3.4E-01	5.5E+03			5.5E+03	
				1.5E-01	I					1.0E+00	1.0E-01				Metolachlor Metribuzin Mineral oils	51218-45-2 21087-64-9 8012-95-1					1.2E+04 2.0E+03 2.3E+05	4.9E+04 8.2E+03		9.5E+03 1.6E+03 2.3E+05	
1.8E+01	C	5.1E-03	C	2.0E-04	I				V	1.0E+00			1.4E+09	8.6E+05	Mirex Molinate Molybdenum	2385-85-5 2212-67-1 7439-98-7	3.9E-02		4.7E-01	3.6E-02	1.6E+01	6.6E+02		1.6E+01 1.3E+02 3.9E+02	
				1.0E-01	I					1.0E+00			1.4E+09		Monochloramine Monomethylaniline N,N-Diphenyl-1,4-benzenediamine	10599-90-3 100-61-8 74-31-7					7.8E+03	6.6E+02	9.9E+01	7.8E+03 1.3E+02 1.9E+01	
1.8E+00	C	0.0E+00	C	2.0E-03	I				V	1.0E+00			1.4E+09	5.7E+04	Naled Naphtha, High Flash Aromatic (HFAN) Naphthylamine, 2-	300-76-5 64742-95-6 91-59-8	3.9E-01	1.4E+00		3.0E-01	1.6E+02 2.3E+03		1.4E+08	1.6E+02 2.3E+03	
				1.0E-01	I					1.0E+00	1.0E-01		1.4E+09		Napropamide Nickel Acetate Nickel Carbonate	15299-99-7 373-02-4 3333-67-3			1.5E+04 1.5E+04	1.5E+04 1.5E+04	7.8E+03 8.6E+02	3.3E+04 3.6E+03	2.0E+04 2.0E+04	6.3E+03 6.7E+02 6.7E+02	
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1.0E+00			1.4E+09		Nickel Carbonyl Nickel Hydroxide Nickel Oxide	13463-39-3 12054-48-7 1313-99-1			1.5E+04 1.5E+04 1.5E+04	1.5E+04 1.5E+04 1.5E+04	8.6E+02 8.6E+02 8.6E+02	2.0E+04 2.0E+04 2.8E+04		8.2E+02 8.2E+02 8.4E+02	
				2.6E-04	C	1.1E-02	C	1.4E-05	C	4.0E-02			1.4E+09		Nickel Refinery Dust Nickel Soluble Salts Nickel Sulfide	NA 7440-02-0 12035-72-2	4.1E-01		1.6E+04 1.5E+04 8.0E+03	1.6E+04 1.5E+04 4.1E-01	8.6E+02 1.6E+03 8.6E+02	2.0E+04 1.3E+05 2.0E+04		8.2E+02 1.5E+03 8.2E+02	
1.7E+00	C	4.8E-04	I	1.1E-02	C	1.4E-05	C	4.0E-02		1.0E+00	1.0E-01		1.4E+09		Nickelocene Nitrate Nitrate + Nitrite (as N)	1271-28-9 14797-55-8 NA		1.5E+04	1.5E+04	8.6E+02	3.6E+03	2.0E+04	6.7E+02	1.3E+05	
2.0E-02	P			1.0E-01	I					1.0E+00			1.4E+09		Nitrite Nitroamine, 2- Nitroamine, 4-	14797-65-0 88-74-4 100-01-6	3.5E+01	1.2E+02		2.7E+01	7.8E+03 7.8E+02 3.1E+02	3.3E+03 3.3E+03	7.1E+04 8.5E+06	7.8E+03 6.3E+02 2.5E+02	
				4.0E-05	I	2.0E-03	I	9.0E-03	I	1.0E+00		3.1E+03	1.4E+09	7.3E+04	Nitrobenzene Nitrocellulose Nitroturanton	98-95-3 9004-70-0 67-20-9			5.1E+00	5.1E+00	1.6E+02	9.9E+08	6.9E+02	1.3E+02 1.9E+08 4.4E+03	
1.3E+00	C	3.7E-04	C	1.0E-04	P					1.0E+00	1.0E-01		1.4E+09		Nitrofurazone Nitroglycerin Nitroguanidine	59-87-0 55-63-0 556-88-7	5.3E-01	1.9E+00	1.0E+04	4.2E-01	7.8E+00 7.8E+03	3.3E+01 3.3E+04		6.3E+00 6.3E+03	
1.7E-02	P			1.0E-01	I					1.0E+00	1.0E-01		1.4E+09		Nitromethane Nitropropane, 2- Nitroso-N-ethylurea, N-	75-52-5 79-46-9 759-73-9			5.4E+00 1.4E-02 1.8E+02	5.4E+00 1.4E-02 4.5E-03	8.8E+01 2.7E+02	8.8E+01 2.7E+02	8.8E+01 2.7E+02		
2.7E+01	C	7.7E-03	C	5.0E-03	P	V			M	1.0E+00		1.8E+04	1.4E+09	1.7E+04											
1.2E+02	C	3.4E-02	C	2.0E-02	I	V				1.0E+00	1.0E-01		1.4E+09		Nitroso-N-methylurea, N- Nitroso-di-N-butylamine, N- Nitroso-di-N-propylamine, N-	684-93-5 924-16-3 621-64-7	1.3E-03 1.3E-01 9.9E-02	5.0E-03 4.3E-01 3.5E-01	4.1E+01 9.9E-02 7.8E-02	1.0E-03 9.9E-02 7.8E-02					
5.4E+00	I	1.6E-03	I						V	1.0E+00			1.4E+09	2.4E+05	Nitrosodiethanolamine, N- Nitrosodiethylamine, N-	1116-54-7 55-18-5	2.5E-01 1.0E-03	8.8E-01 4.0E-03	4.8E+03 3.2E+01	1.9E-01 8.1E-04					
7.0E+00	I	2.0E-03	C	1.0E+00	I				M	1.0E+00	1.0E-01		1.4E+09		Nitrosodimethylamine, N-	62-75-9	3.0E-03		6.0E-03	2.0E-03	6.3E-01		3.4E+00	5.3E-01	
2.8E+00	I	8.0E-04	C						M	1.0E+00	1.0E-01		1.4E+09		Nitrosodiphenylamine, N- Nitrosomethylethylamine, N- Nitrosomorpholine [N-]	86-30-6 10595-95-6 59-89-2	1.4E+02 3.2E-02 1.0E-01	5.0E+02 5.4E-02 3.7E-01	1.5E+06 5.4E-02 2.0E+03	1.1E+02 2.0E-02 8.1E-02					
1.5E+02	I	4.3E-02	I	1.0E+00	I				M	1.0E+00	1.0E-01		1.4E+09		Nitrosopiperidine [N-] Nitrosopyrrolidine, N- Nitrotoluene, m-	100-75-4 930-55-2 99-08-1	7.4E-02 3.3E-01	2.6E-01 1.2E+00	1.4E+03 6.3E+03	5.8E-02 2.6E-01					
5.1E+01	I	1.4E-02	I	8.0E-06	P	4.0E-05	X	V	M	1.0E+00		2.4E+05	1.4E+09	8.2E+04											
4.9E-03	I	2.6E-06	C						V	1.0E+00	1.0E-01		1.4E+09		Nitrosodiphenylamine, N- Nitrosomethylethylamine, N- Nitrosomorpholine [N-]	86-30-6 10595-95-6 59-89-2	1.4E+02 3.2E-02 1.0E-01	5.0E+02 5.4E-02 3.7E-01	1.5E+06 5.4E-02 2.0E+03	1.1E+02 2.0E-02 8.1E-02					
2.2E+01	P	9.0E-04	P	3.0E-04	X	2.0E-02	P	V		1.0E+00		1.5E+03	1.4E+09	1.4E+05	Nitrotoluene, o- Nitrotoluene, p- Nonane, n-	88-72-2 99-99-0 111-84-2	3.2E+00 4.3E+01	1.5E+02		3.2E+00 3.4E+01	7.0E+01 3.1E+02 2.3E+01	1.3E+03	2.2E+01	7.0E+01 2.5E+02 1.1E+01	
6.7E+00	C	1.9E-03	C	4.0E-02	I					1.0E+00	1.0E-01		1.4E+09		Norflurazon Nustar	27314-13-2 85509-19-9					3.1E+03 5.5E+01	1.3E+04 2.3E+02		2.5E+03 4.4E+01	

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1				
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³ -y) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)
				3.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Octabromodiphenyl Ether	32536-52-0					2.3E+02	9.9E+02		1.9E+02
				5.0E-02	I				1.0E+00	6.0E-03		1.4E+09		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				1.0E+00	1.0E-01		1.4E+09		Octamethylpyrophosphoramide	152-16-9					1.6E+02	6.6E+02		1.3E+02
				5.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Oryzalin	19044-88-3					3.9E+03	1.6E+04		3.2E+03
				5.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Oxadiazon	19866-30-9					3.9E+02	1.6E+03		3.2E+02
				2.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Oxamyl	23135-22-0					2.0E+03	8.2E+03		1.6E+03
				1.3E-02	I				1.0E+00	1.0E-01		1.4E+09		Paclobutrazol	76738-62-0					1.0E+03	4.3E+03		8.2E+02
				4.5E-03	I				1.0E+00	1.0E-01		1.4E+09		Paraquat Dichloride	1910-42-5					3.5E+02	1.5E+03		2.8E+02
				6.0E-03	H				1.0E+00	1.0E-01		1.4E+09		Parathion	56-38-2					4.7E+02	2.0E+03		3.8E+02
				5.0E-02	H			V	1.0E+00			4.5E+04		Pebutal	1114-71-2					3.9E+03			3.9E+03
				4.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Pendimethalin	40487-42-1					3.1E+03	1.3E+04		2.5E+03
				2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Pentabromodiphenyl Ether	32534-81-9					1.6E+02	6.6E+02		1.3E+02
				1.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Pentabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-99)	60348-60-9					7.8E+00	3.3E+01		6.3E+00
				8.0E-04	I			V	1.0E+00		4.5E+02	1.4E+09	8.1E+04	Pentachlorobenzene	608-93-5					6.3E+01			6.3E+01
9.0E-02	P							V	1.0E+00		4.5E+02	1.4E+09	9.7E+03	Pentachloroethane	76-01-7	7.7E+00			7.7E+00				
2.6E-01	H			3.0E-03	I			V	1.0E+00		4.3E+05	1.4E+09	4.3E+05	Pentachloronitrobenzene	82-68-8	2.7E+00			2.7E+00				2.3E+02
4.0E-01	I	5.1E-06	C	5.0E-03	I				1.0E+00	2.6E-01		1.4E+09		Pentachlorophenol	87-96-5	1.7E+00	2.5E+00	7.5E+05	1.0E+00	3.9E+02	6.6E+02		2.5E+02
4.0E-03	X			2.0E-03	P				1.0E+00	1.0E-01		1.4E+09		Pentaerythritol tetranitrate (PETN)	78-11-5	1.7E+02	6.2E+02		1.4E+02	1.6E+02	6.6E+02		1.3E+02
						1.0E+00	P	V	1.0E+00		3.9E+02	1.4E+09	7.8E+02	Pentane, n-	109-66-0							8.1E+02	8.1E+02
														Perchlorates									
				7.0E-04	I				1.0E+00			1.4E+09		--Ammonium Perchlorate	7790-98-9					5.5E+01			5.5E+01
				7.0E-04	I				1.0E+00			1.4E+09		--Lithium Perchlorate	7791-03-9					5.5E+01			5.5E+01
				7.0E-04	I				1.0E+00			1.4E+09		--Perchlorate and Perchlorate Salts	14797-73-0					5.5E+01			5.5E+01
				7.0E-04	I				1.0E+00			1.4E+09		--Potassium Perchlorate	11118-14-1					5.5E+01			5.5E+01
				7.0E-04	I				1.0E+00			1.4E+09		--Sodium Perchlorate	1801-89-0					5.5E+01			5.5E+01
				2.0E-02	P			V	1.0E+00			1.4E+09	1.3E+05	Perfluorobutane Sulfonate	375-73-5					1.6E+03			1.6E+03
2.2E-03	C	6.3E-07	C	5.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Permethrin	52645-53-1					3.9E+03	1.6E+04		3.2E+03
									1.0E+00	1.0E-01		1.4E+09		Phenacetin	62-44-2	3.2E+02	1.1E+03	6.1E+06	2.5E+02				
				2.5E-01	I				1.0E+00	1.0E-01		1.4E+09		Phenmedipham	13684-63-4					2.0E+04	8.2E+04		1.6E+04
				3.0E-01	I	2.0E-01	C		1.0E+00	1.0E-01		1.4E+09		Phenol	108-95-2					2.3E+04	9.9E+04	2.8E+08	1.9E+04
				5.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Phentolazone	92-84-2					3.9E+01	1.6E+02		3.2E+01
				6.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Phenylenediamine, m-	108-45-2					4.7E+02	2.0E+03		3.8E+02
4.7E-02	H								1.0E+00	1.0E-01		1.4E+09		Phenylenediamine, o-	95-54-5	1.5E+01	5.3E+01		1.2E+01				
				1.9E-01	H				1.0E+00	1.0E-01		1.4E+09		Phenylenediamine, p-	106-50-3					1.5E+04	6.3E+04		1.2E+04
1.9E-03	H								1.0E+00	1.0E-01		1.4E+09		Phenylphenol, 2-	90-43-7	3.6E+02	1.3E+03		2.8E+02				
				2.0E-04	H				1.0E+00	1.0E-01		1.4E+09		Phosphate	298-02-2					1.6E+01	6.6E+01		1.3E+01
						3.0E-04	I	V	1.0E+00		1.6E+03	1.4E+09	9.8E+02	Phosgene	75-44-5							3.1E-01	3.1E-01
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Phosmet	782-11-6					1.6E+03	6.6E+03		1.3E+03
														Phosphates, Inorganic									
				4.9E+01	P				1.0E+00			1.4E+09		--Aluminum metaphosphate	13176-88-0					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Ammonium polyphosphate	68333-79-9					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Calcium pyrophosphate	7790-76-3					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Diammonium phosphate	7783-28-0					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Dicalcium phosphate	7757-93-9					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Dimagnesium phosphate	7782-75-4					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Dipotassium phosphate	1158-11-4					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Disodium phosphate	7558-79-4					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Monoaluminum phosphate	13530-50-2					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Monoammonium phosphate	7722-76-1					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Monocalcium phosphate	7758-23-8					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Monomagnesium phosphate	7757-86-0					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Monopotassium phosphate	7778-77-0					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Monosodium phosphate	7558-80-7					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Polyphosphoric acid	8017-16-1					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Potassium triphosphate	13845-36-8					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium acid pyrophosphate	7758-16-9					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium aluminum phosphate (acidic)	7785-88-8					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium aluminum phosphate (anhydrous)	10279-59-1					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium aluminum phosphate (tetrahydrate)	10305-76-7					3.8E+06			3.8E+06
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium hexametaph									

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Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³ -y) ⁻¹	k e y	RfD _o (mg/kg-day)	k e y	RfC _i (mg/m ³)	k e y	v o l a t i l e	m u t a g e n	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)	
				4.9E+01	P					1.0E+00			1.4E+09		~Tetrasodium pyrophosphate	7722-88-5					3.8E+06			3.8E+06	
				4.9E+01	P					1.0E+00			1.4E+09		~Trialuminum sodium tetra decahydrooctaorthophosphate (dihydrate)	15136-87-5					3.8E+06			3.8E+06	
				4.9E+01	P					1.0E+00			1.4E+09		~Tricalcium phosphate	7758-87-4					3.8E+06			3.8E+06	
				4.9E+01	P					1.0E+00			1.4E+09		~Trimagnesium phosphate	7757-87-1					3.8E+06			3.8E+06	
				4.9E+01	P					1.0E+00			1.4E+09		~Tripotassium phosphate	7778-53-2					3.8E+06			3.8E+06	
				4.9E+01	P					1.0E+00			1.4E+09		~Trisodium phosphate	7601-54-9					3.8E+06			3.8E+06	
				3.0E-04	I	3.0E-04	I	V		1.0E+00			1.4E+09		Phosphine	7803-51-2					2.3E+01		4.3E+05	2.3E+01	
				4.9E+01	P	1.0E-02	I			1.0E+00			1.4E+09		Phosphoric Acid	7664-38-2					3.8E+06		1.4E+07	3.0E+06	
				2.0E-05	I			V		1.0E+00			1.4E+09	6.9E+03	Phosphorus, White	7723-14-0					1.6E+00			1.6E+00	
															Phthalates										
1.4E-02	I	2.4E-06	C	2.0E-02	I					1.0E+00	1.0E-01		1.4E+09		~Bis(2-ethylhexyl)phthalate	117-81-7	5.0E+01	1.8E+02	1.6E+06	3.9E+01	1.6E+03	6.6E+03		1.3E+03	
				1.0E+00	I					1.0E+00	1.0E-01		1.4E+09		~Butylphthalyl Butylglycolate	85-70-1					7.8E+04	3.3E+05		6.3E+04	
				1.0E-01	I					1.0E+00	1.0E-01		1.4E+09		~Dibutyl Phthalate	84-74-2					7.8E+03	3.3E+04		6.3E+03	
				8.0E-01	I					1.0E+00	1.0E-01		1.4E+09		~Diethyl Phthalate	84-66-2					6.3E+04	2.6E+05		5.1E+04	
				1.0E-01	I			V		1.0E+00			1.4E+09	2.1E+04	~Dimethylterephthalate	120-61-6					7.8E+03			7.8E+03	
				1.0E-02	P					1.0E+00	1.0E-01		1.4E+09		~Octyl Phthalate, di-N-	117-84-0					7.8E+02	3.3E+03		6.3E+02	
				1.0E+00	H					1.0E+00	1.0E-01		1.4E+09		~Phthalic Acid, P-	100-21-0					7.8E+04	3.3E+05		6.3E+04	
				2.0E+00	I	2.0E-02	C			1.0E+00	1.0E-01		1.4E+09		~Phthalic Anhydride	85-44-9					1.6E+05	6.6E+05	2.8E+07	1.3E+05	
				7.0E-02	I					1.0E+00	1.0E-01		1.4E+09		Picloram	1918-02-1					5.5E+03	2.3E+04		4.4E+03	
				1.0E-04	X					1.0E+00	1.0E-01		1.4E+09		Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3					7.8E+00	3.3E+01		6.3E+00	
				1.0E-02	I					1.0E+00	1.0E-01		1.4E+09		Pirimiphos, Methyl	29232-93-7					7.8E+02	3.3E+03		6.3E+02	
3.0E+01	C	8.6E-03	C	7.0E-06	H					1.0E+00	1.0E-01		1.4E+09		Polybrominated Biphenyls	59536-65-1	2.3E-02	8.2E-02	4.4E+02	1.8E-02	5.5E-01	2.3E+00		4.4E-01	
				7.0E-02	S	2.0E-05	S	7.0E-05	I	V		1.4E+09	5.9E+05		Polychlorinated Biphenyls (PCBs)										
				2.0E+00	S	5.7E-04	S			1.0E+00	1.4E-01		1.4E+09	1.1E+05	~Aroclor 1221	11104-28-2	3.5E-01	8.8E-01	5.6E-01	1.7E-01	1.8E+00	5.5E+00	2.8E+03	1.4E+00	
				2.0E+00	S	5.7E-04	S			1.0E+00	1.4E-01		1.4E+09	1.1E+05	~Aroclor 1232	11141-16-5	3.5E-01	8.8E-01	5.5E-01	1.7E-01	1.8E+00	5.5E+00	2.0E+03	1.4E+00	
				2.0E+00	S	5.7E-04	S			1.0E+00	1.4E-01		1.4E+09	7.9E+05	~Aroclor 1242	53469-21-9	3.5E-01	8.8E-01	3.9E+00	2.3E-01	1.8E+00	5.5E+00	2.1E+03	1.4E+00	
				2.0E+00	S	5.7E-04	S			1.0E+00	1.4E-01		1.4E+09	5.1E+05	~Aroclor 1248	12872-29-6	3.5E-01	8.8E-01	2.5E+00	2.3E-01	1.8E+00	5.5E+00	2.1E+03	1.4E+00	
				2.0E+00	S	5.7E-04	S	2.0E-05	I	V		1.0E+00	1.4E+01	1.4E+09	8.4E+05	~Aroclor 1254	11097-69-1	3.5E-01	8.8E-01	4.1E+00	2.4E-01	1.6E+00	4.7E+00		1.2E+00
				2.0E+00	S	5.7E-04	S			1.0E+00	1.4E-01		1.4E+09	1.3E+06	~Aroclor 1260	13966-82-5	3.5E-01	8.8E-01	6.5E+00	2.4E-01	1.8E+00	5.5E+00	2.1E+03	1.4E+00	
				6.0E-04	X			V		1.0E+00	1.4E-01		1.4E+09	7.2E+05	~Aroclor 5460	11126-42-4	1.8E-01	4.5E-01	5.0E+00	1.2E-01	1.8E+00	5.5E+00	2.8E+03	1.4E+00	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.0E+00	1.4E-01		1.4E+09	2.0E+06	~Heptachlorobiphenyl, 2,3,3',4,4',5,5'-(PCB 189)	39636-31-9	1.8E-01	4.5E-01	3.5E+00	1.2E-01	1.8E+00	5.5E+00	2.0E+03	1.4E+00	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.0E+00	1.4E-01		1.4E+09	1.4E+06	~Hexachlorobiphenyl, 2,3',4,4',5,5'-(HCB 167)	52663-72-6	1.8E-01	4.5E-01	3.5E+00	1.2E-01	1.8E+00	5.5E+00	2.0E+03	1.4E+00	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.0E+00	1.4E-01		1.4E+09	1.5E+06	~Hexachlorobiphenyl, 2,3,3',4,4',5'-(PCB 157)	69782-90-7	1.8E-01	4.5E-01	3.6E+00	1.2E-01	1.8E+00	5.5E+00	2.0E+03	1.4E+00	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.0E+00	1.4E-01		1.4E+09	1.5E+06	~Hexachlorobiphenyl, 2,3,3',4,4',5'-(PCB 156)	38380-08-4	1.8E-01	4.5E-01	3.8E+00	1.2E-01	1.8E+00	5.5E+00	2.1E+03	1.4E+00	
3.9E+03	E	1.1E+00	E	2.3E-08	E	1.3E-06	E	V		1.0E+00	1.4E-01		1.4E+09	1.4E+06	~Hexachlorobiphenyl, 3,3',4,4',5,5'-(PCB 169)	32774-16-6	1.8E-04	4.5E-04	3.5E-03	1.2E-04	1.8E-03	5.5E-03	2.0E+00	1.4E-03	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.0E+00	1.4E-01		1.4E+09	1.0E+06	~Pentachlorobiphenyl, 2,3,4,4',5'-(PCB 123)	66510-44-3	1.8E-01	4.5E-01	2.5E+00	1.2E-01	1.8E+00	5.5E+00	1.4E+03	1.4E+00	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.0E+00	1.4E-01		1.4E+09	8.3E+05	~Pentachlorobiphenyl, 2,3',4,4',5'-(PCB 118)	31508-00-6	1.8E-01	4.5E-01	2.0E+00	1.2E-01	1.8E+00	5.5E+00	1.2E+03	1.4E+00	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.0E+00	1.4E-01		1.4E+09	8.5E+05	~Pentachlorobiphenyl, 2,3,3',4,4'-(HCB 165)	32598-14-4	1.8E-01	4.5E-01	2.1E+00	1.2E-01	1.8E+00	5.5E+00	1.2E+03	1.4E+00	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.0E+00	1.4E-01		1.4E+09	1.0E+06	~Pentachlorobiphenyl, 2,3,4,4',5'-(PCB 114)	74472-37-0	1.8E-01	4.5E-01	2.5E+00	1.2E-01	1.8E+00	5.5E+00	1.4E+03	1.4E+00	
1.3E+04	E	3.8E+00	E	7.0E-09	E	4.0E-07	E	V		1.0E+00	1.4E-01		1.4E+09	1.0E+06	~Pentachlorobiphenyl, 3,3',4,4',5'-(PCB 126)	57465-28-8	5.3E-05	1.4E-04	7.5E-04	3.7E-05	5.5E-04	1.6E-03	4.3E-01	4.1E-04	
2.0E+00	I	5.7E-04	I					V		1.0E+00	1.4E-01		1.4E+09	7.9E+05	~Polychlorinated Biphenyls (high risk)	1336-36-3	3.5E-01	8.8E-01	3.9E+00	2.3E-01	1.8E+00	5.5E+00	2.1E+03	1.4E+00	
4.0E-01	I	1.0E-04	I					V		1.0E+00	1.4E-01		1.4E+09		~Polychlorinated Biphenyls (low risk)	1336-36-3									
7.0E-02	I	2.0E-05	I					V		1.0E+00	1.4E-01		1.4E+09		~Polychlorinated Biphenyls (lowest risk)	1336-36-3									
1.3E+01	E	3.8E-03	E	7.0E-06	E	4.0E-04	E			1.0E+00	1.4E-01		1.4E+09		~Tetrachlorobiphenyl, 3,3',4,4'-(HCB 111)	32588-13-3	5.3E-02	1.4E-01	1.0E+03	3.8E-02	5.5E-01	1.6E+00	5.7E+05	4.1E-01	
3.9E+01	E	1.1E-02	E	2.3E-06	E	1.3E-04	E	V		1.0E+00	1.4E-01		1.4E+09	7.3E+05	~Tetrachlorobiphenyl, 3,4,4',5'-(PCB 81)	70389-60-4	1.8E-02	4.5E-02	1.8E-01	1.2E-02	1.8E-01	5.5E-01	1.0E+02	1.4E-01	
				6.0E-04	I			V		1.0E+00	1.0E-01		1.4E+09		Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9							8.5E+05	8.5E+05	
				6.0E-02	I			V		1.0E+00	1.3E-01		1.4E+09	1.4E+05	Polynuclear Aromatic Hydrocarbons (PAHs)										
				3.0E-01	I			V		1.0E+00	1.3E-01		1.4E+09	5.2E+05	~Acenaphthene	83-32-9					4.7E+03	1.5E+04		3.6E+03	
7.3E-01	E	1.1E-04																							

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _i (mg/m ³)	k _e y	v	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)
		3.4E-05	C	4.0E-03 2.0E-02	I I	3.0E-03	I I	V I		1.0E+00 1.0E+00	1.3E-01 1.3E-01		1.4E+09 1.4E+09	5.8E+04 4.6E+04	-Methylnaphthalene, 2- Naphthalene	91-57-6 91-20-3			3.8E+00 3.8E+00		3.1E+02 1.8E+03	1.0E+03 5.1E+03	1.4E+02	2.4E+02 1.3E+02
1.2E+00	C	1.1E-04	C	3.0E-02 2.0E-02	I P			V I		1.0E+00 1.0E+00	1.3E-01 1.0E-01		1.4E+09 1.4E+09	2.4E+06	-Nitropyrene, 4- Pyrene Potassium Perfluorobutane Sulfonate	57835-92-4 129-00-0 29420-49-3	5.8E-01 1.6E+00 3.5E+04		4.2E-01	2.3E+03 1.6E+03	7.6E+03 6.6E+03		1.8E+03 1.3E+03	
1.5E-01	I			9.0E-03 6.0E-03 1.5E-02	I H I			V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09 1.4E+09	4.2E+05	Prochloraz Profuralin Prometon	67747-09-5 26399-36-0 1610-18-0	4.6E+00 1.6E+01		3.6E+00	7.0E+02 4.7E+02 1.2E+03	3.0E+03 4.9E+03		5.7E+02 4.7E+02 9.5E+02	
				4.0E-03 1.3E-02 5.0E-03	I I I			V I I		1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09 1.4E+09		Prometryn Propachlor Propanil	7287-19-6 1918-16-7 709-98-8				3.1E+02 1.0E+03 3.9E+02	1.3E+03 4.3E+03 1.6E+03		2.5E+02 8.2E+02 3.2E+02	
				2.0E-02 2.0E-03 2.0E-02	I I I			V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01	1.1E+05	1.4E+09 1.4E+09	6.3E+04	Propargite Propargyl Alcohol Propazine	2312-35-8 107-19-7 139-40-2				1.6E+03 1.6E+02 1.6E+03	6.6E+03 6.6E+03		1.3E+03 1.6E+02 1.3E+03	
				2.0E-02 1.3E-02	I I			V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09		Propham Propiconazole	122-42-9 60207-90-1				1.6E+03 1.0E+03	6.6E+03 4.3E+03		1.3E+03 8.2E+02	
				8.0E-03	I	V				1.0E+00		3.3E+04	1.4E+09	8.9E+03	Propionaldehyde	123-38-6						7.5E+01		7.5E+01
				1.0E-01 3.0E+00	X C	1.0E+00 C	V V			1.0E+00 1.0E+00		2.6E+02 3.5E+02	1.4E+09 1.4E+09	7.0E+03 7.0E+02	Propyl benzene Propylene Propylene Glycol	103-65-1 115-07-1 57-55-6				7.8E+03 1.6E+06		7.3E+03 6.6E+06		3.8E+03 1.3E+06
				2.7E-04 7.0E-01 7.0E-01	A H H		V I			1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 8.5E+04 1.1E+05	1.6E+05 7.8E+04	Propylene Glycol Dinitrate Propylene Glycol Monoethyl Ether Propylene Glycol Monomethyl Ether	6423-43-4 1569-02-4 10/-98-2				5.5E+04 5.5E+04		3.9E+05 1.6E+05		3.9E+05 4.1E+04
2.4E-01	I	3.7E-06	I	3.0E-02	I	V				1.0E+00		7.8E+04	1.4E+09	1.0E+04	Propylene Oxide	75-56-9	2.9E+00		7.8E+00	2.1E+00			3.2E+02	3.2E+02
				2.5E-01 2.5E-02	I I			V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09		Pursuit Pydrin	81335-77-5 51630-58-1					2.0E+04 2.0E+03	8.2E+04 8.2E+03		1.6E+04 1.6E+03
3.0E+00	I			1.0E-03 5.0E-04	I I			V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01	5.3E+05	1.4E+09 1.4E+09	5.5E+04	Pyridine Quinalphos Quinone	110-86-1 12593-03-8 91-22-5	2.3E-01	8.2E-01	1.8E-01		7.8E+01 3.9E+01	1.6E+02		7.8E+01 3.2E+01
				3.0E-02 5.0E-02	I H			V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09	4.7E+05	Refractory Ceramic Fibers Resmethrin Ronnel	NA 10453-86-8 299-84-3					2.3E+03 3.9E+03	9.9E+03	4.3E+07	4.3E+07 3.9E+03
2.2E-01	C	6.3E-05	C	4.0E-03 2.5E-02	I I			V I	M	1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09		Rotenone Safrole Savex	83-79-4 94-59-7 7858/-05-0	7.0E-01	2.7E+00	2.2E+04	5.5E-01	3.1E+02 2.0E+03	1.3E+03 8.2E+03		2.5E+02 1.6E+03
				5.0E-03 5.0E-03 5.0E-03	I I C	2.0E-02 C				1.0E+00 1.0E+00			1.4E+09 1.4E+09		Selenious Acid Selenium Selenium Sulfide	7783-00-8 7782-49-2 7446-34-6					3.9E+02 3.9E+02 3.9E+02		2.8E+07 2.8E+07	3.9E+02 3.9E+02
				9.0E-02 5.0E-03	I I			V I		1.0E+00 1.0E+00	1.0E-01 4.0E-02		1.4E+09 1.4E+09		Sethoxydim Silica (crystalline, respirable) Silver	74051-80-2 7631-86-9 7440-22-4					7.0E+03 3.9E+02	3.0E+04 4.3E+06		5.7E+03 3.9E+02
1.2E-01	H			5.0E-03 1.3E-02 4.0E-03	I I I			V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09		Simazine Sodium Acifluorfen Sodium Azide	122-34-9 62476-59-9 26828-22-8	5.8E+00	2.1E+01		4.5E+00	3.9E+02 1.0E+03 3.1E+02	1.6E+03 4.3E+03		3.2E+02 8.2E+02 3.1E+02
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M		2.5E-02			1.4E+09		Sodium Dichromate	10588-01-9	3.1E-01		9.2E+00	3.0E-01	1.6E+03		2.8E+05	1.6E+03
2.7E-01	H			3.0E-02 5.0E-02	I A	1.3E-02	C			1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09		Sodium Diethylthiocarbamate Sodium Fluoride	148-18-5 7681-49-4	2.6E+00	9.2E+00	2.0E+00	2.3E+03 3.9E+03	9.9E+03	1.8E+07	1.9E+03 3.9E+03	
2.4E-02	H			2.0E-05 1.0E-03 3.0E-02	I H I			V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09		Sodium Fluoroacetate Sodium Metavanadate Stirofos (Tetrachlorovinphos)	62-74-8 13718-26-8 961-11-5	2.9E+01	1.0E+02		2.3E+01	1.6E+00 2.3E+03 2.3E+03	6.6E+00 7.8E+01 9.9E+03		1.3E+00 7.8E+01 1.9E+03
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M		2.5E-02			1.4E+09		Strontium Chromate	7789-06-2	3.1E-01		9.2E+00	3.0E-01	1.6E+03		2.8E+05	1.6E+03
				6.0E-01 3.0E-04	I I			V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09		Strontium, Stable Strychnine	7440-24-6 57-24-9					4.7E+04 2.3E+01	9.9E+01		4.7E+04 1.9E+01
				2.0E-01 3.0E-03 1.0E-03	I P P	1.0E+00 2.0E-03	I X	V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01	8.7E+02	1.4E+09 1.4E+09	9.4E+03	Styrene Styrene-Acrylonitrile (SAN) Trimer Sulfolane	100-42-5 NA 126-33-0					1.6E+04 2.3E+02 7.8E+01	9.7E+03 9.9E+02 3.3E+02	2.8E+06	6.0E+03 1.9E+02 6.3E+01
				8.0E-04	P			V		1.0E+00	1.0E-01		1.4E+09		Sulfonylbis(4-chlorobenzene), 1,1'- Sulfur Trioxide Sulfuric Acid	80-07-9 7446-11-9 7664-93-9					6.3E+01	2.6E+02	1.4E+06 1.4E+06	5.1E+01 1.4E+06 1.4E+06
				2.5E-02 3.0E-02 7.0E-02	I H I			V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09		Systhane TCMTB Tebuthiuron	88671-89-0 21564-17-0 34014-18-1					2.0E+03 2.3E+03 5.5E+03	8.2E+03 9.9E+03 2.3E+04		1.6E+03 1.9E+03 4.4E+03
				2.0E-02 1.3E-02	H I			V I		1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09		Temphos Terbacil	3383-96-8 5902-51-2					1.6E+03 1.0E+03	6.6E+03 4.3E+03		1.3E+03 8.2E+02

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1									
SFO (mg/kg-day) ⁻¹	k _e (y ⁻¹)	IUR (ug/m ³) ⁻¹	k _e (y ⁻¹)	RfD _o (mg/kg-day)	k _e (y ⁻¹)	RfC _i (mg/m ³)	k _e (y ⁻¹)	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)					
				2.5E-05	H		V		1.0E+00		3.1E+01	1.4E+09	2.6E+05	Terbutolol	13071-79-9									2.0E+00				2.0E+00
				1.0E-03	I				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Terbutryn	886-50-0									7.8E+01	3.3E+02			6.3E+01
				1.0E-04	I				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Tetrabromodiphenyl ether, 2,2',4,4'-(BDE-47)	5436-43-1									7.8E+00	3.3E+01			6.3E+00
				3.0E-04	I		V		1.0E+00		1.4E+09	5.1E+04	1.4E+09	Tetrachlorobenzene, 1,2,4,5-	95-94-3									2.3E+01				2.3E+01
2.6E-02	I	7.4E-06	I	3.0E-02	I		V		1.0E+00		6.8E+02	1.4E+09	5.7E+03	Tetrachloroethane, 1,1,1,2-	630-20-6	2.7E+01		2.2E+00	2.0E+00	2.3E+03								2.3E+03
2.0E-01	I	5.8E-05	C	2.0E-02	I		V		1.0E+00		1.9E+03	1.4E+09	1.5E+04	Tetrachloroethane, 1,1,2,2-	79-34-5	3.5E+00		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	1.0E+00		1.7E+02	1.4E+09	2.4E+03	Tetrachloroethylene	127-18-4	3.3E+02		2.5E+01	2.4E+01	4.7E+02			9.8E+01					8.1E+01
2.0E+01	H			3.0E-02	I				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Tetrachlorophenol, 2,3,4,6-	58-90-2				3.5E-02	3.5E-02	2.3E+03	9.9E+03						1.9E+03
				5.0E-04	I		V		1.0E+00	1.0E-01	1.4E+09	1.1E+05	1.4E+09	Tetrachlorotoluene, p- alpha, alpha- Tetraethyl Dithiopyrophosphate	5216-25-1 3689-24-5	3.5E-02				3.9E+01	1.6E+02							3.2E+01
				8.0E+01	I	V			1.0E+00		1.1E+03	1.4E+09	1.2E+03	Tetrafluoroethane, 1,1,1,2-	811-97-2													1.0E+05
				2.0E-03	P				1.0E+00	6.5E-04	1.4E+09	1.4E+09	1.4E+09	Tetryl (Trinitrophenylmethyltriamine)	479-45-8					1.6E+02	1.0E+05							1.6E+02
				7.0E-06	X				1.0E+00		1.4E+09	1.4E+09	1.4E+09	Thallium (I) Nitrate	10102-45-1					5.5E-01								5.5E-01
				1.0E-05	X				1.0E+00		1.4E+09	1.4E+09	1.4E+09	Thallium (Soluble Salts)	7440-28-0					7.8E-01								7.8E-01
				6.0E-06	X		V		1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Thallium Acetate	563-68-8					4.7E-01	2.0E+00							3.8E-01
				2.0E-05	X				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Thallium Carbonate	6533-73-9					1.6E+00	6.6E+00							1.3E+00
				6.0E-06	X				1.0E+00		1.4E+09	1.4E+09	1.4E+09	Thallium Chloride	7791-12-0					4.7E-01								4.7E-01
				2.0E-05	X				1.0E+00		1.4E+09	1.4E+09	1.4E+09	Thallium Sulfate	7446-18-6					1.6E+00								1.6E+00
				1.0E-02	I				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Thiobencarb	28249-77-6					7.8E+02	3.3E+03							6.3E+02
				7.0E-02	X				1.0E+00	7.5E-03	1.4E+09	1.4E+09	1.4E+09	Thiodiglycol	111-48-8					5.5E+03	3.1E+05							5.4E+03
				3.0E-04	H				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Thiofanox	39196-18-4					2.3E+01	9.9E+01							1.9E+01
				8.0E-02	I				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Thiophanate, Methyl	23564-05-8					6.3E+03	2.6E+04							5.1E+03
				5.0E-03	I				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Thiram	137-26-8					3.9E+02	1.6E+03							3.2E+02
				6.0E-01	H				1.0E+00		1.4E+09	1.4E+09	1.4E+09	Tin	7440-31-5					4.7E+04								4.7E+04
				1.0E-04	A	V			1.0E+00		1.4E+09	1.4E+09	1.4E+09	Titanium tetrachloride	7500-45-0													1.4E+05
1.8E-01	X			2.0E-04	X				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Toluene	108-88-3	3.9E+00	1.4E+01		3.0E+00	6.3E+03	2.2E+04							4.9E+03
3.0E-02	P			4.0E-03	X				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Toluene, 2,5-diamine	95-70-5	2.3E+01	8.2E+01		1.8E+01	1.6E+01	6.6E+01							1.3E+01
				3.0E+00	P		V		1.0E+00		3.4E-01	1.4E+09	1.1E+03	Toluidine, p-	106-49-0					3.1E+02	1.3E+03							2.5E+02
				1.0E-02	X	6.0E-01	P	V	1.0E+00		1.4E+02	1.4E+09	8.3E+02	Total Petroleum Hydrocarbons (Aliphatic High)	NA					2.3E+05								2.3E+05
				1.0E-02	X	1.0E-01	P	V	1.0E+00		6.9E+00	1.4E+09	1.0E+03	Total Petroleum Hydrocarbons (Aliphatic Low)	NA					7.8E+02								5.2E+02
				4.0E-02	P				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Total Petroleum Hydrocarbons (Aliphatic Medium)	NA													9.6E+01
				4.0E-03	P	3.0E-02	P	V	1.0E+00		1.8E+03	1.4E+09	3.5E+03	Total Petroleum Hydrocarbons (Aromatic High)	NA					3.1E+03	1.3E+04							2.5E+03
				4.0E-03	P	3.0E-03	P	V	1.0E+00		1.4E+09	5.2E+04	1.4E+09	Total Petroleum Hydrocarbons (Aromatic Low)	NA					3.1E+02								8.2E+01
				4.0E-03	P	3.0E-03	P	V	1.0E+00		1.4E+09	5.2E+04	1.4E+09	Total Petroleum Hydrocarbons (Aromatic Medium)	NA					3.1E+02								1.1E+02
1.1E+00	I	3.2E-04	I						1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Toxaphene	9001-35-2	6.3E-01	2.2E+00	1.2E+04	4.9E-01	5.9E+02	2.5E+03							4.7E+02
				7.5E-03	I				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Tralomehrin	68841-25-6					2.3E+01								2.3E+01
				3.0E-04	A		V		1.0E+00		1.4E+09	3.4E+03	1.4E+09	Tri-n-butyltin	688-73-3													
				8.0E+01	X				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Triacetin	102-76-1					6.3E+06	2.6E+07							5.1E+06
				1.3E-02	I		V		1.0E+00		1.4E+09	3.6E+05	1.4E+09	Triallate	2303-17-5					1.0E+03								1.0E+03
				1.0E-02	I				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Triasuturon	8209-50-5					7.8E+02	3.3E+03							6.3E+02
9.0E-03	P			5.0E-03	I		V		1.0E+00		1.4E+09	4.5E+04	1.4E+09	Tribromobenzene, 1,2,4-	615-54-3					3.9E+02								3.9E+02
				1.0E-02	P				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Tributyl Phosphate	126-73-8	7.7E+01	2.7E+02		6.0E+01	7.8E+02	3.3E+03							6.3E+02
				3.0E-04	P				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Tributyltin Compounds	NA					2.3E+01	9.9E+01							1.9E+01
				3.0E-04	I				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Tributyltin Oxide	56-35-9					2.3E+01	9.9E+01							1.9E+01
7.0E-02	I			3.0E+01	I	3.0E+01	H	V	1.0E+00		9.1E+02	1.4E+09	1.3E+03	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1					2.3E+06								4.0E+04
				2.0E-02	I				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Trichloroacetic Acid	76-03-9	9.9E+00	3.5E+01		7.8E+00	1.6E+03	6.6E+03							1.3E+03
2.9E-02	H			3.0E-05	X				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Trichloroaniline HCl, 2,4,6-	33663-50-2	2.4E+01	8.5E+01		1.9E+01	2.3E+00	9.9E+00							1.9E+00
7.0E-03	X			8.0E-04	X				1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.4E+09	Trichloroaniline, 2,4,6-	634-93-5	9.9E+01	3.5E+02		7.8E+01	6.3E+01	9.9E+00							6.3E+01

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Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg-day)	k e y	RfC _i (mg/m ³)	k e y	v o l a t i l e	m u t a g e n	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL Child HQ=1 (mg/kg)	Dermal SL Child HQ=1 (mg/kg)	Inhalation SL Child HQ=1 (mg/kg)	Noncarcinogenic SL Child HI=1 (mg/kg)
7.7E-03 2.0E-02	I P			2.0E+00 7.5E-03 1.0E-02	P I P					1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01 1.0E-01		1.4E+09 1.4E+09 1.4E+09	5.1E+05	Triethylene Glycol Trifluralin Trimethyl Phosphate	112-27-6 1582-09-8 512-56-1	9.0E+01 3.5E+01	1.2E+02		9.0E+01 2.7E+01	1.6E+05 5.9E+02 7.8E+02	6.6E+05 3.3E+03		1.3E+05 5.9E+02 6.3E+02
				5.0E-03 7.0E-03 1.0E-02	P V P V X					1.0E+00 1.0E+00 1.0E+00	2.9E+02 2.2E+02 1.8E+02	1.4E+09 1.4E+09 1.4E+09	9.4E+03 7.9E+03 6.6E+03		Trimethylbenzene, 1,2,3- Trimethylbenzene, 1,2,4- Trimethylbenzene, 1,3,5-	526-73-8 95-63-6 108-67-8						4.9E+01 5.8E+01 7.8E+02	4.9E+01 5.8E+01 7.8E+02	
3.0E-02	I			3.0E-02 5.0E-04 2.0E-02	I I P					1.0E+00 1.0E+00 1.0E+00	1.9E-02 3.2E-02 1.0E-01		1.4E+09 1.4E+09 1.4E+09		Trinitrobenzene, 1,3,5- Trinitrotoluene, 2,4,6- Triphenylphosphine Oxide	99-35-4 118-96-7 791-28-6	2.3E+01	2.6E+02	2.1E+01	2.3E+03 3.9E+01 1.6E+03	5.2E+04 5.2E+02 6.6E+03		2.2E+03 3.6E+01 1.3E+03	
2.3E+00	C	6.6E-04	C	2.0E-02 1.0E-02	A X					1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01 1.0E-01	4.7E+02	1.4E+09 1.4E+09 9.0E+05		Tris(1,3-Dichloro-2-propyl) Phosphate Tris(1-chloro-2-propyl)phosphate Tris(2,3-dibromopropyl)phosphate	13874-87-8 13674-84-5 126-72-1	3.0E-01	3.8E+00	2.8E-01	1.6E+03 7.8E+02	6.6E+03 3.3E+03		1.3E+03 6.3E+02	
2.0E-02 3.2E-03	P P			7.0E-03 1.0E-01 3.0E-03	P P I					1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01 4.0E-05		1.4E+09 1.4E+09 1.4E+09		Tris(2-chloroethyl)phosphate Tris(2-ethylhexyl)phosphate Uranium (Soluble Salts)	145-95-8 78-42-2 NA	3.5E+01 2.2E+02	1.2E+02 7.7E+02	2.7E+01 1.7E+02	5.5E+02 7.8E+03 2.3E+02	2.3E+03 3.3E+04		4.4E+02 6.3E+03 2.3E+02	
1.0E+00	C	2.9E-04 8.3E-03	C P	9.0E-03 5.0E-03	I S	7.0E-06 1.0E-04	P A			1.0E+00 2.6E-02 2.6E-02	1.0E-01 1.4E+09 1.4E+09		1.4E+09 1.4E+09 1.4E+09		Urethane Vanadium Pentoxide Vanadium and Compounds	51-79-6 1314-62-1 1440-62-2	1.5E-01	6.0E-01	4.8E+03 4.6E+02	1.2E-01 4.6E+02	7.0E+02 3.9E+02		9.9E+03 1.4E+05 3.9E+02	6.6E+02 3.9E+02
				1.0E-03 2.5E-02 1.0E+00	I I H	2.0E-01 1.0E-01	I V			1.0E+00 1.0E+00 1.0E+00	1.4E+09 1.4E+09 2.8E+03	1.2E+05 1.4E+09 4.4E+03			Verapamil Vinclozolin Vinyl Acetate	1929-77-7 50471-44-8 108-05-4				7.8E+01 2.0E+03 7.8E+04	8.2E+03		7.8E+01 1.6E+03 9.1E+02	
7.2E-01	I	3.2E-05 4.4E-06	H I	3.0E-03 3.0E-04	I I	1.0E-01 1.0E-01	I V M			1.0E+00 1.0E+00 1.0E+00	3.4E+03 3.9E+03	1.4E+09 1.4E+09 9.6E+02	1.4E+03 9.6E+02		Vinyl Bromide Vinyl Chloride Warfarin	593-60-2 75-01-4 81-81-2	9.4E-02		1.2E-01 1.6E-01	1.2E-01 5.9E-02	2.3E+02 2.3E+01	9.9E+01		4.3E+00 7.0E+01 1.9E+01
				2.0E-01 2.0E-01 2.0E-01	S S S	1.0E-01 1.0E-01 1.0E-01	S V S			1.0E+00 1.0E+00 1.0E+00	3.9E+02 3.9E+02 4.3E+02	1.4E+09 1.4E+09 6.5E+03	5.8E+03 5.5E+03		Xylene, p- Xylene, m- Xylene, o-	106-42-3 108-38-3 95-47-6				1.6E+04 1.6E+04 1.6E+04		5.8E+02 5.7E+02 6.7E+02	5.6E+02 5.5E+02 6.5E+02	
				2.0E-01 3.0E-04 3.0E-01	I I I	1.0E-01 1.0E-01	I V			1.0E+00 1.0E+00 1.0E+00	2.6E+02	1.4E+09 1.4E+09 1.4E+09	6.5E+03		Xylenes Zinc Phosphate Zinc and Compounds	1330-20-7 1314-84-7 7440-66-6				1.6E+04 2.3E+01 2.3E+04		6.8E+02	6.5E+02 2.3E+01 2.3E+04	
				5.0E-02 8.0E-05	I X					1.0E+00 1.0E+00	1.0E-01 1.0E-01		1.4E+09 1.4E+09		Zineb Zirconium	12122-67-7 7440-67-7				3.9E+03 6.3E+00	1.6E+04		3.2E+03 6.3E+00	

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Toxicity and Chemical-specific					Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y	v o l a t i l e	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
5.1E-06	C				ALAR	1596-84-5	5.5E-01	
2.2E-06	I	9.0E-03	I	V	Acephate	30560-19-1		
					Acetaldehyde	75-07-0	1.3E+00	9.4E+00
		3.1E+01	A	V	Acetochlor	34256-82-1		
		2.0E-03	X	V	Acetone	67-64-1		3.2E+04
					Acetone Cyanohydrin	75-86-5		2.1E+00
		6.0E-02	I	V	Acetonitrile	75-05-8		6.3E+01
1.3E-03	C			V	Acetophenone	98-86-2	2.2E-03	
					Acetylaminofluorene, 2-	53-96-3		
		2.0E-05	I	V	Acrolein	107-02-8		2.1E-02
1.0E-04	I	6.0E-03	I		Acrylamide	79-06-1	1.0E-02	6.3E+00
		1.0E-03	I	V	Acrylic Acid	79-10-7		1.0E+00
6.8E-05	I	2.0E-03	I	V	Acrylonitrile	107-13-1	4.1E-02	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-04	
		1.0E-04	X	V	Allyl	74223-64-6		
					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		
					Amdro	67485-29-4		
					Ametryn	834-12-8		
6.0E-03	C				Aminobiphenyl, 4-	92-67-1	4.7E-04	
					Aminophenol, m-	591-27-5		
					Aminophenol, p-	123-30-8		
					Amitraz	33089-61-1		
		1.0E-01	I	V	Ammonia	7664-41-7		1.0E+02
		3.0E-03	X	V	Ammonium Sulfamate	7773-06-0		
					Amyl Alcohol, tert-	75-85-4		3.1E+00
1.6E-06	C	1.0E-03	I		Aniline	62-53-3	1.8E+00	1.0E+00
					Anthraquinone, 9,10-	84-85-1		
					Antimony (metallic)	7440-36-0		
					Antimony Pentoxide	1314-60-9		
					Antimony Potassium Tartrate	11071-15-1		
					Antimony Tetroxide	1332-81-6		
		2.0E-04	I		Antimony Trioxide	1309-64-4		2.1E-01
7.1E-06	I				Apollo	74115-24-5		
					Aramite	140-57-8	4.0E-01	
4.3E-03	I	1.5E-05	C		Arsenic, Inorganic	7440-38-2	6.5E-04	1.6E-02
		5.0E-05	I		Arsine	7784-42-1		5.2E-02
					Assure	76578-14-8		
					Asulam	3337-71-1		
2.5E-04	C				Atrazine	1912-24-9	1.1E-02	
					Auramine	492-80-8		
					Avermectin B1	65195-55-3		
3.1E-05	I			V	Azobenzene	103-33-3	9.1E-02	
		7.0E-06	P		Azodicarbonamide	123-77-3		7.3E-03
1.5E-01	C	2.0E-04	C		Barium	7440-39-3	6.8E-06	5.2E-01
					Barium Chromate	10294-40-3		2.1E-01
					Baygon	114-26-1		
					Bayleton	43121-43-3		
				V	Baythroid	68359-37-5		
					Benefin	1861-40-1		
				V	Benomyl	17804-35-2		
					Bentazon	25057-89-0		
					Benzaldehyde	100-52-7		
7.8E-06	I	3.0E-02	I	V	Benzene	71-43-2	3.6E-01	3.1E+01
				V	Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1		
					Benzenethiol	108-98-5		
6.7E-02	I			M	Benzidine	92-87-5	1.5E-05	
				V	Benzoic Acid	65-85-0		
					Benzotrithloride	98-07-7		
					Benzyl Alcohol	100-51-6		
4.9E-05	C	1.0E-03	P	V	Benzyl Chloride	100-44-7	5.7E-02	1.0E+00
2.4E-03	I	2.0E-05	I		Beryllium and compounds	7440-41-7	1.2E-03	2.1E-02
					Bidrin	141-66-2		
					Bifenox	42576-02-3		
					Biphenthrin	82657-04-3		
		4.0E-04	X	V	Biphenyl, 1,1'-	92-52-4		4.2E-01
1.0E-05	H			V	Bis(2-chloro-1-methylethyl) ether	108-60-1	2.8E-01	
					Bis(2-chloroethoxy)methane	111-91-1		

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Toxicity and Chemical-specific						Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³) ³	k v o c	v o l a t i l e	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
3.3E-04	I			V		Bis(2-chloroethyl)ether	111-44-4	8.5E-03	
6.2E-02	I			V		Bis(chloromethyl)ether	542-88-1	4.5E-05	
						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-04	X			V		Bromo-2-chloroethane, 1-	107-04-0	4.7E-03	
		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-02	
1.1E-06	I			V		Bromodiform	75-25-2	2.6E+00	
		5.0E-03	I	V		Bromomethane	74-83-9		5.2E+00
				V		Bromophos	2104-96-3		
				V		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-02	2.1E+00
				V		Butanol, N-	71-36-3		
		3.0E+01	P	V		Butyl Benzyl Phthlate	85-68-7		
				V		Butyl alcohol, sec-	78-92-2		3.1E+04
				V		Butylate	2008-41-5		
5.7E-08	C			V		Butylated hydroxyanisole	25013-16-5	4.9E+01	
				V		Butylated hydroxytoluene	128-37-0		
				V		Butylbenzene, n-	104-51-8		
				V		Butylbenzene, sec-	135-98-8		
				V		Butylbenzene, tert-	98-06-6		
				V		Cacodylic Acid	75-60-5		
1.8E-03	I	1.0E-05	A			Cadmium (Diet)	7440-43-9		
1.8E-03	I	1.0E-05	A			Cadmium (Water)	7440-43-9	1.6E-03	1.0E-02
1.5E-01	C	2.0E-04	C		M	Calcium Chromate	13765-19-0	6.8E-06	2.1E-01
		2.2E-03	C			Caprolactam	105-60-2		2.3E+00
4.3E-05	C					Captafol	2425-06-1	6.5E-02	
6.6E-07	C					Captaf	133-06-2	4.3E+00	
		7.0E-01	I	V		Carbaryl	83-26-2		7.3E+02
						Carboran	1563-66-2		
						Carbon Disulfide	75-15-0		
6.0E-06	I	1.0E-01	I	V		Carbon Tetrachloride	56-23-5	4.7E-01	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		
1.0E-04	I	7.0E-04	I	V		Chloranil	118-75-2	2.8E-02	7.3E-01
4.6E-03	C					Chlorodane	12789-03-6	6.1E-04	
						Chlordecone (Kepone)	143-50-0		
		1.5E-04	A	V		Chlorfenvinphos	470-90-6		1.5E-01
						Chlorimuron, Ethyl-	90982-32-4		
						Chlorine	7782-50-5		
		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-03	2.1E+01
						Chloro-2-methylaniline HCl, 4-	3165-93-3		
7.7E-05	C					Chloro-2-methylaniline, 4-	95-69-2	3.6E-02	
						Chloroacetaldehyde, 2-	107-20-0		
		3.0E-05	I			Chloroacetic Acid	79-11-8		3.1E-02
						Chloroacetophenone, 2-	532-27-4		
3.1E-05	C	5.0E-02	P	V		Chloroaniline, p-	106-47-8		5.2E+01
						Chlorobenzene	108-90-7		
						Chlorobenzilate	510-15-6	9.1E-02	
		3.0E-01	P	V		Chlorobenzoic Acid, p-	74-11-3		3.1E+02
				V		Chlorobenzotrifluoride, 4-	98-56-6		
						Chlorobutane, 1-	109-69-3		
		5.0E+01	I	V		Chlorodifluoromethane	75-45-6		5.2E+04
2.3E-05	I	9.8E-02	A	V		Chloroethanol, 2-	107-07-3		1.0E+02
						Chloroform	67-66-3	1.2E-01	
6.9E-04	C	9.0E-02	I	V		Chloromethane	74-87-3		9.4E+01
		1.0E-05	X			Chloromethyl Methyl Ether	107-30-2	4.1E-03	
						Chloronitrobenzene, o-	88-73-3		1.0E-02
		6.0E-04	P			Chloronitrobenzene, p-	100-00-5		6.3E-01
				V		Chlorophenol, 2-	95-57-8		
		4.0E-04	C	V		Chloropicrin	76-06-2		4.2E-01
8.9E-07	C			V		Chlorothalonil	1897-45-6	3.2E+00	
				V		Chlorotoluene, o-	95-49-8		
				V		Chlorotoluene, p-	106-43-4		

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Toxicity and Chemical-specific				Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1		
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y	v o l u t i l e	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
6.9E-02	C					Chlorozotocin Chlorpropham Chlorpyrifos	54749-90-5 101-21-3 2921-88-2	4.1E-05	
						Chlorpyrifos Methyl Chlorsulfuron Chlorthiophos	5598-13-0 64902-72-3 60238-56-4		
8.4E-02	S	1.0E-04	I		M	Chromium(III), Insoluble Salts Chromium(VI) Chromium, Total	16065-83-1 18540-29-9 7440-47-3	1.2E-05	1.0E-01
9.0E-03	P	6.0E-06	P			Cobalt	7440-48-4	3.1E-04	6.3E-03
6.2E-04	I		V		M	Coke Oven Emissions Copper	8007-45-2 7440-50-8	1.6E-03	
		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
		6.0E-01	C			Cresol, p-chloro-m-	59-50-7		6.3E+02
				V		Cresols Crotonaldehyde, trans-	1319-77-3 123-73-9		6.3E+02
6.3E-05	C	4.0E-01	I	V		Cumene Cupferron Cyanazine	98-82-8 135-20-6 21725-46-2	4.5E-02	4.2E+02
						Cyanides ~Calcium Cyanide ~Copper Cyanide	592-01-8 544-92-3		
8.0E-04	S		V			~Cyanide (CN-) ~Cyanogen ~Cyanogen Bromide	57-12-5 460-19-5 506-68-3		8.3E-01
						~Cyanogen Chloride ~Hydrogen Cyanide ~Potassium Cyanide	506-77-4 74-90-8 151-50-8		8.3E-01
						~Potassium Silver Cyanide ~Silver Cyanide ~Sodium Cyanide	506-61-6 506-64-9 143-33-9		
					Y	~Thiocyanates ~Thiocyanic Acid ~Zinc Cyanide	NA 463-56-9 567-21-1		
6.0E+00	I		V			Cyclohexane	110-82-7		6.3E+03
		7.0E-01	P	V		Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro- Cyclohexanone	87-84-3 108-94-1		7.3E+02
		1.0E+00	X	V		Cyclohexene Cyclohexylamine Cyhalothrin/karate	110-83-8 108-91-8 68085-85-8		1.0E+03
						Cypermethrin Cyrromazine DDD	52315-07-8 66215-27-8 72-54-8	4.1E-02	
6.9E-05	C					DDE, p,p'- DDT Dacthal	72-55-9 50-29-3 1861-32-1	2.9E-02 2.9E-02	
						Dalapon Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209) Demeton	75-99-0 1163-19-5 8065-48-3		
						Di(2-ethylhexyl)adipate Diallate Diazinon	103-23-1 2303-16-4 333-41-5		
6.0E-03	P	2.0E-04	I	V	M	Dibenzothiophene Dibromo-3-chloropropane, 1,2- Dibromobenzene, 1,3-	132-65-0 96-12-8 108-36-1	1.7E-04	2.1E-01
						Dibromobenzene, 1,4- Dibromochloromethane Dibromoethane, 1,2-	106-37-6 124-48-1 106-93-4	1.0E-01 4.7E-03	9.4E+00
		4.0E-03	X	V		Dibromomethane (Methylene Bromide) Dibutyltin Compounds Dicamba	74-95-3 NA 1918-00-9		4.2E+00
4.2E-03	P		V			Dichloro-2-butene, 1,4-	764-41-0	6.7E-04	
4.2E-03	P		V			Dichloro-2-butene, cis-1,4-	1476-11-5	6.7E-04	
4.2E-03	P		V			Dichloro-2-butene, trans-1,4-	110-57-6	6.7E-04	
		2.0E-01	H	V		Dichloroacetic Acid	79-43-6		2.1E+02
1.1E-05	C	8.0E-01	I	V		Dichlorobenzene, 1,2- Dichlorobenzene, 1,4-	95-50-1 106-46-7	2.6E-01	8.3E+02
3.4E-04	C					Dichlorobenzidine, 3,3'- Dichlorobenzophenone, 4,4'- Dichlorodifluoromethane	91-94-1 90-98-2 75-71-8	8.3E-03	1.0E+02
1.6E-06	C			V		Dichloroethane, 1,1-	75-34-3	1.8E+00	
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

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Toxicity and Chemical-specific						Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³) ⁻¹	k e y	v o l a t i l e	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
				V		Dichloroethylene, 1,2-cis-	156-59-2		
				V		Dichloroethylene, 1,2-trans-	156-60-5		
						Dichlorophenol, 2,4-	120-83-2		
1.0E-05	C	4.0E-03	I	V		Dichlorophenoxy Acetic Acid, 2,4-	94-75-7		
						Dichlorophenoxybutyric Acid, 4-(2,4-	94-82-6		
						Dichloropropane, 1,2-	78-87-5	2.8E-01	4.2E+00
4.0E-06	I	2.0E-02	I	V		Dichloropropane, 1,3-	142-28-9		
						Dichloropropanol, 2,3-	616-23-9		
						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-02	5.2E-01
						Dicyclopentadiene	77-73-6		3.1E-01
4.6E-03	I	3.0E-04	X	V		Dieldrin	60-57-1	6.1E-04	
3.0E-04	C	5.0E-03	I			Diesel Engine Exhaust	NA	9.4E-03	5.2E+00
						Diethanolamine	111-42-2		2.1E-01
						Diethylene Glycol Monobutyl Ether	112-34-5		1.0E-01
						Diethylene Glycol Monoethyl Ether	111-90-0		3.1E-01
1.0E-01	C	3.0E-04	P	V		Diethylformamide	617-84-5		
						Diethylstilbestrol	56-53-1	2.8E-05	
4.0E+01	I	4.0E+01	V			Difenzoquat	43222-48-6		
						Difflubenzuron	35367-38-5		
						Diffluoroethane, 1,1-	75-37-6		4.2E+04
1.3E-05	C	7.0E-01	P	V		Dihydrosafrole	94-58-6	2.2E-01	
						Diisopropyl Ether	108-20-3		7.3E+02
						Diisopropyl Methylphosphonate	1445-75-6		
1.3E-03	C					Dimethipin	55290-64-7		
						Dimethoate	60-51-5		
						Dimethoxybenzidine, 3,3'-	119-90-4		
						Dimethyl methylphosphonate	756-79-6		
						Dimethylamino azobenzene [p-]	60-11-7	2.2E-03	
						Dimethylaniline HCl, 2,4-	21436-96-4		
						Dimethylaniline, 2,4-	95-68-1		
						Dimethylaniline, N,N-	121-69-7		
						Dimethylbenzidine, 3,3'-	119-93-7		
3.0E-02	I	3.0E-02	V			Dimethylformamide	68-12-2		3.1E+01
2.0E-06	X	2.0E-06	V			Dimethylhydrazine, 1,1-	57-14-7		2.1E-03
1.6E-01	C	1.6E-01	V			Dimethylhydrazine, 1,2-	540-73-8	1.8E-05	
						Dimethylphenol, 2,4-	105-67-9		
						Dimethylphenol, 2,6-	576-26-1		
						Dimethylphenol, 3,4-	95-65-8		
1.3E-05	C	1.3E-05	V			Dimethylvinylchloride	513-37-1	2.2E-01	
						Dinitro- <i>o</i> -cresol, 4,6-	534-52-1		
						Dinitro- <i>p</i> -cyclohexyl Phenol, 4,6-	131-89-5		
						Dinitrobenzene, 1,2-	528-29-0		
						Dinitrobenzene, 1,3-	99-65-0		
						Dinitrobenzene, 1,4-	100-25-4		
8.9E-05	C	8.9E-05				Dinitrophenol, 2,4-	51-28-5		
						Dinitrotoluene Mixture, 2,4/2,6-	NA		
						Dinitrotoluene, 2,4-	121-14-2	3.2E-02	
						Dinitrotoluene, 2,6-	606-20-2		
						Dinitrotoluene, 2-Amino-4,6-	35572-78-2		
						Dinitrotoluene, 4-Amino-2,6-	19400-51-0		
5.0E-06	I	3.0E-02	I	V		Dinitrotoluene, Technical grade	25321-14-6		
						Dinoseb	88-85-7		
						Dioxane, 1,4-	123-91-1	5.6E-01	3.1E+01
1.3E+00	I	1.3E+00				Dioxins			
3.8E+01	C	4.0E-08	C	V		~Hexachlorodibenzo-p-dioxin, Mixture	NA	2.2E-06	
						~TCDD, 2,3,7,8-	1746-01-6	7.4E-08	4.2E-05
2.2E-04	I	2.2E-04				Diphenamid	957-51-7		
						Diphenyl Sulfone	127-63-9		
						Diphenylamine	122-39-4		
1.4E-01	C	1.4E-01				Diphenylhydrazine, 1,2-	122-66-7	1.3E-02	
						Diquat	85-00-7		
1.4E-01	C	1.4E-01				Direct Black 38	1937-37-7	2.0E-05	
1.4E-01	C	1.4E-01				Direct Blue 6	2602-46-2	2.0E-05	
						Direct Brown 95	16071-86-6	2.0E-05	
						Disulfoton	298-04-4		
						Dithiane, 1,4-	505-29-3		
						Diuron	330-54-1		
						Dodine	2439-10-3		
						EPTC	759-94-4		
						Endosulfan	115-29-7		
						Endothall	145-73-3		
1.2E-06	I	1.0E-03	I	V		Endrin	72-20-8		1.0E+00
						Epichlorohydrin	106-89-8	2.3E+00	2.1E+01
						Epoxybutane, 1,2-	106-88-7		

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Toxicity and Chemical-specific		Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1		
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³) ³	k e y c mutagen	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
6.0E-02	P	V		Ethephon	16672-87-0		
				Ethion	563-12-2		
				Ethoxyethanol Acetate, 2-	111-15-9		6.3E+01
2.0E-01	I	V		Ethoxyethanol, 2-	110-80-5		2.1E+02
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
1.0E+01	I	V		Ethyl Chloride (Chloroethane)	75-00-3		1.0E+04
				Ethyl Ether	60-29-7		
3.0E-01	P	V		Ethyl Methacrylate	97-63-2		3.1E+02
2.5E-06	C	1.0E+00	I	Ethyl-p-nitrophenyl Phosphonate	2104-64-5	1.1E+00	1.0E+03
				Ethylbenzene	100-41-4		
				Ethylene Cyanohydrin	109-78-4		
				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
8.8E-05	C	3.0E-02	C	Ethylene Oxide	75-21-8	3.2E-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-04	
				Ethylphthalyl Ethyl Glycolate	84-72-0		
				Express	101200-48-0		
				Fenamiphos	22224-92-6		
1.3E-02	C			Fenproprathrin	39515-41-8		
				Fluometuron	2164-17-2		
1.3E-02	C			Fluoride (Soluble Fluoride)	7782-41-4		1.4E+01
				Fluridone	59756-60-4		
				Flurprimidol	56425-91-3		
				Flutolanil	66332-96-5		
				Fluvalinate	69409-94-5		
				Folpet	133-07-3		
1.3E-05	I	9.8E-03	A	Fomesafen	72178-02-0	2.2E-01	1.0E+01
				Fonofos	944-22-9		
				Formaldehyde	50-00-0		
3.0E-04	X	V		Formic Acid	64-18-6		3.1E-01
				FosetylAL	39148-24-8		
				Furans			
				~Dibenzofuran	132-64-9		
				~Furan	110-00-9		
2.0E+00	I	V		~Tetrahydrofuran	109-99-9		2.1E+03
				Furazolidone	67-45-8		
4.3E-04	C	5.0E-02	H	Furfural	98-01-1	6.5E-03	5.2E+01
				Furium	531-82-8		
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	H	V		Glycidyl	765-34-4		1.0E+00
				Glyphosate	1071-83-6		
				Goal	42874-03-3		
				Guanidine	113-00-8		
1.0E-02	A			Guanidine Chloride	50-01-1		1.0E+01
				Guthion	86-50-0		
1.3E-03	I		V	Haloxyp, Methyl	69806-40-2	2.2E-03	
				Harmony	79277-27-3		
2.6E-03	I		V	Heptachlor	76-44-8		
				Heptachlor Epoxide	1024-57-3	1.1E-03	
				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-03	
2.2E-05	I		V	Hexachlorobutadiene	87-68-3	1.3E-01	
1.8E-03	I			Hexachlorocyclohexane, Alpha-	319-84-6	1.6E-03	
5.3E-04	I			Hexachlorocyclohexane, Beta-	319-85-7	5.3E-03	
3.1E-04	C			Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	9.1E-03	
5.1E-04	I			Hexachlorocyclohexane, Technical	608-73-1	5.5E-03	
2.0E-04	I	V		Hexachlorocyclopentadiene	77-47-4		2.1E-01
1.1E-05	C	3.0E-02	I	Hexachloroethane	67-72-1	2.6E-01	3.1E+01
				Hexachlorophene	70-30-4		
1.0E-05	I	V		Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4		1.0E-02
				Hexamethylene Diisocyanate, 1,6-	822-06-0		
				Hexamethylphosphoramide	680-31-9		
7.0E-01	I	V		Hexane, N-	110-54-3		7.3E+02
				Hexanedioic Acid	124-04-9		
3.0E-02	I	V		Hexanone, 2-	591-78-6		3.1E+01
4.9E-03	I	3.0E-05	P	Hexazinone	51235-04-2	5.7E-04	3.1E-02
4.9E-03	I			Hydrazine	302-01-2	5.7E-04	
				Hydrazine Sulfate	10034-93-2		

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Toxicity and Chemical-specific		Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1		
IUR (ug/m ³ -y)	k e y	RfC _i (mg/m ³ -y)	k e y c m u t a g e n	Analyte CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)	
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			
				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 NA		3.1E+02	
				Kerb 23950-58-5			
				Lactofen 77501-63-4			
				Lead Compounds			
1.5E-01	C	2.0E-04	C	M	~Lead Chromate 7758-97-6	6.8E-06	2.1E-01
1.2E-05	C				~Lead Phosphate 7446-27-7	2.3E-01	
8.0E-05	C				~Lead acetate 301-04-2	3.5E-02	
1.2E-05	C				~Lead and Compounds 7439-92-1		1.5E-01
					~Lead subacetate 1335-32-6	2.3E-01	
	V				~Tetraethyl Lead 78-00-2		
					Linuron 330-55-2		
					Lithium 7439-93-2		
					Londax 83055-99-6		
					MCPA 94-74-6		
					MCPB 94-81-5		
					MCPD 93-65-2		
7.0E-04	C				Malathion 121-75-5		7.3E-01
					Maleic Anhydride 108-31-6		
					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		
					Mephsolan 950-10-7		
					Mepiquat Chloride 24307-26-4		
					Mercury Compounds		
3.0E-04	S				~Mercuric Chloride (and other Mercury salts) 7487-94-7		3.1E-01
3.0E-04	I	Y			~Mercury (elemental) 7439-97-6		3.1E-01
					~Methyl Mercury 22967-92-6		
	Y				~Phenylmercuric Acetate 62-38-4		
					Merphos 150-50-5		
					Merphos Oxide 78-48-8		
3.0E-02	P	Y			Metalaxyl 57837-19-1		3.1E+01
					Methacrylonitrile 126-98-7		
					Methamidophos 10265-92-6		
2.0E+01	I	V			Mevanone 67-56-1		2.1E+04
					Methidathion 950-37-8		
					Methomyl 16752-77-5		
1.4E-05	C				Methoxy-5-nitroaniline, 2- 99-59-2	2.0E-01	
					Methoxychlor 72-43-5		
1.0E-03	P	V			Methoxyethanol Acetate, 2- 110-49-6		1.0E+00
2.0E-02	I	V			Methoxyethanol, 2- 109-86-4		2.1E+01
	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-03	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		
4.0E-02	H	V			Methyl Phosphonic Acid 993-13-5		4.2E+01
2.8E-05	C				Methyl Styrene (Mixed Isomers) 25013-15-4		
					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+01	3.1E+03
					Methyl-1,4-benzenediamine dihydrochloride, 2- 615-45-2		
					Methyl-5-Nitroaniline, 2- 99-55-8		
2.4E-03	C				Methyl-N-nitro-N-nitrosoguanidine, N- 70-25-7	1.2E-03	
3.7E-05	C				Methylaniline Hydrochloride, 2- 636-21-5	7.6E-02	
					Methylarsonic acid 124-58-3		

Toxicity and Chemical-specific						Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1
IUR (ug/m ³ -1)	k e y	RfC _i (mg/m ³)	k e y	v o l a t i l e	m u t a g e n	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncancer Hazard SL HI=1 (ug/m ³)
6.3E-03	C				M	Methylbenzene, 1,4-diamine monohydrochloride, 2- Methylbenzene-1,4-diamine sulfate, 2- Methylcholanthrene, 3-	74612-12-7 615-50-9 56-49-5	1.6E-04	
1.0E-08	I	6.0E-01	I	V	M	Methylene Chloride	75-09-2	1.0E+02	6.3E+02
4.3E-04	C				M	Methylene-bis(2-chloroaniline), 4,4'- Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-14-4 101-61-1	2.4E-03 2.2E-01	
4.6E-04	C	2.0E-02	C			Methylenedibenzeneamine, 4,4'- Methylenediphenyl Diisocyanate Methylstyrene, Alpha-	101-77-9 101-68-8 98-83-9	6.1E-03	2.1E+01 6.3E-01
				V		Metolachlor Metribuzin Mineral oils	51218-45-2 21087-64-9 8012-95-1		
5.1E-03	C			V		Mirex Molinate Molybdenum	2385-85-5 2212-67-1 7439-98-7	5.5E-04	
				V		Monochloramine Monomethylaniline N,N'-Diphenyl-1,4-benzenediamine	10599-90-3 100-61-8 74-31-7		
0.0E+00	C	1.0E-01	P	V		Naled Naphtha, High Flash Aromatic (HFAN) Naphthylamine, 2-	300-76-5 64742-95-6 91-59-8		1.0E+02
2.6E-04	C	1.4E-05	C			Napropamide	15299-99-7		
2.6E-04	C	1.4E-05	C			Nickel Acetate Nickel Carbonate	373-02-4 3333-67-3	1.1E-02 1.1E-02	1.5E-02 1.5E-02
2.6E-04	C	1.4E-05	C	V		Nickel Carbonyl	13463-39-3	1.1E-02	1.5E-02
2.6E-04	C	1.4E-05	C			Nickel Hydroxide	12054-48-7	1.1E-02	1.5E-02
2.6E-04	C	2.0E-05	C			Nickel Oxide	1313-99-1	1.1E-02	2.1E-02
2.4E-04	I	1.4E-05	C			Nickel Refinery Dust	NA	1.2E-02	1.5E-02
2.6E-04	C	9.0E-05	A			Nickel Soluble Salts	7440-02-0	1.1E-02	9.4E-02
4.8E-04	I	1.4E-05	C			Nickel Subsulfide	12035-72-2	5.8E-03	1.5E-02
2.6E-04	C	1.4E-05	C			Nickelocene Nitrate Nitrate + Nitrite (as N)	1271-28-9 14797-55-8 NA	1.1E-02	1.5E-02
		5.0E-05	X			Nitrite	14797-65-0		
		6.0E-03	P			Nitroaniline, 2- Nitroaniline, 4-	88-74-4 100-01-6		5.2E-02 6.3E+00
4.0E-05	I	9.0E-03	I	V		Nitrobenzene Nitrocellulose Nitrofurantoin	98-95-3 9004-70-0 67-20-9	7.0E-02	9.4E+00
3.7E-04	C					Nitrofurazone Nitroglycerin Nitroguanidine	59-87-0 55-63-0 556-88-7	7.6E-03	
8.8E-06	P	5.0E-03	P	V		Nitromethane	75-52-5	3.2E-01	5.2E+00
2.7E-03	H	2.0E-02	I	V		Nitropropane, 2-	79-46-9	1.0E-03	2.1E+01
7.7E-03	C				M	Nitroso-N-ethylurea, N-	759-73-9	1.3E-04	
3.4E-02	C				M	Nitroso-N-methylurea, N-	684-93-5	3.0E-05	
1.6E-03	I			V		Nitroso-di-N-butylamine, N-	924-16-3	1.8E-03	
2.0E-03	C					Nitroso-di-N-propylamine, N-	621-64-7	1.4E-03	
8.0E-04	C					Nitrosodiethanolamine, N-	1116-54-7	3.5E-03	
4.3E-02	I				M	Nitrosodiethylamine, N-	55-18-5	2.4E-05	
1.4E-02	I	4.0E-05	X	V	M	Nitrosodimethylamine, N-	62-75-9	7.2E-05	4.2E-02
2.6E-06	C					Nitrosodiphenylamine, N-	00-30-0	1.1E+00	
6.3E-03	C			V		Nitrosomethylethylamine, N-	10595-95-6	4.5E-04	
1.9E-03	C					Nitrosomorpholine [N-]	59-89-2	1.5E-03	
2.7E-03	C					Nitrosopiperidine [N-]	100-75-4	1.0E-03	
6.1E-04	I					Nitrosopyrrolidine, N- Nitrotoluene, m-	930-55-2 99-08-1	4.6E-03	
				V		Nitrotoluene, o- Nitrotoluene, p- Nonane, n-	88-72-2 99-99-0 111-84-2		2.1E+01
						Norflurazon Nustar Octabromodiphenyl Ether	27314-13-2 85509-19-9 32536-52-0		
						Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX) Octamethylpyrophosphoramide Oryzalin	2691-41-0 152-16-9 19044-88-3		
						Oxadiazon Oxamyl Paclitaxel	19666-30-9 23135-22-0 76738-62-0		
				V		Paraquat Dichloride Parathion Pebulate	1910-42-5 56-38-2 1114-71-2		
						Pendimethalin Pentabromodiphenyl Ether Pentabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-99)	40487-42-1 32534-81-9 60348-60-9		

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Toxicity and Chemical-specific		Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1		
IUR (ug/m ³ -y) ⁻¹	k e y	RfC _i (mg/m ³ -y)	k e y c mutagen	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
			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	
		1.0E+00	P V	Pentaerythritol tetranitrate (PETN)	78-11-5		1.0E+03
				Pentane, n-	109-66-0		
				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		
			V	Perfluorobutane Sulfonate	375-73-5		
6.3E-07	C			Permethrin	52645-53-1	4.5E+00	
				Phenacetin	62-44-2		
		2.0E-01	C	Phenmedipham	13684-63-4		2.1E+02
				Phenol	108-95-2		
				Phenothiazine	92-84-2		
				Phenylenediamine, m-	108-45-2		
				Phenylenediamine, o-	95-54-5		
				Phenylenediamine, p-	106-50-3		
		3.0E-04	I V	Phenylphenol, 2-	90-43-7		
				Phorate	298-02-2		
				Phosgene	75-44-5		3.1E-01
				Phosmet	732-11-6		
				Phosphates, Inorganic			
				~Aluminum metaphosphate	13776-88-0		
				~Ammonium polyphosphate	68333-79-9		
				~Calcium pyrophosphate	7790-76-3		
				~Diammonium phosphate	7783-28-0		
				~Dicalcium phosphate	7757-93-9		
				~Dimagnesium phosphate	7782-75-4		
				~Dipotassium phosphate	7758-11-4		
				~Disodium phosphate	7558-79-4		
				~Monoaluminum phosphate	13530-50-2		
				~Monoammonium phosphate	7722-76-1		
				~Monocalcium phosphate	7758-23-8		
				~Monomagnesium phosphate	7757-86-0		
				~Monopotassium phosphate	7778-77-0		
				~Monosodium phosphate	7558-80-7		
				~Polyphosphoric acid	8017-16-1		
				~Potassium tripolyphosphate	13845-36-8		
				~Sodium acid pyrophosphate	7758-16-9		
				~Sodium aluminum phosphate (acidic)	7785-88-8		
				~Sodium aluminum phosphate (anhydrous)	10279-59-1		
				~Sodium aluminum phosphate (tetrahydrate)	10305-76-7		
				~Sodium hexametaphosphate	10124-56-8		
				~Sodium polyphosphate	68915-31-1		
				~Sodium trimetaphosphate	7785-84-4		
				~Sodium tripolyphosphate	7758-29-4		
				~Tetrapotassium phosphate	7320-34-5		
				~Tetrasodium pyrophosphate	7722-88-5		
				~Trialuminum sodium tetra decahydrogenoctaorthophosphate (dihydrate)	15136-87-5		
				~Tricalcium phosphate	7758-87-4		
				~Trimagnesium phosphate	7757-87-1		
				~Tripotassium phosphate	7778-53-2		
				~Trisodium phosphate	7601-54-9		
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+00	
				~Butylphthalyl Butylglycolate	85-70-1		
				~Dibutyl Phthalate	84-74-2		
			V	~Diethyl Phthalate	84-66-2		
				~Dimethylterephthalate	120-61-6		
				~Octyl Phthalate, di-N-	117-84-0		
2.0E-02	C			~Phthalic Acid, P-	100-21-0		
				~Phthalic Anhydride	85-44-9		2.1E+01
				Picloram	1918-02-1		
				Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3		
				Pirimiphos, Methyl	29232-93-7		
8.6E-03	C			Polybrominated Biphenyls	59536-65-1	3.3E-04	
2.0E-05	S		V	Polychlorinated Biphenyls (PCBs)			
				~Aroclor 1016	12674-11-2	1.4E-01	

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Toxicity and Chemical-specific				Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y c m u t a g e n	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
5.7E-04	S		V	~Aroclor 1221	11104-28-2	4.9E-03	
5.7E-04	S		V	~Aroclor 1232	11141-16-5	4.9E-03	
5.7E-04	S		V	~Aroclor 1242	53469-21-9	4.9E-03	
5.7E-04	S		V	~Aroclor 1248	12672-29-6	4.9E-03	
5.7E-04	S		V	~Aroclor 1254	11097-69-1	4.9E-03	
5.7E-04	S		V	~Aroclor 1260	11096-82-5	4.9E-03	
			V	~Aroclor 5460	11126-42-4		
1.1E-03	E	1.3E-03	E V	~Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	2.5E-03	1.4E+00
1.1E-03	E	1.3E-03	E V	~Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167)	52663-72-6	2.5E-03	1.4E+00
1.1E-03	E	1.3E-03	E V	~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	2.5E-03	1.4E+00
1.1E-03	E	1.3E-03	E V	~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 156)	38380-08-4	2.5E-03	1.4E+00
1.1E+00	E	1.3E-06	E V	~Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	2.5E-06	1.4E-03
1.1E-03	E	1.3E-03	E V	~Pentachlorobiphenyl, 2',3,4,4',5- (PCB 123)	65510-44-3	2.5E-03	1.4E+00
1.1E-03	E	1.3E-03	E V	~Pentachlorobiphenyl, 2,3',4,4',5- (PCB 118)	31508-00-6	2.5E-03	1.4E+00
1.1E-03	E	1.3E-03	E V	~Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	2.5E-03	1.4E+00
1.1E-03	E	1.3E-03	E V	~Pentachlorobiphenyl, 2,3,4,4',5- (PCB 114)	74472-37-0	2.5E-03	1.4E+00
3.8E+00	E	4.0E-07	E V	~Pentachlorobiphenyl, 3,3',4,4',5- (PCB 126)	57465-28-8	7.4E-07	4.2E-04
5.7E-04	I		V	~Polychlorinated Biphenyls (high risk)	1336-36-3	4.9E-03	
1.0E-04	I		V	~Polychlorinated Biphenyls (low risk)	1336-36-3	2.8E-02	
2.0E-05	I		V	~Polychlorinated Biphenyls (lowest risk)	1336-36-3	1.4E-01	
3.8E-03	E	4.0E-04	E	~Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	7.4E-04	4.2E-01
1.1E-02	E	1.3E-04	E V	~Tetrachlorobiphenyl, 3,4,4',5- (PCB 81)	70362-50-4	2.5E-04	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		
1.1E-04	C		V M	~Benz[a]anthracene	56-55-3	9.2E-03	
1.1E-04	C			~Benzo[j]fluoranthene	205-82-3	2.6E-02	
1.1E-03	C		M	~Benzo[a]pyrene	50-32-8	9.2E-04	
1.1E-04	C		M	~Benzo[b]fluoranthene	205-99-2	9.2E-03	
1.1E-04	C		M	~Benzo[k]fluoranthene	207-08-9	9.2E-03	
			V	~Chloronaphthalene, Beta-	91-58-7		
1.1E-05	C		M	~Chrysene	218-01-9	9.2E-02	
1.2E-03	C		M	~Dibenzo[a,h]anthracene	53-70-3	8.4E-04	
1.1E-03	C			~Dibenzo(a,e)pyrene	192-85-4	2.6E-03	
7.1E-02	C		M	~Dimethylbenz[a]anthracene, 7,12-	57-97-6	1.4E-05	
			V	~Fluoranthene	206-44-0		
1.1E-04	C		M	~Fluorene	86-73-7		
			V	~Indeno[1,2,3-cd]pyrene	193-39-5	9.2E-03	
			V	~Methylnaphthalene, 1-	90-12-0		
3.4E-05	C	3.0E-03	I V	~Methylnaphthalene, 2-	91-57-6		
			V	~Naphthalene	91-20-3	8.3E-02	3.1E+00
1.1E-04	C		V	~Nitropyrene, 4-	57835-92-4	2.6E-02	
			V	~Pyrene	129-00-0		
			V	Potassium Perfluorobutane Sulfonate	29420-49-3		
			V	Prochloraz	67747-09-5		
			V	Profluralin	26399-36-0		
			V	Prometon	1610-18-0		
			V	Prometryn	7287-19-6		
			V	Propachlor	1918-16-7		
			V	Propanil	709-98-8		
			V	Propargite	2312-35-8		
			V	Propargyl Alcohol	107-19-7		
			V	Propazine	139-40-2		
			V	Propham	122-42-9		
8.0E-03	I	V		Propiconazole	60207-90-1		
			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
			V	Propylene Glycol	57-55-6		
2.7E-04	A		V	Propylene Glycol Dinitrate	6423-43-4		2.8E-01
2.0E+00	I	V		Propylene Glycol Monoethyl Ether	1569-02-4		2.1E+03
			V	Propylene Glycol Monomethyl Ether	107-98-2		
3.7E-06	I	3.0E-02	I V	Propylene Oxide	75-56-9	7.6E-01	3.1E+01
			V	Pursuit	81335-77-5		
			V	Pydrin	51630-58-1		
			V	Pyridine	110-86-1		
			V	Quinalphos	13593-03-8		
			V	Quinoline	91-22-5		
3.0E-02	A			Refractory Ceramic Fibers	NA		3.1E+01
			V	Resmethrin	10453-86-8		
			V	Ronnel	299-84-3		
6.3E-05	C		M	Rotenone	83-79-4	1.6E-02	
			V	Safrole	94-59-7		
			V	Savay	78587-05-0		

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Toxicity and Chemical-specific						Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1
IUR (ug/m ³ -y) ⁻¹	k e y	RfC _i (mg/m ³ -y)	k e y	v o l a t i l e	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
2.0E-02		C				Selenious Acid	7783-00-8		
2.0E-02		C				Selenium	7782-49-2		2.1E+01
						Selenium Sulfide	7446-34-6		2.1E+01
3.0E-03		C				Sethoxydim	74051-80-2		
						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		
1.5E-01	C	2.0E-04	C		M	Sodium Dichromate	10588-01-9	6.8E-06	2.1E-01
						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		
						Stirofos (Tetrachlorovinphos)	961-11-5		
1.5E-01	C	2.0E-04	C		M	Strontium Chromate	7789-06-2	6.8E-06	2.1E-01
						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	NA		
2.0E-03		X				Sulfolane	126-33-0		2.1E+00
1.0E-03		C	V			Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9		
1.0E-03		C				Sulfur Trioxide	7446-11-9		1.0E+00
						Sulfuric Acid	7664-93-9		1.0E+00
						Systhane	88671-89-0		
						TCMTB	21564-17-0		
						Tebuthiuron	34014-18-1		
						Temephos	3383-96-8		
						Terbacil	5902-51-2		
						Terbufos	13071-79-9		
						Terbutryn	886-50-0		
						Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1		
						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-02	
2.6E-07	I	4.0E-02	I	V		Tetrachloroethylene	127-18-4	1.1E+01	4.2E+01
						Tetrachlorophenol, 2,3,4,6-	58-90-2		
						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
						Tetryl (Trinitrophenylmethyl nitramine)	479-45-8		
						Thallium (I) Nitrate	10102-45-1		
						Thallium (Soluble Salts)	7440-28-0		
						Thallium Acetate	563-68-8		
						Thallium Carbonate	6533-73-9		
						Thallium Chloride	7791-12-0		
						Thallium Sulfate	7446-18-6		
						Thiobencarb	28249-77-6		
						Thiodiglycol	111-48-8		
						Thiofanox	39196-18-4		
						Thiophanate, Methyl	23564-05-8		
						Thiram	137-26-8		
1.0E-04	A	V				Tin	7440-31-5		1.0E-01
5.0E+00	I	V				Titanium Tetrachloride	7550-45-0		
						Toluene	108-88-3		5.2E+03
						Toluene-2,5-diamine	95-70-5		
						Toluidine, p-	106-49-0		
						Total Petroleum Hydrocarbons (Aliphatic High)	NA		
6.0E-01	P	V				Total Petroleum Hydrocarbons (Aliphatic Low)	NA		6.3E+02
1.0E-01	P	V				Total Petroleum Hydrocarbons (Aliphatic Medium)	NA		1.0E+02
						Total Petroleum Hydrocarbons (Aromatic High)	NA		
3.0E-02	P	V				Total Petroleum Hydrocarbons (Aromatic Low)	NA		3.1E+01
3.0E-03	P	V				Total Petroleum Hydrocarbons (Aromatic Medium)	NA		3.1E+00
3.2E-04	I					Toxaphene	8001-35-2	8.8E-03	
						Tralometrin	66841-25-6		
						Tri-n-butyltin	688-73-3		
						Triacetin	102-76-1		
						Triallate	2303-17-5		
						Triasulfuron	82097-50-5		
						Tribromobenzene, 1,2,4-	615-54-3		
						Tributyl Phosphate	126-73-8		
						Tributyltin Compounds	NA		
3.0E+01	H	V				Tributyltin Oxide	56-35-9		
						Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1		3.1E+04
						Trichloroacetic Acid	76-03-9		

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Toxicity and Chemical-specific				Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1		
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³) ³	k e y	v o l a t i l e	mutagen	CAS No.	Analyte	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
						33663-50-2	Trichloroaniline HCl, 2,4,6-		
						634-93-5	Trichloroaniline, 2,4,6-		
						87-61-6	Trichlorobenzene, 1,2,3-		
						120-82-1	Trichlorobenzene, 1,2,4-		2.1E+00
						71-55-6	Trichloroethane, 1,1,1-		5.2E+03
						79-00-5	Trichloroethane, 1,1,2-	1.8E-01	2.1E-01
						79-01-6	Trichloroethylene	4.8E-01	2.1E+00
						75-69-4	Trichlorofluoromethane		7.3E+02
						95-95-4	Trichlorophenol, 2,4,5-		
						88-06-2	Trichlorophenol, 2,4,6-	9.1E-01	
						93-76-5	Trichlorophenoxyacetic Acid, 2,4,5-		
						93-72-1	Trichlorophenoxypropionic acid, -2,4,5		
						598-77-6	Trichloropropane, 1,1,2-		
						96-18-4	Trichloropropane, 1,2,3-		3.1E-01
						96-19-5	Trichloropropene, 1,2,3-		3.1E-01
						1330-78-5	Tricresyl Phosphate (TCP)		
						58138-08-2	Tridiphane		
						121-44-8	Triethylamine		7.3E+00
						112-27-6	Triethylene Glycol		
						1682-09-8	Trifluralin		
						512-56-1	Trimethyl Phosphate		
						526-73-8	Trimethylbenzene, 1,2,3-		5.2E+00
						95-63-6	Trimethylbenzene, 1,2,4-		7.3E+00
						108-67-8	Trimethylbenzene, 1,3,5-		
						99-35-4	Trinitrobenzene, 1,3,5-		
						118-96-7	Trinitrotoluene, 2,4,6-		
						791-28-6	Triphenylphosphine Oxide		
						13674-87-8	Tris(1,3-Dichloro-2-propyl) Phosphate		
						13674-84-5	Tris(1-chloro-2-propyl)phosphate	4.3E-03	
						126-72-7	Tris(2,3-dibromopropyl)phosphate		
						115-96-8	Tris(2-chloroethyl)phosphate		
						78-42-2	Tris(2-ethylhexyl)phosphate		4.2E-02
						NA	Uranium (Soluble Salts)		
						51-79-6	Urethane	3.5E-03	
						1314-62-1	Vanadium Pentoxide	3.4E-04	7.3E-03
						7440-62-2	Vanadium and Compounds		1.0E-01
						1929-77-7	Vermolate		
						50471-44-8	Vinclozolin		2.1E+02
						108-05-4	Vinyl Acetate		
						593-60-2	Vinyl Bromide	8.8E-02	3.1E+00
						75-01-4	Vinyl Chloride	1.7E-01	1.0E+02
						81-81-2	Warfarin		
						106-42-3	Xylene, P-		1.0E+02
						108-38-3	Xylene, m-		1.0E+02
						95-47-6	Xylene, o-		1.0E+02
						1330-20-7	Xylenes		1.0E+02
						1314-84-7	Zinc Phosphide		
						7440-66-6	Zinc and Compounds		
						12122-67-7	Zineb		
						7440-67-7	Zirconium		

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Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHILD Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k _e (y ⁻¹)	IUR (ug/m ³ -y ⁻¹)	k _e (y ⁻¹)	RfD _c (mg/kg-day)	k _e (y ⁻¹)	RfC _c (mg/m ³)	k _e (y ⁻¹)	v _o	mutagen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)
1.8E-02	C	5.1E-06	C	1.5E-01	I					-1.5	1	1	Yes	ALAR	1596-84-5	4.3E+00	1.3E+04		4.3E+00	3.0E+03	1.0E+07		3.0E+03	
8.7E-03	I			4.0E-03	I					-0.85	1	1	Yes	Acephate	30560-19-1	9.0E+00	1.1E+04		8.9E+00	8.0E+01	1.1E+05		8.0E+01	
		2.2E-06	I							-0.34	1	1	Yes	Acetaldehyde	75-07-0			2.6E+00	2.6E+00			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	3.1E+01	A	V		-0.24	1	1	Yes	Acetone	67-64-1					1.8E+04	4.4E+06	6.4E+04	1.4E+04	
						2.0E-03	X	V		-0.03	1	1	Yes	Acetone Cyanohydrin	75-96-5							4.2E+00	4.2E+00	
						6.0E-02	I	V		-0.34	1	1	Yes	Acetonitrile	75-05-8								1.3E+02	1.3E+02
3.8E+00	C	1.3E-03	C	1.0E-01	I					1.58	1	1	Yes	Acetophenone	98-86-2					2.0E+03	4.6E+04		1.9E+03	
										3.12	1	1	Yes	Acetylaminofluorene, 2-	53-96-3	2.1E-02	6.4E-02		1.6E-02					
				5.0E-04	I	2.0E-05	I	V		-0.01	1	1	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	Yes	Acrylamide	79-06-1	5.0E-02	2.3E+01		5.0E-02	4.0E+01	2.1E+04		4.0E+01	
				5.0E-01	I	1.0E-03	I	V		0.35	1	1	Yes	Acrylic Acid	79-10-7					1.0E+04	1.1E+06	2.1E+00	2.1E+00	
5.4E-01	I	6.8E-05	I	4.0E-02	A	2.0E-03	I	V		0.25	1	1	Yes	Acrylonitrile	107-13-1	1.4E-01	1.4E+01	8.3E-02	5.2E-02	8.0E+02	8.8E+04	4.2E+00	4.1E+00	
						6.0E-03	P			-0.32	1	1	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+00	4.2E+00		1.0E+00	2.0E+02	6.9E+02		1.6E+02	2.0E+00
				1.0E-03	I					1.13	1	1	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	Yes	Aldicarb Sulfone	1646-88-4					2.0E+01	2.4E+04		2.0E+01	2.0E+00
										-0.78	1	1	Yes	Aldicarb sulfoxide	1646-87-3								2.0E+01	4.0E+00
1.7E+01	I	4.9E-03	I	3.0E-05	I					6.5	1	1	No	Aldrin	309-00-2	4.6E-03		1.1E-03	9.2E-04	6.0E-01			6.0E-01	
				2.5E-01	I					2.2	1	1	Yes	Allyl	74223-64-6					5.0E+03	2.4E+05		4.9E+03	
				5.0E-03	I	1.0E-04	X	V		0.17	1	1	Yes	Allyl Alcohol	107-18-6					1.0E+02	1.3E+04	2.1E-01	2.1E-01	
2.1E-02	C	6.0E-06	C	1.0E+00	P	5.0E-03	P			1.93	1	1	Yes	Allyl Chloride	107-05-1	3.7E+00	3.3E+01	9.4E-01	7.3E-01	2.0E+04	4.5E+06		2.0E+04	2.1E+00
				4.0E-04	I						1	1	Yes	Aluminum	7429-90-5					8.0E+00	1.8E+03		8.0E+00	
											1	1	Yes	Aluminum Hydroxide	20899-73-8					6.0E+00	5.1E+02		5.9E+00	
				3.0E-04	I					2.31	1	1	Yes	Amdb	67485-29-4					1.8E+02	9.7E+02		1.5E+02	
2.1E+01	C	6.0E-03	C	9.0E-03	I					2.98	1	1	Yes	Ametyln	834-12-8					3.7E-03	1.5E-02		3.0E-03	
										2.86	1	1	Yes	Amindiphenyl, 4-	92-67-1									
				8.0E-02	P					0.21	1	1	Yes	Aminophenol, m-	591-27-5					1.6E+03	2.8E+05		1.6E+03	
				2.0E-02	P					0.04	1	1	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	Amtraz	33089-61-1					5.0E+01	9.7E+00		8.2E+00	
						1.0E-01	I	V		0.23	1	1	Yes	Ammonia	7664-41-7					4.0E+03	9.1E+05		4.0E+03	
				2.0E-01	I						1	1	Yes	Ammonium Sulfamate	7773-06-0							6.3E+00	6.3E+00	
						3.0E-03	X	V		0.89	1	1	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	Yes	Aniline	62-53-3	1.4E+01	6.6E+02		1.3E+01	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+00	4.9E+00		1.4E+00	4.0E+01	1.1E+02		3.0E+01	
				4.0E-04	I					0.15	1	1	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	1	Yes	Antimony Pentoxide	1314-60-9					1.0E+01	3.4E+02		9.7E+00	
				9.0E-04	H					-7.28	0.15	1	No	Antimony Potassium Tartrate	11071-15-1					1.8E+01			1.8E+01	
				4.0E-04	H					0.15	1	1	Yes	Antimony Tetroxide	1332-81-6					8.0E+00	2.7E+02		7.8E+00	
						2.0E-04	I			0.15	1	1	Yes	Antimony Trioxide	1309-64-4									
2.5E-02	I	7.1E-06	I	1.3E-02	I					3.1	1	0.9	Yes	Apache	74115-24-5	3.1E+00	2.3E+00		1.3E+00	2.6E+02	2.1E+03		2.3E+02	
				5.0E-02	H					4.82	1	0.8	Yes	Aramite	1440-57-8					1.0E+03	8.2E+02		4.5E+02	
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C				1	1	Yes	Arsenic, Inorganic	7440-38-2	5.2E-02	9.3E+00		5.2E-02	6.0E+00	1.4E+03		6.0E+00	1.0E+01
				3.5E-06	C	5.0E-05	I				1	1	Yes	Arsine	7784-42-1					7.0E-02	1.6E+01		7.0E-02	
				9.0E-03	I					4.28	1	0.9	Yes	Assure	76578-14-8					1.8E+02	3.8E+02		1.2E+02	
				5.0E-02	I					-0.27	1	1	Yes	Asulam	3337-71-1					1.0E+03	8.0E+05		1.0E+03	
2.3E-01	C			3.5E-02	I					2.61	1	1	Yes	Atrazine	1912-24-9	3.4E-01	2.6E+00		3.0E-01	7.0E+02	6.2E+03		6.3E+02	3.0E+00
8.8E-01	C	2.5E-04	C							2.98	1	0.9	Yes	Auramine	492-80-8	8.9E-02	2.6E-01		6.6E-02					
1.1E-01	I	3.1E-05	I	4.0E-04	I					4.48	1	1	No	Avermectin B1	65195-55-3					8.0E+00			8.0E+00	
										3.82	1	1	Yes	Azobenzene	103-33-3	7.1E-01	7.0E-01	1.8E-01	1.2E-01	2.0E+04	6.8E+07		2.0E+04	
				1.0E+00	P	7.0E-06	P			-1.7	1	1	Yes	Azodicarbonamide	123-77-3									
				2.0E-01	I	5.0E-04	H			0.07	1	1	Yes	Barium	7440-39-3					4.0E+03	6.4E+04		3.8E+03	2.0E+03
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M		0.025	1	1	Yes	Barium Chromate	10294-40-3	5.0E-02	2.3E-01		4.1E-02	4.0E+02	2.3E+03		3.4E+02	
				4.0E-03	I					1.52	1	1	Yes	Baygon	114-26-1					8.0E+01	3.6E+03		7.8E+01	
				3.0E-02	I					2.77	1	1	Yes	Bayleton	43121-43-3					6.0E+02	6.9E+03		5.5E+02	
				2.5E-02	I					5.95	1	0.7	Yes	Baythroid	68359-37-5					5.0E+02	1.6E+02		1.2E+02	
				3.0E-01	I					5.29	1	0.8	Yes	Benefin	1861-40-1					6.0E+03	2.4E+03		1.7E+03	
				5.0E-02	I					2.12	1	1	Yes											

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Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHILD Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³ -y) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _i (mg/m ³ -y)	k _e y	v	mutagen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)
1.3E+01	I			4.0E+00	I					1.87	1	1	Yes	Benzoic Acid	65-85-0	6.0E-03	5.7E-03		2.9E-03	8.0E+04	1.2E+06		7.5E+04	
										3.9	1	1	Yes	Benzoic Acid	98-07-7									
1.7E-01	I	4.9E-05	C	1.0E-01	P	2.0E-03	P	1.0E-03	P	1.1	1	1	Yes	Benzyl Alcohol	100-51-6	4.6E-01	3.2E+00	1.1E-01	8.9E-02	2.0E+03	8.9E+04		2.0E+03	
		2.4E-03	I	2.0E-03	I	2.0E-05	I			2.3	1	1	Yes	Benzyl Chloride	100-44-7					4.0E+01	3.2E+02	2.1E+00	2.0E+00	
										0.007	1	Yes	Beryllium and compounds	7440-41-7					4.0E+01	6.4E+01		2.5E+01	4.0E+00	
				1.0E-04	I					0	1	1	Yes	Bidrin	141-66-2					2.0E+00	1.1E+03		2.0E+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					8.15	1	0	No	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	Yes	Biphenyl, 1,1'-	92-52-4	9.7E+00	6.3E+00		3.8E+00	1.0E+04	7.3E+03	8.3E-01	8.3E-01	
7.0E-02	H	1.0E-05	H	4.0E-02	I					2.48	1	1	Yes	Bis(2-chloro-1-methylethyl) ether	108-60-1	1.1E+00	7.9E+00	5.6E-01	3.6E-01	8.0E+02	6.5E+03		7.1E+02	
				3.0E-03	P					1.3	1	1	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							1.29	1	1	Yes	Bis(2-chloroethyl)ether	111-44-4	7.1E-02	2.6E+00	1.7E-02	1.4E-02					
2.2E+02	I	6.2E-02	I							0.57	1	1	Yes	Bis(chloromethyl)ether	542-88-1	3.5E-04	3.2E-02	9.1E-05	7.2E-05					
				5.0E-02	I					3.32	1	1	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	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	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	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	Yes	Bromate	15641-45-4	1.1E-01	2.0E+01		1.1E-01	8.0E+01	1.8E+04		8.0E+01	1.0E+01
2.0E+00	X	6.0E-04	X								1.92	1	1	Bromo-2-chloroethane, 1-	107-04-0	3.9E-02	5.5E-01	9.4E-03	7.4E-03					
				8.0E-03	I	6.0E-02	I			2.99	1	1	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	Yes	Bromochloromethane	74-97-5							8.3E+01	8.3E+01	
6.2E-02	I	3.7E-05	C	2.0E-02	I					2	1	1	Yes	Bromodichloromethane	75-27-4	1.3E+00	1.8E+01	1.5E-01	1.3E-01	4.0E+02	6.4E+03		3.8E+02	8.0E+01(F)
7.9E-03	I	1.1E-06	I	2.0E-02	I					2.4	1	1	Yes	Bromoforn	75-25-2	9.9E+00	1.4E+02	5.1E+00	3.3E+00	4.0E+02	6.2E+03		3.8E+02	8.0E+01(F)
				1.4E-03	I	5.0E-03	I	V		1.19	1	1	Yes	Bromomethane	74-83-9					2.8E+01	1.0E+03	1.0E+01	7.5E+00	
				5.0E-03	H					5.21	1	0.8	Yes	Bromophos	2104-96-3					1.0E+02	5.5E+01		3.5E+01	
				2.0E-02	I					3.39	1	0.9	Yes	Bromoxynil	1689-84-5					4.0E+02	1.8E+03		3.3E+02	
3.4E+00	C	3.0E-05	I	2.0E-02	I					5.4	1	0.8	Yes	Bromoxynil Octanoate	1689-89-2	2.3E-02	1.6E-01	1.9E-01	1.8E-02	4.0E+02	2.1E+02		1.4E+02	
				2.0E-03	I	V				1.99	1	1	Yes	Butadiene, 1,3-	106-99-0					2.0E+03	1.0E+05	4.2E+00	4.2E+00	
				1.0E-01	I					0.88	1	1	Yes	Butanol, N-	71-36-3					2.0E+03	1.0E+05		2.0E+03	
1.9E-03	P			2.0E-01	I					4.73	1	0.9	Yes	Butyl Benzyl Phosphate	85-88-7	4.1E+01	2.6E+01		1.6E+01	4.0E+03	2.9E+03		1.7E+03	
				2.0E+00	P	3.0E+01	P	V		0.61	1	1	Yes	Butyl alcohol, sec-	78-92-2					4.0E+04	3.0E+06	6.3E+04	2.4E+04	
				5.0E-02	I					4.15	1	1	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	1	Yes	Butylated hydroxyanisole	25013-16-5	3.9E+02	6.2E+02		2.4E+02					
3.6E-03	P			3.0E-01	P					5.1	1	1	Yes	Butylated hydroxytoluene	128-37-0	2.2E+01	3.8E+00		3.3E+00	6.0E+03	1.2E+03		1.0E+03	
				5.0E-02	P					4.38	1	1	No	Butylbenzene, n-	104-51-8					1.0E+03			1.0E+03	
				1.0E-01	X					4.57	1	1	No	Butylbenzene, sec-	135-98-8					2.0E+03			2.0E+03	
				1.0E-01	X					4.11	1	1	Yes	Butylbenzene, tert-	98-06-6					2.0E+03	1.1E+03		6.9E+02	
				2.0E-02	A					0.36	1	1	Yes	Cacodylic Acid	75-60-5					4.0E+02	6.7E+04		4.0E+02	
1.8E-03	I	1.0E-03	I	1.0E-05	A					0.025	1	1	Yes	Cadmium (Diet)	7440-43-9					1.0E+01	1.1E+02		9.2E+00	
1.8E-03	I	5.0E-04	I	1.0E-05	A					0.05	1	1	Yes	Cadmium (Water)	7440-43-9					4.0E+02	2.3E+03		3.4E+02	5.0E+00
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M		0.025	1	1	Yes	Calcium Chromate	13765-19-0	5.0E-02	2.3E-01		4.1E-02					
1.5E-01	C	4.3E-05	C	5.0E-01	I	2.2E-03	C			-0.19	1	1	Yes	Caprolactam	105-60-2					1.0E+04	9.0E+05		9.9E+03	
2.3E-03	C	6.6E-07	C	2.0E-03	I					3.8	1	0.9	Yes	Captadol	2425-06-1	5.2E-01	1.7E+00		4.0E-01	4.0E+01	1.5E+02		3.2E+01	
				1.3E-01	I					2.8	1	1	Yes	Captan	133-06-2	3.4E+01	3.4E+02		3.1E+01	2.6E+03	3.0E+04		2.4E+03	
				1.0E-01	I					2.36	1	1	Yes	Carbaryl	83-25-2					2.0E+03	2.4E+04		1.8E+03	
				5.0E-03	I					2.32	1	1	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	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	Yes	Carbon Tetrachloride	56-23-5	1.1E+00	4.2E+00	9.4E-01	4.5E-01	8.0E+01	3.4E+02	2.1E+02	4.9E+01	5.0E+00
				1.0E-02	I					3.81	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	Yes	Carboxin	5234-68-4					2.0E+03	4.1E+04		1.9E+03	
				9.0E-04	I						1	1	Yes	Ceric oxide	1306-38-3									
4.0E-01	H			1.0E-01	I					0.99	1	1	Yes	Chloral Hydrate	302-17-0					2.0E+03	1.5E+05		2.0E+03	
				1.5E-02	I					1.9	1	1	Yes	Chloramben	133-90-4					3.0E+02	7.4E+03		2.9E+02	
3.5E-01	I	1.0E-04	I	5.0E-04	I	7.0E-04	I	V		2.22	1	1	Yes	Chloranil	118-75-2	1.9E-01	3.4E+00		1.8E-01	1.0E+01		1.5E+00	1.3E+00	2.0E+00
1.0E+01	I	4.6E-03	C	3.0E-04	I					5.41	1	0.8	Yes	Chlordane	12789-03-6	2.2E-01		5.6E-02	4.5E-02	6.0E+00	5.4E+00		2.9E+00	
				7.0E-04	A					3.81	1	0.9	Yes	Chlordecone (Kepone)	143-50-0	7.8E-03	6.2E-03		3.5E-03				</	

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	ke	IUR (ug/m ³ -y) ⁻¹	ke	RfD _c (mg/kg-day)	ke	RfC _c (mg/m ³)	ke	v	muta-	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)
1.0E-01	P	7.7E-05	C	3.0E-03	X					2.27	1	1	Yes	Chloro-2-methylaniline, 4-	95-69-2	7.8E-01	6.4E+00		6.9E-01	6.0E+01	5.5E+02		5.4E+01	
2.7E-01	X			2.0E-03	H					0.09	1	1	Yes	Chloroacetaldehyde, 2-	107-20-0	2.9E-01	4.4E+01		2.9E-01	4.0E+01	6.3E+03		4.0E+01	6.0E+01
						3.0E-05	I			0.22	1	1	Yes	Chloroacetic Acid	79-11-8									
										1.93	1	1	Yes	Chloroacetophenone, 2-	532-27-4									
2.0E-01	P			4.0E-03	I					1.83	1	1	Yes	Chloroaniline, p-	106-47-8	3.9E-01	5.7E+00		3.6E-01	8.0E+01	1.3E+03		7.6E+01	
				2.0E-02	I	5.0E-02	P	V		2.84	1	1	Yes	Chlorobenzene	108-90-7					4.0E+02	1.3E+03	1.0E+02		7.8E+01
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	5.4E-01		3.1E-01	4.0E+02	3.5E+02		1.9E+02	1.0E+02
				3.0E-02	X					2.65	1	1	Yes	Chlorobenzoic Acid, p-	74-11-3					6.0E+02	3.4E+03			5.1E+02
				3.0E-03	P	3.0E-01	P	V		3.6	1	1	Yes	Chlorobenzotrifluoride, 4-	98-56-6					6.0E+01	9.3E+01	6.3E+02		3.5E+01
				4.0E-02	P		V			2.64	1	1	Yes	Chlorobutane, 1-	109-69-3					8.0E+02	3.0E+03			6.4E+02
				5.0E+01	I	V				1.08	1	1	Yes	Chlorodifluoromethane	75-45-6							1.0E+05		1.0E+05
3.1E-02	C	2.3E-05	I	1.0E-02	I	9.8E-02	A	V		0.03	1	1	Yes	Chloroethanol, 2-	107-07-3	2.5E+00	2.8E+01	2.4E-01	2.2E-01	4.0E+02	7.7E+04	2.0E+02		4.0E+02
				9.0E-02	I	V				1.97	1	1	Yes	Chloroform	67-66-3					2.0E+02	2.5E+03			9.7E+01
				9.0E-02	I	V				0.91	1	1	Yes	Chloromethane	74-87-3							1.9E+02		1.9E+02
2.4E+00	C	6.9E-04	C	3.0E-03	P	1.0E-05	X	V		0.32	1	1	Yes	Chloromethyl Methyl Ether	107-30-2	3.2E-02	3.5E+00	8.1E-03	6.5E-03	6.0E+01	6.4E+02			5.5E+01
3.0E-01	P			3.0E-03	P	1.0E-05	X	V		2.24	1	1	Yes	Chloronitrobenzene, o-	88-73-3	2.6E-01	2.5E+00		2.3E-01	6.0E+01	6.4E+02			5.5E+01
6.3E-03	P			1.0E-03	P	6.0E-04	P			2.39	1	1	Yes	Chloronitrobenzene, p-	100-00-5	1.2E+01	9.3E+01		1.1E+01	2.0E+01	1.7E+02			1.8E+01
				5.0E-03	I		V			2.15	1	1	Yes	Chlorophenol, 2-	95-57-8					1.0E+02	1.0E+03			9.1E+01
				4.0E-04	C	V				2.09	1	1	Yes	Chloropicrin	76-06-2							8.3E-01		8.3E-01
3.1E-03	C	8.9E-07	C	1.5E-02	I					3.05	1	0.9	Yes	Chlorothalonil	1897-45-6	2.5E+01	1.5E+02		2.2E+01	3.0E+02	2.1E+03			2.6E+02
				2.0E-02	I	V				3.42	1	1	Yes	Chlorotoluene, o-	95-49-8					4.0E+02	5.8E+02			2.4E+02
				2.0E-02	X	V				3.33	1	1	Yes	Chlorotoluene, p-	106-43-4					4.0E+02	6.6E+02			2.5E+02
2.4E+02	C	6.9E-02	C	2.0E-01	I					-1.02	1	1	Yes	Chlorozotocin	54749-90-5	3.2E-04	7.1E-01		3.2E-04	4.0E+03	9.8E+03			2.8E+03
				1.0E-03	A					3.51	1	0.9	Yes	Chlorpropham	101-21-3					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	I					2	1	1	Yes	Chlorsulfuron	64902-72-3					1.0E+03	5.7E+04			9.9E+02
				8.0E-04	H					5.8	1	0.8	Yes	Chlorthophos	60238-56-4					1.6E+01	3.4E+00			2.8E+00
5.0E-01	J	8.4E-02	S	1.5E+00	I	1.0E-04	I	M		0.013	1	1	Yes	Chromium(III), Insoluble Salts	16065-83-1	5.0E-02	1.1E-01		3.5E-02	3.0E+04	8.9E+04			2.2E+04
				3.0E-03	I					0.025	1	1	Yes	Chromium(VI)	18540-29-9					6.0E+01	1.7E+02			4.4E+01
				3.0E-03	I					0.013	1	1	Yes	Chromium, Total	7440-47-3					6.0E+01	1.7E+02			4.4E+01
				9.0E-03	P	3.0E-04	P	6.0E-06	P	1	1	1	Yes	Cobalt	7440-48-4					6.0E+00	3.4E+03			6.0E+00
				6.2E-04	I			V	M	1	0	0	Yes	Coke Oven Emissions	8007-45-2					8.0E+02	1.8E+05			8.0E+02
				4.0E-02	H					1	1	1	Yes	Copper	7440-50-8					8.0E+02	1.8E+05			8.0E+02
				5.0E-02	I	6.0E-01	C			1.96	1	1	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	Yes	Cresol, o-	95-48-7					1.0E+03	1.2E+04			9.3E+02
				1.0E-01	A	6.0E-01	C			1.94	1	1	Yes	Cresol, p-	106-44-5					2.0E+03	2.5E+04			1.9E+03
1.9E+00	H			1.0E-01	A					3.1	1	1	Yes	Cresol, ip-chloro-m-	99-50-7	4.1E-02	2.6E+00		4.0E-02	2.0E+03	5.2E+03			1.4E+03
				1.0E-01	A	6.0E-01	C			1.95	1	1	Yes	Cresols	1319-77-3					2.0E+03	2.4E+04			1.9E+03
				1.0E-03	P		V			0.6	1	1	Yes	Crotonaldehyde, trans-	123-73-9					2.0E+01	1.5E+03			2.0E+01
2.2E-01	C	6.3E-05	C	1.0E-01	I	4.0E-01	I	V		3.66	1	1	Yes	Cumene	98-82-8	3.5E-01	1.5E+00		3.5E-01	2.0E+03	1.9E+03	8.3E+02		4.5E+02
8.4E-01	H			2.0E-03	H					-3.16	1	1	No	Cupferron	135-20-6	9.3E-02			8.7E-02	4.0E+01	7.5E+02			3.8E+01
				2.0E-03	H					2.22	1	1	Yes	Cyanazine	21725-46-2					4.0E+01	7.5E+02			3.8E+01
				1.0E-03	I						1	1	Yes	Cyanides										
				5.0E-03	I						1	1	Yes	-Calcium Cyanide	592-01-8					2.0E+01	4.5E+03			2.0E+01
				6.0E-04	I	8.0E-04	S	V			1	1	Yes	-Copper Cyanide	544-92-3					1.0E+02	2.3E+04			1.0E+02
				1.0E-03	I		V			0.07	1	1	Yes	-Cyanide (CN)	57-12-5					1.2E+01	2.7E+03	1.7E+00		1.5E+00
				9.0E-02	I		V				1	1	Yes	-Cyanogen	460-19-5					2.0E+01	5.1E+03			2.0E+01
				5.0E-02	I		V				1	1	Yes	-Cyanogen Bromide	506-68-3					1.8E+03	1.6E+06			1.8E+03
				6.0E-04	I	8.0E-04	I	V		-0.25	1	1	Yes	-Cyanogen Chloride	506-77-4					1.0E+03	5.8E+05			1.0E+03
				2.0E-03	I		V				1	1	Yes	-Hydrogen Cyanide	74-90-8					1.2E+01	2.7E+03	1.7E+00		1.5E+00
				5.0E-03	I						1	1	Yes	-Potassium Cyanide	151-50-8					4.0E+01	4.5E+03			4.0E+01
				1.0E-01	I					0.04	1	1	Yes	-Potassium Silver Cyanide	506-61-6					1.0E+02	4.5E+02			8.2E+01
				1.0E-03	I						0.04	1	Yes	-Silver Cyanide	506-64-9					2.0E+03	1.8E+04			1.8E+03
				2.0E-04	P						1	0	Yes	-Sodium Cyanide	143-33-9					2.0E+01	4.5E+03			2.0E+01
				2.0E-04	P						1	0	Yes	-Thiocyanates	NA					4.0E+00	9.1E+02			4.0E+00
				2.0E-04	X		V			0.58	1	1	Yes	-Thiocyanic Acid	463-56-9					4.0E+00	9.1E+02			4.0E+00
				5.0E-02	I						1	1	Yes	-Zinc Cyanide	557-21-1					1.0E+03	3.8E+05			1.0E+03
2.3E-02	H			5.0E+00	I	7.0E-01	P	V		4.72	1	0.9	Yes	Cyclohexane	110-82-7	3.4E+								

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -y)	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³)	k _e (y)	v	muta-gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	
2.4E-01	I	6.9E-05	C	1.0E-02 7.5E-03	I					6.6 0.96 6.02	1 1 1	0.7 1 0.8	No Yes Yes	Cypermethrin Cyromazine DDD	52315-07-8 66215-27-8 72-54-8	3.2E-01	3.4E-02		3.1E-02	2.0E+02 1.5E+02			2.0E+02 1.5E+02		
3.4E-01	I	9.7E-05	C					V		6.51	1	0.8	No	DDE, p,p'-	72-55-9	2.3E-01		5.8E-02	4.6E-02						
3.4E-01	I	9.7E-05	I	5.0E-04 1.0E-02	I					6.91 4.28	1 1	0.7 0.9	No Yes	DDT Dacthal	50-29-3 1861-32-1	2.3E-01			2.3E-01	1.0E+01 2.0E+02			1.0E+01 1.2E+02		
7.0E-04	I			3.0E-02 7.0E-03 4.0E-05	I					0.78 12.11 3.21	1 1 1	1 0 0.9	Yes No No	Dalapon Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209) Demeton	75-99-0 1163-19-5 8065-48-3	1.1E+02			1.1E+02	6.0E+02 1.4E+02 8.0E-01	5.5E+04 4.1E+00		6.0E+02 1.4E+02 6.7E-01	2.0E+02	
1.2E-03	I			6.0E-01	I					8.12	1	0	No	Di(2-ethylhexyl)adipate	103-23-1	6.5E+01			6.5E+01	1.2E+04			1.2E+04	4.0E+02	
6.1E-02	H									4.49 3.81	1 1	0.9 0.9	Yes Yes	Diallate Diazinon	2303-16-4 333-41-5	1.3E+00	8.9E-01		5.2E-01	1.4E+01	3.9E+01		1.0E+01		
8.0E-01	P	6.0E-03	P	1.0E-02 2.0E-04 4.0E-04	X P X	2.0E-04	I	V	M	4.38 2.96 3.75	1 1 1	1 1 0.9	Yes Yes Yes	Dibenzothiophene Dibromo-3-chloropropane, 1,2- Dibromobenzene, 1,3-	132-65-0 96-12-8 108-36-1	3.1E-02	1.6E-01	3.4E-04	3.3E-04	2.0E+02 4.0E+00 8.0E+00	9.6E+01 2.4E+01 1.6E+01		4.2E-01	6.5E+01 3.7E-01 5.3E+00	2.0E-01
8.4E-02	I	2.7E-05	C	1.0E-02 2.0E-02	I					3.79 2.16	1 1	0.9 1	Yes Yes	Dibromobenzene, 1,4- Dibromochloromethane Dibromoethane, 1,2-	106-37-6 124-48-1 106-93-4	9.3E-01 3.9E-02	1.4E+01 6.9E-01	2.1E-01 9.4E-03	1.7E-01 7.5E-03	2.0E+02 4.0E+02 1.8E+02	3.7E+02 6.7E+03 3.6E+03	1.9E+01	3.8E+02 1.7E+01	8.0E+01(F) 5.0E-02	
2.0E+00	I	6.0E-04	I	1.0E-02 9.0E-03	H A	4.0E-03	X	V		1.7 2.21	1 1	1 0	Yes No Yes	Dibromomethane (Methylene Bromide) Dibutyltin Compounds Dicamba	74-95-3 NA 1918-00-9					2.0E+02 6.0E+00 6.0E+02	5.4E+03 1.0E+04	8.3E+00	8.0E+00 6.0E+00 5.7E+02		
4.2E-03	P									2.6	1	1	Yes	Dichloro-2-butene, 1,4-	764-41-0			1.3E-03	1.3E-03						
4.2E-03	P									2.6	1	1	Yes	Dichloro-2-butene, cis-1,4-	1476-11-5			1.3E-03	1.3E-03						
4.2E-03	P									2.6	1	1	Yes	Dichloro-2-butene, trans-1,4-	110-57-6			1.3E-03	1.3E-03						
5.0E-02	I			4.0E-03 9.0E-02	I					0.92 3.43	1 1	1 1	Yes Yes	Dichloroacetic Acid Dichlorobenzene, 1,2-	79-43-6 95-50-1	1.6E+00	9.2E+01		1.5E+00	8.0E+01 2.9E+03	5.4E+03 2.4E+01		4.2E+02	7.9E+01 3.0E+02	6.0E+01 6.0E+02
5.4E-03	C	1.1E-05	C	7.0E-02	A	8.0E-01	I	V		3.44	1	1	Yes	Dichlorobenzene, 1,4-	106-46-7	1.4E+01	2.0E+01	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	1.0E-02 9.0E-03 2.0E-01	I					3.51 4.44 2.16	1 1 1	1 0.9 1	Yes Yes Yes	Dichlorobenzidine, 3,3'- Dichlorobenzophenone, 4,4'- Dichlorodifluoromethane	91-94-1 90-98-2 75-71-9	1.7E-01	4.3E-01		1.2E-01	1.8E+02 4.0E+03	1.4E+02 3.8E+04	2.1E+02	7.8E+01 2.0E+02		
5.7E-03	C	1.6E-06	C	2.0E-01	P					1.79	1	1	Yes	Dichloroethane, 1,1-	78-34-3	1.4E+01	1.8E+02	3.5E+00	2.7E+00	4.0E+03	5.8E+04		3.8E+03		
9.1E-02	I	2.6E-05	I	6.0E-03 5.0E-02	X I	7.0E-03	P	V		1.48 2.13	1 1	1 1	Yes Yes	Dichloroethane, 1,2- Dichloroethylene, 1,1-	107-06-2 75-35-4	8.6E-01	1.8E+01	2.2E-01	1.7E-01	1.2E+02 1.0E+03	2.8E+03 8.5E+03	1.5E+01 4.2E+02	1.3E+01 2.8E+02	5.0E+00 7.0E+00	
				2.0E-03	I					1.86	1	1	Yes	Dichloroethylene, 1,2-cis-	156-59-2					4.0E+01	3.6E+02		3.6E+01	7.0E+01	
				2.0E-02	I					2.09	1	1	Yes	Dichloroethylene, 1,2-trans-	156-60-5					4.0E+02	3.6E+03		3.6E+02	1.0E+02	
				3.0E-03	I					3.06	1	1	Yes	Dichlorophenyl(2,4-	120-83-2					6.0E+01	1.9E+02		4.6E+01		
3.6E-02	C	1.0E-05	C	1.0E-02 8.0E-03	I					2.81 3.53	1 1	1 0.9	Yes Yes	Dichlorophenoxy Acetic Acid, 2,4- Dichlorophenoxybutyric Acid, 4-(2,4-	94-75-7 94-82-6	2.2E+00	2.3E+01	5.6E-01	4.4E-01	2.0E+02 1.6E+02	1.3E+03 4.8E+02		1.7E+02 1.2E+02	7.0E+01	
				9.0E-02	A	4.0E-03	I	V		1.98	1	1	Yes	Dichloropropane, 1,2-	78-87-5					1.8E+03	2.1E+04	8.3E+00	8.3E+00	5.0E+00	
1.0E-01	I	4.0E-06	I	2.0E-02 3.0E-03 3.0E-02	P I I	2.0E-02	I	V		2 0.78 2.04	1 1 1	1 1 1	Yes Yes Yes	Dichloropropane, 1,3- Dichloropropanol, 2,3- Dichloropropene, 1,3-	142-28-9 616-23-9 542-75-6	7.8E-01	7.5E+00	1.4E+00	4.7E-01	4.0E+02 6.0E+01 6.0E+02	4.6E+03 4.9E+03 6.5E+03	4.2E+01	3.7E+02 5.9E+01 3.9E+01		
2.9E-01	I	8.3E-05	C	5.0E-04	I					1.43	1	1	Yes	Dichlorvos	62-73-7	2.7E-01	1.3E+01		2.6E-01	1.0E+01	5.6E+02		9.9E+00		
1.6E+01	I	4.6E-03	I	8.0E-02 5.0E-05	P I	3.0E-04	X	V		3.51 5.4	1 1	0.8	Yes Yes	Dicyclopentadiene Dieldrin	77-73-6 60-57-1	4.9E-03	2.6E-03		1.7E-03	1.6E+03 1.0E+00	3.5E+03 6.1E-01	6.3E-01	6.3E-01 3.8E-01		
				3.0E-04	C					5.0E-03	1	0		Diesel Engine Exhaust	NA					4.0E+01	8.4E+04		4.0E+01		
				2.0E-03 3.0E-02	P P	2.0E-04	P			-1.43 0.56	1 1	1	Yes Yes	Diethanolamine Diethylene Glycol Monobutyl Ether	111-42-2 112-34-5					6.0E+02	8.4E+04		6.0E+02		
3.5E+02	C	1.0E-01	C	6.0E-02 1.0E-03	P P	3.0E-04	P			-0.54 0.05	1 1	1	Yes Yes	Diethylene Glycol Monoethyl Ether Diethylformamide	111-90-0 617-84-5	2.2E-04	6.3E-05		4.9E-05	1.2E+03 2.0E+01	7.8E+05 4.2E+03		1.2E+03 2.0E+01		
				8.0E-02	I					0.65	1	1	Yes	Difenzoat	43222-48-6					1.6E+03	7.3E+05		1.6E+03		
				2.0E-02	I					3.88	1	0.9	Yes	Diffubenzuron	35367-38-5					4.0E+02	1.0E+03		2.9E+02		
4.4E-02	C	1.3E-05	C	4.0E+01	I	V				0.75	1	1	Yes	Diffuroethane, 1,1-	75-37-6							8.3E+04	8.3E+04		
										3.38	1	1	Yes	Dihydroxafrole	94-58-6	1.8E+00	2.2E+00	4.3E-01	3.0E-01	1.6E+03	1.3E+05		1.5E+03	1.5E+03	
				7.0E-01	P	V				1.52	1	1	Yes	Diisopropyl Ether	108-20-3							1.5E+03	1.6E+03		
				8.0E-02	I					1.03	1	1	Yes	Diisopropyl Methylphosphonate	1445-75-6					1.6E+03	1.3E+05		1.5E+03		
1.6E+00	P			2.0E-02 2.0E-04	I					-0.17 0.78	1 1	1	Yes Yes	Dimethipin Dimethoate	55290-64-7 60-51-5					4.0E+02 4.0E+00	2.4E+05 6.4E+02		4.0E+02 4.0E+00		
1.7E-03	P			6.0E-02	P					-0.61	1	1	Yes	Dimethyl methylphosphonate	756-79-6	4.6E+01	2.7E+04		4.6E+01	1.2E+03	8.1E+05		1.2E+03		
4.6E+00	C	1.3E-03	C							4.58	1	1	Yes	Dimethylamino azobenzene [p-]	60-11-7	1.7E-02	6.9E-03		4.9E-03						
5.8E-01	H									-1.51	1	1	Yes	Dimethylamine HCl, 2,4-	21436-96-4	1.3E-01	4.0E+02		1.3E-01						
2.0E-01	P			2.0E-03	X					1.68	1	1	Yes	Dimethylamine, 2,4-	95-68-1	3.9E-01	6.8E+00		3.7E-01	4.0E+01	8.0E+02		3.8E+01		

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k _e (y ⁻¹)	IUR (ug/m ³) ⁻¹	k _e (y ⁻¹)	RfD _c (mg/kg-day)	k _e (y ⁻¹)	RfC _i (mg/m ³)	k _e (y ⁻¹)	v _o	muta-gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)
1.1E+01	P			2.0E-03	I			V		2.31	1	1	Yes	Dimethylaniline, N,N-Dimethylbenzidine, 3,3'-	121-69-7 119-93-7	7.1E-03	8.2E-02		6.5E-03	4.0E+01	3.0E+02		3.5E+01	
				1.0E-01	P	3.0E-02	I	V		-1.01	1	1	Yes	Dimethylformamide	68-12-2					2.0E+03	1.8E+06	6.3E+01	6.1E+01	
5.5E+02	C	1.6E-01	C		X	2.0E-06	X	V		-1.19	1	1	Yes	Dimethylhydrazine, 1,1-Dimethylhydrazine, 1,2-	57-14-7 540-73-8	1.4E-04	4.8E-02	3.5E-05	2.8E-05	2.0E+00	3.5E+03	4.2E-03	4.2E-03	
				2.0E-02	I					2.3	1	1	Yes	Dimethylphenol, 2,4-Dimethylphenol, 2,6-Dimethylphenol, 3,4-	105-67-9 576-26-1 95-65-8					4.0E+02	3.1E+03		3.6E+02	
4.5E-02	C	1.3E-05	C					V		2.58	1	1	Yes	Dimethylvinylchloride	513-37-1	1.7E+00	6.3E+00	4.3E-01	3.3E-01	1.2E+01	8.5E+01	1.7E+02	1.8E+01	
				8.0E-05	X					2.13	1	1	Yes	Dinitro-o-cresol, 4,6-Dinitro-o-cyclohexyl Phenol, 4,6-	534-52-1 131-89-5					1.6E+00	2.6E+01		1.5E+00	
				1.0E-04	P					1.69	1	1	Yes	Dinitrobenzene, 1,2-Dinitrobenzene, 1,3-Dinitrobenzene, 1,4-	528-29-0 99-65-0 100-25-4					2.0E+00	5.3E+01		1.9E+00	
6.8E-01	I									2.18	1	1	Yes	Dinitrophenol, 2,4-Dinitrotoluene Mixture, 2,4/2,6-Dinitrotoluene, 2,4-	NA 121-14-2	1.1E-01 2.5E-01	1.4E+00 4.1E+00		1.1E-01 2.4E-01	4.0E+01	7.5E+02		3.8E+01	
3.1E-01	C	8.9E-05	C			2.0E-03	I			1.98	1	1	Yes	Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 4-Amino-2,6-	606-20-2 35572-78-2 19406-51-0	5.2E-02	7.1E-01		4.8E-02	6.0E+00	9.3E+01		5.7E+00	
4.5E-01	X			9.0E-04	X					2.18	1	1	Yes	Dinitrotoluene, Technical grade	25321-14-6	1.7E-01	2.1E+00		1.6E-01	1.8E+01	2.5E+02		1.7E+01	
1.0E-01	I	5.0E-06	I			1.0E-03	I			3.56	1	0.9	Yes	Dinoseb	88-85-7	7.8E-01	2.2E+02	1.1E+00	4.6E-01	2.0E+01	5.4E+01		1.5E+01	7.0E+00
				3.0E-02	I	3.0E-02	I	V		-0.27	1	1	Yes	Dioxane, 1,4-Dioxins	123-91-1					6.0E+02	1.9E+05	6.3E+01	5.7E+01	
6.2E+03	I	1.3E+00	I							8.21	1	0	No	Hexachlorodibenzo-p-dioxin, Mixture	NA	1.3E-05			1.3E-05					
1.3E+05	C	3.8E+01	C			7.0E-10	I	4.0E-08	C	V		0.5	No	-TCDD, 2,3,7,8-	1746-01-6	6.0E-07		1.5E-07	1.2E-07	1.4E-05		8.3E-05	1.2E-05	3.0E-05
				3.0E-02	I					2.86	1	1	Yes	Diphenamid	957-51-7					6.0E+02	4.2E+03		5.3E+02	
				8.0E-04	X					2.4	1	1	Yes	Diphenyl Sulfone	127-63-9					1.6E+01	2.0E+02		1.5E+01	
				2.5E-02	I					3.5	1	1	Yes	Diphenylamine	122-39-4					5.0E+02	8.4E+02		3.1E+02	
8.0E-01	I	2.2E-04	I			2.2E-03	I			2.94	1	1	Yes	Diphenylhydrazine, 1,2-Diquat	122-667 85-00-7	9.7E-02	3.7E-01		7.7E-02	4.4E+01			4.4E+01	2.0E+01
7.1E+00	C	1.4E-01	C							4.9	1	1	No	Direct Black 38	1937-37-1	1.1E-02			1.1E-02					
7.4E+00	C	1.4E-01	C							-2.03	1	1	No	Direct Blue 6	2602-46-2	1.1E-02			1.1E-02					
6.7E+00	C	1.4E-01	C							-6.53	1	1	No	Direct Brown 95	16071-86-6	1.2E-02			1.2E-02					
				4.0E-05	I					4.02	1	0.9	Yes	Diquat	298-04-4					8.0E-01	1.3E+00		5.0E-01	
				1.0E-02	I			V		0.77	1	1	Yes	Dithiane, 1,4-Dithiane	505-29-3					2.0E+02	1.6E+04		2.0E+02	
				2.0E-03	I					2.68	1	1	Yes	Diuron	330-54-1					4.0E+01	3.6E+02		3.6E+01	
				4.0E-03	I					1.15	1	1	Yes	Dodine	2439-10-3					8.0E+01	1.1E+04		8.0E+01	
				2.5E-02	I			V		3.21	1	1	Yes	EPTC	759-94-4					5.0E+02	1.5E+03		3.8E+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	
				2.0E-02	I					1.91	1	1	Yes	Endothal	145-73-3					4.0E+02	8.5E+03		3.8E+02	1.0E+02
9.9E-03	I	1.2E-06	I			3.0E-04	I			5.2	1	0.8	Yes	Endrin	72-20-8				2.9E+00	6.0E+00	3.7E+00		2.3E+00	2.0E+00
				6.0E-03	P	1.0E-03	I	V		0.45	1	1	Yes	Epichlorohydrin	106-89-8	7.9E+00	7.5E+02	4.7E+00	2.9E+00	1.2E+02	1.3E+04	2.1E+00	2.0E+00	
				2.0E-02	I	V				0.86	1	1	Yes	Epoxybutane, 1,2-Ethephon	106-88-7 16672-87-0					2.0E+03	2.3E+05	1.3E+02	1.2E+02	
				5.0E-03	I					-0.22	1	1	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	Yes	Ethoxyethanol Acetate, 2-Ethoxyethanol, 2-Ethoxyethanol, 2-	111-15-9					2.0E+03	2.3E+05	1.3E+02	1.2E+02	
4.8E-02	H			9.0E-02	P	2.0E-01	I	V		-0.32	1	1	Yes	Ethoxyethanol, 2-Ethyl Acetate	110-80-5 141-78-6	1.6E+00	4.3E+01		1.6E+00	1.8E+04	1.2E+06	1.5E+02	1.4E+02	
				5.0E-03	P	8.0E-03	P	V		1.32	1	1	Yes	Ethyl Acrylate	140-88-5					1.0E+02	3.0E+03	1.7E+01	1.4E+01	
				1.0E+01	I	V				1.43	1	1	Yes	Ethyl Chloride (Chloroethane)	75-00-3					4.0E+03	2.0E+05		2.1E+04	2.1E+04
				2.0E-01	I	V				0.89	1	1	Yes	Ethyl Ether	60-29-7					1.8E+03	2.3E+04	6.3E+02	4.6E+02	
				9.0E-02	H	3.0E-01	P	V		1.94	1	1	Yes	Ethyl Methacrylate	97-63-2					1.8E+03	2.3E+04	6.3E+02	4.6E+02	
1.1E-02	C	2.5E-06	C			1.0E-05	I			4.78	1	0.8	Yes	Ethyl-p-nitrophenyl Phosphonate	2104-64-5	7.1E+00	1.2E+01	2.2E+00	1.5E+00	2.0E-01	1.6E-01		8.9E-02	
				1.0E-01	I	1.0E+00	I	V		3.15	1	1	Yes	Ethylbenzene	100-41-4					2.0E+03	3.8E+03	2.1E+03	8.1E+02	7.0E+02
				7.0E-02	P					-0.94	1	1	Yes	Ethylene Cyanohydrin	109-78-4					1.4E+03	1.1E+06		1.4E+03	
				9.0E-02	P			V		-2.04	1	1	No	Ethylene Diamine	107-15-3					1.8E+03			1.8E+03	
				2.0E+00	I	4.0E-01	C			-1.36	1	1	Yes	Ethylene Glycol	107-21-1					4.0E+04	5.7E+07		4.0E+04	
				1.0E-01	I	1.6E+00	I			0.83	1	1	Yes	Ethylene Glycol Monobutyl Ether	111-76-2					2.0E+03	1.4E+05		2.0E+03	
3.1E-01	C	8.8E-05	C			3.0E-02	C	V		-0.3	1	1	Yes	Ethylene Oxide	75-21-8	2.5E-01	5.2E+01	6.4E-02	5.1E-02				6.3E+01	6.3E+01
4.5E-02	C	1.3E-05	C			8.0E-05	I			-0.66	1	1	Yes	Ethylene Thiourea	96-45-7	1.7E+00	9.7E+02		1.7E+00	1.6E+00	1.0E+03		1.6E+00	
6.5E+01	C	1.9E-02	C					V		-0.28	1	1	Yes	Ethyleneimine	151-56-4	1.2E-03	2.4E-01	3.0E-04	2.4E-04					1.6E+00
				3.0E+00	I					2.19	1	1	Yes	Ethylphthalyl Ethyl Glycolate	84-72-0					6.0E+04	1.5E+06		5.8E+04	
				8.0E-03	I					2.55	1	1	Yes	Express	101200-48-0					1.6E+02	5.0E+03		1.6E+02	

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -y) ⁻¹	k _e (y)	RfD _c (mg/kg-day)	k _e (y)	RfC _i (mg/m ³)	k _e (y)	v	mutagen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)
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	
1.3E-02	I									2.42	1	1	Yes	Fluometuron	2164-17-2					2.6E+02	3.4E+03		2.4E+02	
4.0E-02	C	1.3E-02	C										Yes	Fluoride	16984-48-8					8.0E+02	1.8E+05		8.0E+02	
6.0E-02	I	1.3E-02	C								1	1	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	
2.0E-02	I									3.34	1	0.9	Yes	Flurprimidol	56425-91-3					4.0E+02	2.4E+03		3.4E+02	
6.0E-02	I									3.7	1	0.9	Yes	Flutolanil	66332-96-5					1.2E+03	4.5E+03		9.5E+02	
1.0E-02	I									6.81	1	0.6	No	Fluvalinate	69409-94-5					2.0E+02			2.0E+02	
3.5E-03	I	1.0E-01	I							2.85	1	1	Yes	Folpet	133-07-3	2.2E+01	2.0E+02		2.0E+01	2.0E+03	2.1E+04		1.8E+03	
1.9E-01	I									2.9	1	1	Yes	Fomesafen	72178-02-0	4.1E-01	8.7E+00		3.9E-01					
		2.0E-03	I							3.94	1	0.9	Yes	Fonofos	944-22-9					4.0E+01	6.3E+01		2.4E+01	
1.3E-05	I	2.0E-01	I	9.8E-03	A	V				0.35	1	1	Yes	Formaldehyde	50-00-0			4.3E-01	4.3E-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	Yes	Formic Acid	64-18-6					1.8E+04	6.3E+06	6.3E-01	6.3E-01	
3.0E+00	I									-2.4	1	1	No	Fosetyl-AL	39148-24-8					6.0E+04			6.0E+04	
														Furans										
1.0E-03	X		V							4.12	1	1	Yes	~Dibenzofuran	132-64-9					2.0E+01	1.3E+01		7.9E+00	
1.0E-03	I		V							1.34	1	1	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	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	Yes	Furazolidone	67-45-8	2.1E-02	9.8E+00		2.0E-02					
1.5E+00	C	4.3E-04	C							0.41	1	1	Yes	Furfural	98-01-1					6.0E+01	7.1E+03	1.0E+02	3.8E+01	
3.0E-02	I	8.6E-06	C							1.8	1	1	Yes	Furium	531-82-8	5.2E-02	1.8E+00		5.0E-02					
										4.38	1	0.9	Yes	Furmecycloz	60568-05-0	2.6E+00	1.9E+00		1.1E+00					
4.0E-04	I									-5.34	1	1	No	Glufosinate, Ammonium	77182-82-2					8.0E+00			8.0E+00	
										-0.18	1	1	Yes	Glutaraldehyde	111-30-8									
4.0E-04	I	1.0E-03	H	V						-0.12	1	1	Yes	Glycidyl	765-34-4					8.0E+00	1.8E+03	2.1E+00	1.7E+00	
1.0E-01	I									-3.4	1	1	No	Glyphosate, (Salt)	1071-83-6					2.0E+03			2.0E+03	7.0E+02
3.0E-03	I									4.73	1	0.8	Yes	Guanidine	42874-03-3					6.0E+01	6.6E+01		3.2E+01	
1.0E-02	X		V							-1.63	1	1	Yes	Guanidine	133-00-8					2.0E+02	4.2E+05		2.0E+02	
2.0E-02	P									-1.7	1	1	Yes	Guanidine Chloride	50-01-1					4.0E+02	1.0E+09		4.0E+02	
3.0E-03	A	1.0E-02	A							2.75	1	1	Yes	Guthrie	88-50-0					6.0E+01	8.3E+02		5.6E+01	
5.0E-05	I									4.07	1	0.9	Yes	Haloxyp, Methyl	69806-40-2					1.0E+00	3.1E+00		7.6E-01	
1.3E-02	I									1.56	1	1	Yes	Harmory	79277-27-3					2.6E+02	3.5E+04		2.6E+02	
4.5E+00	I	1.3E-03	I	5.0E-04	I	V				6.1	1	0.8	Yes	Heptachlor	76-44-8	1.7E-02	2.2E-03	4.3E-03	1.4E-03	1.0E+01	1.5E+00		1.3E+00	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	8.8E-03	6.8E-03	2.2E-03	1.4E-03	2.6E-01	2.4E-01		1.2E-01	2.0E-01
										6.07	1	0.7	No	Hexabromobenzene	87-82-1					4.0E+01			4.0E+01	
											1	0	No	Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BDE-163)	68631-49-2					4.0E+00			4.0E+00	
1.6E+00	I	4.6E-04	I	8.0E-04	I	V				5.73	1	0.9	No	Hexachlorobenzene	118-74-1	4.9E-02		1.2E-02	9.8E-03	1.6E+01			1.6E+01	1.0E+00
7.8E-02	I	2.2E-05	I	1.0E-03	P	V				4.78	1	0.9	Yes	Hexachlorobutadiene	67-68-3	1.0E+00	4.2E-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-02	1.7E-02		7.1E-03	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-02	5.9E-02		2.5E-02				3.6E+00	
1.1E+00	C	3.1E-04	C	3.0E-04	I					3.72	1	0.9	Yes	Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	7.1E-02	9.6E-02		4.1E-02	6.0E+00	9.2E+00			2.0E-01
1.8E+00	I	5.1E-04	I							4.14	1	0.9	Yes	Hexachlorocyclohexane, Technical	608-73-1	4.3E-02	5.9E-02		2.5E-02					
4.0E-02	I	1.1E-05	C	7.0E-04	I	3.0E-02	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
										4.14	1	1	Yes	Hexachloroethane	67-72-1	1.9E+00	1.7E+00	5.1E-01	3.3E-01	1.4E+01	1.4E+01	6.3E+01	6.2E+00	
										7.54	1	0	No	Hexachlorophene	70-30-4					6.0E+00			6.0E+00	
1.1E-01	I									0.87	1	1	Yes	Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	7.1E-01	8.3E+01		7.0E-01	6.0E+01	7.9E+03		6.0E+01	
										3.2	1	1	Yes	Hexamethylene Diisocyanate, 1,6-	822-06-0								2.1E-02	
										0.28	1	1	Yes	Hexamethylphosphoramide	680-31-9					8.0E+00	2.0E+03		8.0E+00	
6.0E-02	H	7.0E-01	I	V						3.9	1	1	Yes	Hexane, N-	110-54-3					1.2E+03	6.4E+02	1.5E+03	3.2E+02	
2.0E+00	P									0.08	1	1	Yes	Hexanedioic Acid	124-04-9					4.0E+04	1.1E+07		4.0E+04	
5.0E-03	I	3.0E-02	I	V						1.38	1	1	Yes	Hexanone, 2-	591-78-6					1.0E+02	2.7E+03	6.3E+01	3.8E+01	
3.0E+00	I	4.9E-03	I							1.85	1	1	Yes	Hexazinone	51235-04-2					6.6E+02	2.4E+04		6.4E+02	
3.0E+00	I	4.9E-03	I							-2.07	1	1	Yes	Hydrazine	302-01-2	2.6E-02	1.1E+02	1.1E-03	1.1E-03			6.3E-02	6.3E-02	
											1	1	Yes	Hydrazine Sulfate	10034-93-2	2.6E-02	4.7E+00		2.6E-02					
											1	1	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	Yes	Hydrogen Fluoride	7664-39-3					8.0E+02	1.8E+05		2.9E+01	2.8E+01
										0.23	1	1	Yes	Hydrogen Sulfide	7783-06-4							4.2E+00	4.2E+00	
6.0E-02																								

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³ -y) ⁻¹	k _e y	RfD _c (mg/kg-day)	k _e y	RfC _c (mg/m ³ -y)	k _v y	o	muta- gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)
9.5E-04	I			3.0E-01	I			V		0.76	1	1	Yes	Isobutyl Alcohol	78-83-1				7.8E+01	6.0E+03	3.6E+05		5.9E+03	
				2.0E-01	I	2.0E+00	C			1.7	1	1	Yes	Isophorone	78-59-1	8.2E+01	1.6E+03			4.0E+03	3.6E+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	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	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	No	JP-7	NA							6.3E+02	6.3E+02	
				7.5E-02	I					3.43	1	0.9	Yes	Kerb	23950-58-5					1.5E+03	5.5E+03		1.2E+03	
				2.0E-03	I					4.81	1	0.9	Yes	Lactofen	77501-63-4					4.0E+01	6.7E+01		2.5E+01	
														Lead Compounds										
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C		M	0.025	1	1	Yes	-Lead Chromate	7758-97-6	5.0E-02	2.3E-01		4.1E-02	4.0E+02	2.3E+03		3.4E+02	
8.5E-03	C	1.2E-05	C								1	0.8	Yes	-Lead Phosphate	7446-27-7	9.2E+00	1.6E+03		9.1E+00					
2.8E-01	C	8.0E-05	C							-0.08	1	1	Yes	-Lead acetate	301-04-2	2.8E-01	2.7E+02		2.8E-01					
8.5E-03	C	1.2E-05	C								1	1	Yes	-Lead and Compounds	7439-92-1								1.5E+01	1.5E+01
										-4	1	1	No	-Lead subacetate	1335-32-6	9.2E+00			9.2E+00					
				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	
				2.0E-03	I					3.2	1	0.9	Yes	Linuron	330-55-2					4.0E+01	2.0E+02		3.3E+01	
				2.0E-03	P						1	1	Yes	Lithium	7439-93-2					4.0E+01	9.1E+03		4.0E+01	
				2.0E-01	I					2.18	1	1	Yes	Londax	83055-99-6					4.0E+03	2.4E+05		3.9E+03	
				5.0E-04	I					3.25	1	1	Yes	MCPA	94-74-6					1.0E+01	3.0E+01		7.5E+00	
				1.0E-02	I					3.5	1	0.9	Yes	MCPB	94-81-5					2.0E+02	5.5E+02		1.5E+02	
				1.0E-03	I					3.13	1	1	Yes	MCPP	93-65-2					2.0E+01	7.1E+01		1.6E+01	
				2.0E-02	I					2.36	1	1	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	Yes	Maleic Anhydride	108-31-6					2.0E+03	3.8E+04		1.9E+03	
				5.0E-01	I					-0.84	1	1	Yes	Maleic Hydraside	123-33-1					1.0E+04	8.9E+06		1.0E+04	
				1.0E-04	P					-0.6	1	1	Yes	Malononitrile	109-77-3					2.0E+00	9.1E+02		2.0E+00	
				3.0E-02	H					1.33	1	0.9	Yes	Maprocebol	8018-01-7					6.0E+02	4.9E+03		5.4E+02	
				5.0E-03	I					0.62	1	1	Yes	Maneb	12427-38-2					1.0E+02	4.4E+03		9.8E+01	
				1.4E-01	I	5.0E-05	I				1	1	Yes	Manganese (Di-)	7439-96-5					4.8E+02	4.4E+03		4.3E+02	
				2.4E-02	S	5.0E-05	I			0.04	1	1	Yes	Manganese (Non-di-)	7439-96-5					4.8E+02	4.4E+03		4.3E+02	
				9.0E-05	H					1.04	1	1	Yes	Mephsolan	950-10-7					1.8E+00	2.5E+02		1.8E+00	
				3.0E-02	I					-2.82	1	1	No	Mepiquat Chloride	24307-26-4					6.0E+02			6.0E+02	
														Mercury Compounds										
				3.0E-04	I	3.0E-04	S			-0.22	0.07	1	Yes	-Mercuric Chloride (and other Mercury salts)	7487-94-7					6.0E+00	9.5E+01		5.7E+00	2.0E+00
						3.0E-04	I	V		0.62	1	1	Yes	-Mercury (elemental)	7439-97-6							6.3E-01	6.3E-01	2.0E+00
				1.0E-04	I						1	1	Yes	-Methyl Mercury	22967-92-6					2.0E+00	4.5E+02		2.0E+00	
				8.0E-05	I					0.71	1	1	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	
				3.0E-05	I					5.7	1	0.9	Yes	Merphos Oxide	78-48-8					6.0E-01	9.9E-02		8.5E-02	
				6.0E-02	I					1.65	1	1	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	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	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	Yes	Methanol	67-56-1					4.0E+04	1.8E+07	4.2E+04	2.0E+04	
				1.0E-03	I					2.2	1	1	Yes	Methidathion	950-37-8					2.0E+01	5.8E+02		1.9E+01	
				2.5E-02	I					0.6	1	1	Yes	Methomyl	16752-77-5					5.0E+02	6.8E+04		5.0E+02	
4.9E-02	C	1.4E-05	C							1.47	1	1	Yes	Methoxy-5-nitroaniline, 2-	99-59-2	1.6E+00	5.2E+01		1.5E+00					
				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	Yes	Methoxyethanol Acetate, 2-	110-49-6					1.6E+02	3.5E+04	2.1E+00	2.1E+00	
				5.0E-03	P	2.0E-02	I	V		-0.77	1	1	Yes	Methoxyethanol, 2-	109-86-4					1.0E+02	6.3E+04	4.2E+01	2.9E+01	
				1.0E+00	X			V		0.18	1	1	Yes	Methyl Acetate	79-20-9					2.0E+04	2.9E+06		2.0E+04	
				3.0E-02	H	2.0E-02	P	V		0.8	1	1	Yes	Methyl Acrylate	96-33-3					6.0E+02	3.7E+04	4.2E+01	3.9E+01	
				6.0E-01	I	5.0E+00	I	V		0.29	1	1	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	2.0E-05	X	V		-1.05	1	1	Yes	Methyl Hydrazine	60-34-4			5.6E-03	5.6E-03	2.0E+01	1.5E+04	4.2E-02	4.2E-02	
				8.0E-02	H	3.0E+00	I	V		1.31	1	1	Yes	Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1					1.6E+03	4.9E+04	6.3E+03	1.2E+03	
						1.0E-03	C	V		0.79	1	1	Yes	Methyl Isocyanate	624-83-9								2.1E+00	
				1.4E+00	I	7.0E-01	I	V		1.38	1	1	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	Yes	Methyl Parathion	298-00-0					5.0E+00	4.1E+01		4.5E+00	
				6.0E-02	X					-1	1	1	Yes	Methyl Phosphonic Acid	993-13-5					1.2E+03	1.2E+06		1.2E+03	
9.9E-02	C	2.8E-05	C			4.0E-02	H	V		3.44	1	1	Yes	Methyl Styrene (Mixed Isomers)	25013-15-4					1.2E+02	1.6E+02	8.3E+01	3.8E+01	
										-0.66	1	1	Yes	Methyl methanesulfonate	66-27-3	7.9E-01	4.6E+02		7.9E-01					
1.8E-03	C	2.6E-07	C			3.0E+00	I	V		0.94	1													

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Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k _e (y ⁻¹)	IUR (ug/m ³ -y) ⁻¹	k _e (y ⁻¹)	RfD _c (mg/kg-day)	k _e (y ⁻¹)	RfC _c (mg/m ³ -y)	k _e (y ⁻¹)	v	mutagen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)
1.3E-01	C	3.7E-05	C	1.0E-02	A					-2.06	1	1	Yes	Methylaniline Hydrochloride, 2-Methylarsonic acid	636-21-5 124-58-3	6.0E-01	3.7E+03		6.0E-01	2.0E+02	3.6E+05		2.0E+02	
1.0E-01	X	3.0E-04	X	2.0E-04	X					1	0	No	Methylbenzene, 1,4-diamine monohydrochloride, 2-Methylbenzene-1,4-diamine sulfate, 2-Methylcholanthrene, 3-	74612-12-7 615-50-9 56-49-5	7.8E-01 1.1E-03			7.8E-01	4.0E+00	6.0E+00		4.0E+00 6.0E+00		
2.0E-03	I	1.0E-08	I	6.0E-03	I	6.0E-01	V	M	1.25	1	1	Yes	Methylene Chloride	75-09-2	1.3E+01	3.4E+02	2.0E+02	1.1E+01	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	Yes	Methylene-bis(2-chloroaniline), 4,4'-Methylene-bis(N,N-dimethyl) Aniline, 4,4'-Methylolanthrene, 3-	101-14-4 101-61-1 56-49-5	2.5E-01 1.7E+00	4.2E-01 6.4E-01		1.6E-01 4.6E-01	4.0E+01 7.5E+01	2.6E+01		2.6E+01		
1.6E+00	C	4.6E-04	C	2.0E-02	C	6.0E-04	I	V	1.59	1	1	Yes	Methylenebisbenzenamine, 4,4'-Methylenediphenyl Diisocyanate	101-77-9 101-68-8 98-83-9	4.9E-02	1.6E+00		4.7E-02	1.4E+03	1.7E+03		7.8E+02		
				7.0E-02	H				5.22	1	0.9	Yes	Methylstyrene, Alpha-	101-68-8 98-83-9					1.4E+03	1.7E+03		7.8E+02		
				1.5E-01	I				3.13	1	1	Yes	Metolachlor	51218-45-2					3.0E+03	2.6E+04		2.7E+03		
				2.5E-02	I				1.7	1	1	Yes	Metribuzin	21087-64-9					5.0E+02	1.8E+04		4.9E+02		
				3.0E+00	P			V	6.1	1	1	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-03		1.1E-03	8.8E-04	4.0E+00				4.0E+00	
				2.0E-03	I				3.21	1	1	Yes	Molinate	2212-67-1					4.0E+01	1.2E+02		3.0E+01		
				5.0E-03	I				1	1	1	Yes	Molybdenum	7439-98-7					1.0E+02	2.3E+04		1.0E+02		
				1.0E-01	I				1	1	1	Yes	Monochloramine	10599-90-3					2.0E+03	4.5E+05		2.0E+03	4.0E+03	
				2.0E-03	P				1.66	1	1	Yes	Monomethylaniline	100-61-8					4.0E+01	7.5E+02		3.8E+01		
				3.0E-04	X				4.04	1	0.9	Yes	N,N-Diphenyl-1,4-benzenediamine	74-31-7					6.0E+00	8.8E+00		3.6E+00		
				2.0E-03	I			V	1.38	1	1	Yes	Naled	300-76-5					4.0E+01	6.8E+03		4.0E+01		
1.8E+00	C	0.0E+00	C	3.0E-02	X	1.0E-01	P	V	2.28	1	0	No	Naphtha, High Flash Aromatic (HFAN)	64742-95-6 91-59-8	4.3E-02	3.5E-01		3.9E-02	6.0E+02		2.1E+02	1.5E+02		
				1.0E-01	I				3.36	1	0.9	Yes	Naphthylamine, 2-	15299-99-7					2.0E+03	8.9E+03		1.6E+03		
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1	1	Yes	Napropamide	373-02-4					2.2E+02	6.8E+05		2.2E+02		
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1	1	Yes	Nickel Acetate	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	Yes	Nickel Carbonyl	13463-39-3			2.2E-02	2.2E-02		2.9E-02	2.9E-02		
				2.6E-04	C	1.1E-02	C	1.4E-05	C	0.04	1	Yes	Nickel Hydroxide	12954-48-7			2.2E-02	2.2E-02	2.2E+02	2.0E+03		2.0E+02		
				2.6E-04	C	1.1E-02	C	2.0E-05	C	0.04	1	Yes	Nickel Oxide	1313-99-1			2.2E-02	2.2E-02	2.2E+02	2.0E+03		2.0E+02		
				2.4E-04	I	1.1E-02	C	1.4E-05	C	0.04	0	Yes	Nickel Refinery Dust	NA					2.2E+02	1.0E+04		2.2E+02		
1.7E+00	C	4.8E-04	I	1.1E-02	C	1.4E-05	C		0.04	1	1	Yes	Nickel Soluble Salts	7440-02-0	4.0E-02	1.6E+00		4.5E-02	4.0E+02	1.8E+04		3.9E+02		
				2.6E-04	C	1.1E-02	C	1.4E-05	C	0.04	1	Yes	Nickel Sulfide	12035-72-2	4.0E-02	1.6E+00		4.5E-02	2.2E+02	1.0E+04		2.2E+02		
				1.6E+00	I				1	1	1	Yes	Nickelocene	1271-28-9					2.2E+02			2.2E+02		
									1	1	1	Yes	Nitrate	14797-55-8					3.2E+04	7.3E+06		3.2E+04	1.0E+04	
									1	0	Yes	Nitrate + Nitrite (as N)	NA										1.0E+04	
				1.0E-01	I				1	1	1	Yes	Nitrite	14797-85-0					2.0E+03	4.5E+05		2.0E+03	1.0E+03	
2.0E-02	P	4.0E-05	I	1.0E-02	X	5.0E-05	X		1.85	1	1	Yes	Nitroaniline, 2-	88-74-4	3.9E+00	1.2E+02		3.8E+00	2.0E+02	3.4E+03		1.9E+02		
				4.0E-03	P	6.0E-03	P		1.39	1	1	Yes	Nitroaniline, 4-	100-01-6					8.0E+01	2.8E+03		7.8E+01		
				2.0E-03	I	9.0E-03	I	V	1.85	1	1	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	No	Nitrocellulose	9004-70-0					6.0E+07			6.0E+07		
				7.0E-02	H				-0.47	1	1	Yes	Nitrofurantoin	67-20-9					1.4E+03	1.6E+06		1.4E+03		
1.3E+00	C	3.7E-04	C	1.0E-04	P				0.23	1	1	Yes	Nitrofurazone	59-87-0	6.0E-02	1.6E+01		6.0E-02	2.0E+00	8.7E+01		2.0E+00		
1.7E-02	P	1.0E-01	I	1.0E-04	P				1.62	1	1	Yes	Nitroglycerin	55-63-0	4.6E+00	1.8E+02		4.5E+00	2.0E+03	1.8E+06		2.0E+03		
				8.8E-06	P	5.0E-03	P	V	-0.35	1	1	Yes	Nitromethane	75-52-5			6.4E-01	6.4E-01				1.0E+01	1.0E+01	
				2.7E-03	H	2.0E-02	I	V	0.93	1	1	Yes	Nitropropane, 2-	79-46-9			2.1E-03	2.1E-03				4.2E+01	4.2E+01	
2.7E+01	C	7.7E-03	C	1.0E-01	I				0.23	1	1	Yes	Nitroso-N-ethylurea, N-	759-73-9	9.3E-04	1.5E-01		9.2E-04						
1.2E+02	C	3.4E-02	C	1.0E-01	I				-0.03	1	1	Yes	Nitroso-N-methylurea, N-	684-93-5	2.1E-04	4.5E-02		2.1E-04						
5.4E+00	I	1.6E-03	I	1.0E-01	I			V	2.63	1	1	Yes	Nitroso-di-N-butylamine, N-	924-16-3	1.4E-02	7.6E-02	3.5E-03	2.7E-03						
7.0E+00	I	2.0E-03	C	1.0E-01	I				1.36	1	1	Yes	Nitroso-di-N-propylamine, N-	621-64-7	1.1E-02	3.4E-01		1.1E-02						
2.8E+00	I	8.0E-04	C	1.0E-01	I				-1.28	1	1	Yes	Nitrosodiethanolamine, N-	1116-64-7	2.8E-02	7.8E+01		2.8E-02						
1.5E+02	I	4.3E-02	I	1.0E-01	I				0.48	1	1	Yes	Nitrosodiethylamine, N-	55-18-5	1.7E-04	1.6E-02		1.7E-04						
5.1E+01	I	1.4E-02	I	8.0E-06	P	4.0E-05	X	V	-0.57	1	1	Yes	Nitrosodimethylamine, N-	62-75-9	4.9E-04	1.9E-01	1.5E-04	1.1E-04	1.6E-01	7.4E+01	8.3E-02	5.5E-02		
4.9E-03	I	2.6E-06	C	1.0E-01	I				3.13	1	1	Yes	Nitrosodiphenylamine, N-	86-30-6	1.6E+01	5.0E+01		1.2E+01						
2.2E+01	I	6.3E-03	C	1.0E-01	I			V	0.04	1	1	Yes	Nitrosomethylethylamine, N-	10595-95-6	3.5E-03	6.2E-01	8.9E-04	7.1E-04						
6.7E+00	C	1.9E-03	C	1.0E-01	I				-0.44	1	1	Yes	Nitrosomorpholine [N-]	59-89-2	1.2E-02	5.1E+00		1.2E-02						
9.4E+00	C	2.7E-03	C	1.0E-01	I				0.36	1	1	Yes	Nitrosopiperidine [N-]	100-75-4	8.3E-03	1.0E+00		8.2E-03						
2.1E+00	I	6.1E-04	I	1.0E-04	X				-0.19	1	1	Yes	Nitrosopyrrolidine, N-	930-55-2	3.7E-02	9.9E+00		3.7E-02	2.0E+00	1.4E+01		1.7E+00		
				1.0E-04	X				2.45	1	1	Yes	Nitrotoluene, m-	99-08-1										
2.2E-01	P	9.0E-04	P	1.0E-04	P			V	2.3	1	1	Yes	Nitrotoluene, o-	88-72-2	3.5E-01	2.7E+00		3.1E-01	1.8E+01	1.5E+02		1.6E+01		
1.6E-02	P	4.0E-03	P	1.0E-04	P				2.37	1	1	Yes	Nitrotoluene, p-	99-99-0	4.9E+00	3.3E+01		4.2E+00	8.0E+01	6.2E+02		7.1E+01		
				3.0E-04	X	2.0E-02	P	V	5.65	1	1	No	Nonane, n-	111-84-2					6.0E+00		4.2E+01	5.3E+00		

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Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³) ⁻¹	k _e y	RfD _c (mg/kg-day)	k _e y	RfC _i (mg/m ³)	k _e y	v	muta- gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)		
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	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	Yes	Octamethylpyrophosphoramide	152-16-9					4.0E+01	1.4E+05		4.0E+01			
5.0E-02	I									3.73	1	0.9	Yes	Oryzalin	19044-88-3					1.0E+03	4.1E+03		8.1E+02			
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	Yes	Oxaryl	23135-22-0					5.0E+02	5.0E+05		5.0E+02	2.0E+02		
1.3E-02	I									3.2	1	0.9	Yes	Paclitaxel	76738-62-0					2.6E+02	1.7E+03		2.3E+02			
4.5E-03	I									-4.5	1	1	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	Yes	Pebutate	1114-71-2					1.0E+03	1.3E+03		5.6E+02			
4.0E-02	I									5.18	1	0.9	Yes	Pendimethalin	40487-42-1					8.0E+02	2.3E+02		1.8E+02			
2.0E-03	I									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			
9.0E-02	P									8.0E-04	I	V	5.17	1	0.9	Yes	Pentachlorobenzene	608-93-5				1.6E+01	3.9E+00		3.2E+00	
2.6E-01	H									3.0E-03	I	V	3.22	1	1	Yes	Pentachloroethane	76-01-7	8.7E-01	2.4E+00		6.4E-01	6.0E+01	4.4E+01	2.5E+01	
4.0E-01	I	5.1E-06	C	5.0E-03	I					4.64	1	0.9	Yes	Pentachloronitrobenzene	82-68-8	3.0E-01	1.9E-01		1.2E-01	6.0E+01	4.4E+01		2.5E+01			
4.0E-03	X			2.0E-03	P					5.12	1	0.9	Yes	Pentachlorophenol	87-86-5	1.9E-01	5.0E-02		4.0E-02	1.0E+02	2.9E+01		2.3E+01	1.0E+00		
				1.0E+00	P V					2.38	1	1	Yes	Pentaerythritol tetranitrate (PETN)	78-11-5	1.9E+01	4.1E+02		1.9E+01	4.0E+01	9.6E+02		3.9E+01			
										3.39	1	1	Yes	Pentane, n-	109-66-0							2.1E+03	2.1E+03			
														Perchlorates												
7.0E-04	I										1	1	Yes	-Ammonium Perchlorate	7790-98-9					1.4E+01	3.2E+03		1.4E+01			
7.0E-04	I										1	1	Yes	-Lithium Perchlorate	7791-03-9					1.4E+01	3.2E+03		1.4E+01			
7.0E-04	I										1	1	Yes	-Perchlorate and Perchlorate Salts	14797-73-0					1.4E+01	3.2E+03		1.4E+01	1.5E+01(F)		
7.0E-04	I										1	1	Yes	-Potassium Perchlorate	7778-74-7					1.4E+01	1.6E+03		1.4E+01			
7.0E-04	I										1	1	Yes	-Sodium Perchlorate	7601-89-0					1.4E+01	3.2E+03		1.4E+01			
2.0E-02	P									1.8173	1	1	Yes	Perfluorobutane Sulfonate	375-73-5					4.0E+02	8.3E+03		3.8E+02			
2.2E-03	C	6.3E-07	C	5.0E-02	I					6.5	1	0.6	No	Permethrin	52645-53-1	3.5E+01	1.1E+03		3.4E+01	1.0E+03			1.0E+03			
										1.58	1	1	Yes	Phenacetin	62-44-2								1.0E+03			
2.5E-01	I									3.59	1	0.9	Yes	Phenacetin	13684-63-4					5.0E+03	1.9E+04		4.0E+03			
3.0E-01	I	2.0E-01	C							1.46	1	1	Yes	Phenol	108-95-2					6.0E+03	1.4E+05		5.8E+03			
5.0E-04	X									4.15	1	1	Yes	Phenothiazine	92-84-2					1.0E+01	7.5E+00		4.3E+00			
4.7E-02	H									6.0E-03	I		Yes	Phenylenediamine, m-	108-45-2					1.2E+02	4.8E+04		1.2E+02			
										0.15	1	1	Yes	Phenylenediamine, o-	95-54-5	1.7E+00	2.8E+02		1.6E+00				1.2E+02			
										-0.3	1	1	Yes	Phenylenediamine, p-	106-50-3					3.8E+03	1.4E+06		3.8E+03			
1.9E-03	H									3.09	1	1	Yes	Phenol, 2-	90-43-7	4.0E+01	1.1E+02		3.0E+01	4.0E+00	1.2E+01		3.0E+00			
										3.56	1	0.9	Yes	Phorate	298-02-2									3.0E+00		
										-0.71	1	1	Yes	Phosgene	75-44-5											
2.0E-02	I									2.78	1	1	Yes	Phosnet	32-11-8					4.0E+02	5.3E+03		3.7E+02			
4.9E+01	P												Yes	Phosphates, Inorganic												
													Yes	-Aluminum metaphosphate	13176-88-0					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Ammonium polyphosphate	68333-79-9					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Calcium pyrophosphate	7790-76-3					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Diammonium phosphate	7783-28-0					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Dicalcium phosphate	7757-93-9					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Dimagnesium phosphate	7782-75-4					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Dipotassium phosphate	7758-11-4					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Disodium phosphate	7558-79-4					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Monoaluminum phosphate	13530-50-2					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Monoammonium phosphate	7722-76-1					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Monocalcium phosphate	7758-23-8					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Monomagnesium phosphate	7757-86-0					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Monopotassium phosphate	7778-77-0					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Monosodium phosphate	7558-80-7					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Polyphosphoric acid	8017-16-1					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Potassium triphosphate	13845-36-8					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Sodium acid pyrophosphate	7758-16-9					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Sodium aluminum phosphate (acidic)	7785-88-8					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Sodium aluminum phosphate (anhydrous)	10279-59-1					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Sodium aluminum phosphate (tetrahydrate)	10305-76-7					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Sodium hexametaphosphate	10124-56-8					9.7E+05	2.2E+08		9.7E+05			
4.9E+01	P												Yes	-Sodium polyphosphate	68915-31-1					9.7E+05	2.2E+08		9.7E+05			

Toxicity and Chemical-specific Information															Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncarcinogenic CHLD Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³ -y) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _o (mg/m ³ -y)	k _v y	muta- gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	
				4.9E+01	P					1	1	Yes	~Tetrasodium pyrophosphate	7722-88-5					9.7E+05	2.2E+08		9.7E+05		
				4.9E+01	P					1	0.8	Yes	~Trialuminum sodium tetra decahydrogenoctaorthophosphate (dihydrate)	15136-87-5					9.7E+05	2.2E+08		9.7E+05		
				4.9E+01	P					1	1	Yes	~Tricalcium phosphate	7758-87-4					9.7E+05	2.2E+08		9.7E+05		
				4.9E+01	P					1	1	Yes	~Trimagnesium phosphate	7757-87-1					9.7E+05	2.2E+08		9.7E+05		
				4.9E+01	P					1	1	Yes	~Tripotassium phosphate	7778-63-2					9.7E+05	2.2E+08		9.7E+05		
				4.9E+01	P					1	1	Yes	~Trisodium phosphate	7601-54-9					9.7E+05	2.2E+08		9.7E+05		
				3.0E-04	I	3.0E-04	I	V	-0.27	1	1	Yes	Phosphine	7803-51-2					6.0E+00	1.4E+03	6.3E-01	5.7E-01		
				4.9E+01	P	1.0E-02	I			1	1	Yes	Phosphoric Acid	7664-38-2					9.7E+05	2.2E+08		9.7E+05		
				2.0E-05	I		V		3.08	1	1	Yes	Phosphorus, White	7723-14-0					4.0E-01	9.1E+01		4.0E-01		
1.4E-02	I	2.4E-06	C	2.0E-02	I	1.0E+00	I		7.6	1	0.8	No	Phthalates		5.6E+00		5.6E+00		4.0E+02			4.0E+02	6.0E+00	
									4.15	1	0.9	Yes	~Bis(2-ethylhexyl)phthalate	117-81-7					2.0E+04	4.1E+04		1.3E+04		
												Yes	~Butylphthalyl Butylglycolate	85-70-1					2.0E+03	1.6E+03		9.0E+02		
									4.5	1	0.9	Yes	~Diethyl Phthalate	84-74-2					1.6E+04	2.0E+05		1.5E+04		
									2.42	1	1	Yes	~Dimethylterephthalate	120-61-6					2.0E+03	2.7E+04		1.9E+03		
									2.25	1	1	Yes	~Octyl Phthalate, di-N-	117-84-0					2.0E+02			2.0E+02		
									8.1	1	0	No	~Phthalic Acid, P-	100-21-0					2.0E+04	3.3E+05		1.9E+04		
									2	1	1	Yes	~Phthalic Anhydride	85-44-9					4.0E+04	1.1E+06		3.9E+04		
									1.6	1	1	Yes												
									1.9	1	1	Yes	Picloram	1918-02-1					1.4E+03	4.3E+04		1.4E+03	5.0E+02	
									0.93	1	1	Yes	Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3					2.0E+00	2.1E+02		2.0E+00		
									4.2	1	0.9	Yes	Pirimiphos, Methyl	29232-93-7					2.0E+02	3.1E+02		1.2E+02		
3.0E+01	C	8.6E-03	C	7.0E-06	H					1	0	No	Polybrominated Biphenyls	59538-65-1	2.6E-03			2.6E-03	1.4E-01			1.4E-01		
												No	Polychlorinated Biphenyls (PCBs)											
									5.69	1	0.9	No	~Aroclor 1016	12674-11-2	1.1E+00		2.8E-01	2.2E-01	1.4E+00			1.4E+00		
2.0E+00	S	5.7E-04	S						4.65	1	1	Yes	~Aroclor 1221	11104-28-2	3.9E-02	1.1E-02	9.8E-03	4.6E-03						
2.0E+00	S	5.7E-04	S						4.4	1	1	Yes	~Aroclor 1232	11441-16-5	3.9E-02	1.1E-02	9.8E-03	4.6E-03						
2.0E+00	S	5.7E-04	S						6.34	1	0.7	No	~Aroclor 1242	10489-21-9	3.9E-02		9.8E-03	7.8E-03						
2.0E+00	S	5.7E-04	S						6.2	1	0.7	No	~Aroclor 1248	12672-29-6	3.9E-02		9.8E-03	7.8E-03						
2.0E+00	S	5.7E-04	S	2.0E-05	I				6.5	1	0.5	No	~Aroclor 1254	11097-89-1	3.9E-02		9.8E-03	7.8E-03	4.0E-01			4.0E-01		
2.0E+00	S	5.7E-04	S						7.55	1	0	No	~Aroclor 1260	11096-82-5	3.9E-02		9.8E-03	7.8E-03						
3.9E+00	E	1.1E-03	E	6.0E-04	X				6.34	1	0.7	No	~Aroclor 5460	11126-42-4					1.2E+01			1.2E+01		
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	8.27	1	0	No	~Heptachlorobiphenyl, 2,3,3',4,4',5,5'-(PCB 189)	39635-31-9	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.5	1	0	No	~Hexachlorobiphenyl, 2,3',4,4',5,5'-(PCB 167)	52963-72-6	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.6	1	0	No	~Hexachlorobiphenyl, 2,3,3',4,4',5,5'-(PCB 157)	69782-90-7	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.6	1	0	No	~Hexachlorobiphenyl, 2,3,3',4,4',5,5'-(PCB 156)	39380-08-4	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		
3.9E+03	E	1.1E+00	E	2.3E-08	E	1.3E-06	E	V	7.41	1	0.1	No	~Hexachlorobiphenyl, 3,3',4,4',5,5'-(PCB 169)	32774-16-6	2.0E-05		4.9E-06	4.0E-06	4.7E-04		2.8E-03	4.0E-04		
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	6.98	1	0.4	No	~Pentachlorobiphenyl, 2,3,4,4',5-(PCB 123)	85510-44-3	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.12	1	0.3	No	~Pentachlorobiphenyl, 2,3,3',4,4',5-(PCB 118)	31508-00-6	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	6.79	1	0.5	No	~Pentachlorobiphenyl, 2,3,3',4,4'-(PCB 105)	32598-14-4	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	6.98	1	0.4	No	~Pentachlorobiphenyl, 2,3,4,4',5-(PCB 114)	74472-37-0	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		
1.3E+04	E	3.8E+00	E	7.0E-09	E	4.0E-07	E	V	6.98	1	0.4	No	~Pentachlorobiphenyl, 3,3',4,4',5-(PCB 126)	57465-28-8	6.0E-06		1.5E-06	1.2E-06	1.4E-04		8.3E-04	1.2E-04		
2.0E+00	I	5.7E-04	I						7.1	1	0.7	No	~Polychlorinated Biphenyls (high risk)	1336-36-3										
4.0E-01	I	1.0E-04	I						7.1	1	0.7	No	~Polychlorinated Biphenyls (low risk)	1336-36-3	1.9E-01		5.6E-02	4.4E-02					5.0E-01	
7.0E-02	I	2.0E-05	I						7.1	1	0.7	No	~Polychlorinated Biphenyls (lowest risk)	1336-36-3										
1.3E+01	E	3.8E-03	E	7.0E-06	E	4.0E-04	E		6.63	1	0.6	No	~Tetrachlorobiphenyl, 3,3',4,4'-(PCB 77)	32598-13-3	6.0E-03			6.0E-03	1.4E-01			1.4E-01		
3.9E+01	E	1.1E-02	E	2.3E-06	E	1.3E-04	E	V	6.34	1	0.7	No	~Tetrachlorobiphenyl, 3,4,4',5-(PCB 81)	70362-50-4	2.0E-03		4.9E-04	4.0E-04	4.7E-02		2.8E-01	4.0E-02		
									10.46	1	0	No	Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9										
												No	Polynuclear Aromatic Hydrocarbons (PAHs)											
									3.92	1	1	Yes	~Acenaphthene	83-32-9					1.2E+03	9.6E+02		5.3E+02		
									4.45	1	1	Yes	~Anthracene	120-12-7					6.0E+03	2.5E+03		1.8E+03		
7.3E-01	E	1.1E-04	C						5.76	1	1	No	~Benz[a]anthracene	56-55-3	3.4E-02		1.8E-02	1.2E-02						
1.2E+00	C	1.1E-04	C						6.11	1	0.9	No	~Benzo[<i>b</i>]fluoranthene	205-82-3	6.5E-02			6.5E-02						
7.3E+00	I	1.1E-03	C						6.13	1	1	No	~Benzo[<i>a</i>]pyrene	50-32-8	3.4E-03			3.4E-03					2.0E-01	
7.3E-01	E	1.1E-04	C						5.78	1	1	No	~Benzo[<i>b</i>]fluoranthene	205-99-2	3.4E-02			3.4E-02						
7.3E-02	E	1.1E-04	C						6.11	1	0.9	No	~Benzo[<i>k</i>]fluoranthene	207-08-9	3.4E-01			3.4E-01						
7.3E-03	E	1.1E-05	C	8.0E-02	I				3.9	1	1	Yes	~Chloronaphthalene, Beta-	91-58-7					1.6E+03	1.4E+03		7.5E+02		
									5.81	1	1	No	~Chrysene	218-01-9	3.4E+00			3.4E+00						
7.3E+00	E	1.2E-03	C						6.75	1	0.6	No	~Dibenz[<i>a,h</i>]anthracene	53-70-3	3.4E-03			3.4E-03						

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Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -y) ⁻¹	k _e (y)	RfD _c (mg/kg-day)	k _e (y)	RfC _c (mg/m ³ -y)	k _e (y)	v	mutagen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	
		3.4E-05	C	4.0E-03 2.0E-02	I	3.0E-03	I	V		3.86 3.3	1	1	Yes	-Methylnaphthalene, 2- -Naphthalene	91-57-6 91-20-3			1.7E-01	1.7E-01		8.0E+1 4.0E+2	6.5E+1 7.0E+2	6.3E+00 6.3E+00	3.6E+01 6.1E+00	
1.2E+00	C	1.1E-04	C	3.0E-02 2.0E-02	I			V		4.75 4.88 -0.33	1	0.9	Yes	-Nitropyrene, 4- -Pyrene Potassium Perfluorobutane Sulfonate	57835-92-4 129-00-0 29420-49-3	6.5E-02	2.6E-02		1.9E-02		6.0E+2 4.0E+2	1.5E+2 2.8E+05		1.2E+2 4.0E+2	
1.5E-01	I			9.0E-03 6.0E-03 1.5E-02	I			H		4.1 5.58 2.99	1	0.9	Yes	Prochloraz Profuralin Prometon	67747-09-5 26399-36-0 1610-18-0	5.2E-01	1.3E+00		3.7E-01		1.8E+2 1.2E+2 3.0E+2	5.1E+2 3.3E+1 1.6E+03		1.3E+2 2.6E+1 2.5E+2	
				4.0E-03 1.3E-02 5.0E-03	I					3.51 2.18 3.07	1	0.9	Yes	Prometryn Propachlor Propanil	7287-19-6 1918-16-7 709-98-8						8.0E+1 2.6E+2 1.0E+2	2.3E+2 4.3E+03 4.4E+2		6.0E+1 2.5E+2 8.2E+1	
				2.0E-02 2.0E-03 2.0E-02	I				V	5 -0.38 2.93	1	0.8	Yes	Propargite Propargyl Alcohol Propazine	2312-35-8 107-19-7 139-40-2						4.0E+2 4.0E+1 4.0E+2	2.7E+2 1.2E+4 2.4E+03		1.6E+2 4.0E+1 3.4E+2	
				2.0E-02 1.3E-02	I					2.6 3.72	1	1	Yes	Propam Propiconazole	122-42-9 60207-90-1						4.0E+2 2.6E+2	2.8E+03 1.1E+03		3.5E+2 2.1E+2	
				1.0E-01	X	1.0E+00	X	V		3.69	1	1	Yes	Propyl benzene	103-65-1						2.0E+3	1.8E+3	2.1E+03	6.6E+2	
				2.0E+01	P	3.0E+00	C	V		1.77 -0.92	1	1	Yes	Propylene Propylene Glycol	115-07-1 57-55-6						4.0E+05	3.2E+08	6.3E+03	4.0E+05	
				7.0E-01 7.0E-01	H	2.0E+00	I	V		1.59 0.002 -0.49	1	1	Yes	Propylene Glycol Dinitrate Propylene Glycol Monomethyl Ether Propylene Glycol Monomethyl Ether	6423-43-4 1569-02-4 107-98-2						1.4E+4 1.4E+4	3.3E+6 3.9E+6	4.2E+03	1.4E+4 3.2E+03	
2.4E-01	I	3.7E-06	I	2.5E-01 2.5E-02	I	3.0E-02	I	V		0.03 2.6 6.2	1	1	Yes	Propylene Oxide Pursuit Pyridin	75-56-9 81335-77-5 51630-58-1	3.2E-01	4.5E+01	1.5E+00	2.7E-01		5.0E+3 5.0E+2	7.2E+4	6.3E+01	4.7E+03 5.0E+2	
3.0E+00	I			1.0E-03 5.0E-04	I			V		0.65 4.44 2.03	1	0.9	Yes	Pyridine Quinalphos Quindole	110-86-1 3393-03-8 91-22-5	2.6E-02	2.8E-01		2.4E-02		2.0E+1 1.0E+1	1.5E+3 1.0E+1		2.0E+1 5.1E+00	
				3.0E-02 5.0E-02	I	3.0E-02	A			6.14 4.88	1	0	Yes	Refractory Ceramic Fibers Resmethrin	NA 10453-86-8						6.0E+2 1.0E+3	7.6E+1 6.8E+2		6.7E+1 4.1E+2	
2.2E-01	C	6.3E-05	C	4.0E-03 2.5E-02	I				M	4.1 3.45 5.57	1	0.9	Yes	Rotenone Safrole Savay	83-79-4 94-59-7 78587-05-0	1.1E-01	5.9E-01		9.5E-02		8.0E+1 5.0E+2	2.6E+2 1.4E+2		6.1E+1 1.1E+2	
				5.0E-03 5.0E-03 5.0E-03	I	2.0E-02	C			1 1 1	1	1	Yes	Selenious Acid Selenium Selenium Sulfide	7783-00-8 782-49-2 7446-34-6						1.0E+2 1.0E+2 1.0E+2	2.3E+4 2.3E+4 2.3E+4		1.0E+2 1.0E+2 1.0E+2	5.0E+1
				9.0E-02 5.0E-03	I	3.0E-03	C			4.38 1 0.04	1	0.9	Yes	Sethoxydim Silica (crystalline, respirable) Silver	74051-80-2 7631-96-9 7440-22-4						1.8E+3 1.0E+2	2.4E+3		1.0E+3 9.4E+1	
1.2E-01	H			1.3E-02 4.0E-03	I					2.18 0.37 1	1	1	Yes	Simazine Sodium Acifluorfen Sodium Azide	122-34-9 62476-59-9 26628-22-8	6.5E-01	8.9E+00		6.1E-01		1.0E+2 2.6E+2 8.0E+1	1.6E+3 2.1E+05 1.8E+4		9.4E+1 2.6E+2 8.0E+1	4.0E+00
5.0E-01	C	1.5E-01	C	2.0E-02 3.0E-02 5.0E-02	C	2.0E-04	C		M	0.025 -1.431 1	1	1	Yes	Sodium Dichromate Sodium Diethyldithiocarbamate Sodium Fluoride	10588-01-9 148-18-5 7681-49-4	5.0E-02	2.3E-01		4.1E-02		4.0E+2 6.0E+2 1.0E+3	2.3E+3 1.9E+06 2.3E+05		3.4E+2 6.0E+2 1.0E+3	
2.4E-02	H			2.0E-05 1.0E-03 3.0E-02	I					-3.78 1 3.53	1	1	No	Sodium Fluoroacetate Sodium Metavanadate Stirofos (Tetrachlorovinphos)	62-74-8 13718-26-8 961-11-5	3.2E+00	1.8E+01		2.8E+00		4.0E-01 2.0E+1 6.0E+2	4.5E+3 4.5E+3 3.8E+03		4.0E-01 2.0E+1 5.2E+2	
5.0E-01	C	1.5E-01	C	2.0E-02 6.0E-01 3.0E-04	C	2.0E-04	C		M	0.025 1 1.93	1	1	Yes	Strontium Chromate Strontium, Stable Strychnine	7789-06-2 7440-24-6 57-24-9	5.0E-02	2.3E-01		4.1E-02		4.0E+2 1.2E+4 6.0E+00	2.3E+3 2.7E+06 3.2E+02		3.4E+2 1.2E+4 5.9E+00	
				2.0E-01 3.0E-03 1.0E-03	I	1.0E+00	I	V		2.95 3.1 -0.77	1	1	Yes	Styrene Styrene-Acrylonitrile (SAN) Trimer Sulfolane	100-42-5 NA 126-33-0						4.0E+3 6.0E+1 2.0E+1	1.0E+4 2.4E+2 1.7E+4	2.1E+03	1.2E+3 4.8E+1 2.0E+1	1.0E+2
				8.0E-04	P	1.0E-03	C	V		3.9	1	0.9	Yes	Sulfonylbis(4-chlorobenzene), 1,1'- Sulfur Trioxide Sulfuric Acid	80-07-9 7446-11-9 7664-93-9						1.6E+1	3.5E+1	2.1E+00	1.1E+1 2.1E+00	
				2.5E-02 3.0E-02 7.0E-02	I					2.94 3.3 1.79	1	0.9	Yes	Systhane TCMTB Tebuthiuron	88671-89-0 21564-17-0 34014-18-1						5.0E+2 6.0E+2 1.4E+3	4.8E+3 2.4E+3 4.7E+4		4.5E+2 4.8E+2 1.4E+3	
				2.0E-02 1.3E-02	H					5.96 1.89	1	0.7	No	Temphos Terbacil	3383-96-8 5902-51-2						4.0E+2 2.6E+2	7.0E+3		4.0E+2 2.5E+2	

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Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -1)	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³ -1)	k _e (y)	v	mutagen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)
				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	
				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-04	I		V			4.64	1	1	Yes	Tetrachlorobenzene, 1,2,4,5-	95-94-3					6.0E+00	2.4E+00		1.7E+00	
2.6E-02	I	7.4E-06	I	3.0E-02	I		V			2.93	1	1	Yes	Tetrachloroethane, 1,1,1,2-	630-20-6	3.0E+00	1.0E+01	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	Yes	Tetrachloroethane, 1,1,2,2-	79-34-5	3.9E-01	3.1E+00	9.7E-02	7.6E-02	4.0E+02	3.6E+03		3.6E+02	
2.1E-03	I	2.6E-07	I	6.0E-03	I		V			3.4	1	1	Yes	Tetrachloroethylene	127-18-4	3.7E+01	6.3E+01	2.2E+01	1.1E+01	1.2E+02	2.3E+02	8.3E+01	4.1E+01	5.0E+00
2.0E+01	H			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	
				5.0E-04	I		V			4.54	1	0.9	Yes	Tetrachlorotoluene, p- alpha, alpha, alpha-Tetraethyl Dithiopyrophosphate	5216-25-1 3689-24-5	3.9E-03	1.9E-03		1.3E-03				7.1E+00	
				8.0E+01	I	V				1.68	1	1	Yes	Tetrafluoroethane, 1,1,1,2-	811-97-2							1.7E+05		1.7E+05
				2.0E-03	P					1.64	1	1	Yes	Tetryl (Trinitrophenylmethyl nitramine)	479-45-8					4.0E+01	2.5E+03		3.9E+01	
				7.0E-06	X						1	1	Yes	Thallium (I) Nitrate	10102-45-1					1.4E-01	3.2E+01		1.4E-01	
				1.0E-05	X						1	1	Yes	Thallium (Soluble Salts)	7440-28-0					2.0E-01	4.5E+01		2.0E-01	2.0E+00
				6.0E-06	X		V			-0.17	1	1	Yes	Thallium Acetate	563-68-8					1.2E-01	1.0E+02		1.2E-01	
				2.0E-05	X						1	1	Yes	Thallium Carbonate	6533-73-9					4.0E-01	3.7E+03		4.0E-01	
				6.0E-06	X						1	1	Yes	Thallium Chloride	7791-12-0					1.2E-01	2.7E+01		1.2E-01	
				2.0E-05	X						1	0.9	Yes	Thallium Sulfate	7446-18-6					4.0E-01	9.1E+01		4.0E-01	
				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	Yes	Thiodiglycol	111-48-8					1.4E+03	9.6E+05		1.4E+03	
				3.0E-04	H					2.16	1	1	Yes	Thiofanox	39196-18-4					6.0E+00	4.4E+01		5.3E+00	
				8.0E-02	I					1.4	1	1	Yes	Thiophanate, Methyl	23564-05-8					1.6E+03	2.0E+05		1.6E+03	
				5.0E-03	I					1.73	1	1	Yes	Thiram	137-26-8					1.0E+02	4.0E+03		9.8E+01	
				6.0E-01	H						1	1	Yes	Tin	7440-31-5					1.2E+04	2.7E+06		1.2E+04	
				1.0E-04	A	V					1	1	Yes	Titanium Tetrachloride	7550-45-0							2.1E-01	2.1E-01	
1.8E-01	X			8.0E-02	I	5.0E+00	I	V		2.73	1	1	Yes	Toluene	108-88-3					1.6E+03	5.3E+03	1.0E+04	1.1E+03	1.0E+03
3.0E-02	P			2.0E-04	X					0.16	1	1	Yes	Toluene-2,5-diamine	95-70-5	4.3E-01	7.9E+01		4.3E-01	4.0E+00	8.3E+02		4.0E+00	
				4.0E-03	X					1.39	1	1	Yes	Toluidine, p-	106-49-0	2.6E+00	6.5E+01		2.5E+00	8.0E+01	2.3E+03		7.7E+01	
				3.0E+00	P		V			6.1	1	1	No	Total Petroleum Hydrocarbons (Aliphatic High)	NA					6.0E+04			6.0E+04	
				6.0E-01	P	V				3.9	1	1	Yes	Total Petroleum Hydrocarbons (Aliphatic Low)	NA							1.3E+03	1.3E+03	
				1.0E-02	X	1.0E-01	P	V		5.65	1	1	No	Total Petroleum Hydrocarbons (Aliphatic Medium)	NA					2.0E+02		2.1E+02	1.0E+02	
				4.0E-02	P					5.16	1	1	No	Total Petroleum Hydrocarbons (Aromatic High)	NA					8.0E+02			8.0E+02	
				4.0E-03	P	3.0E-02	P	V		2.13	1	1	Yes	Total Petroleum Hydrocarbons (Aromatic Low)	NA				8.0E+01	6.0E+02	6.3E+01	3.3E+01	3.3E+01	
				4.0E-03	P	3.0E-03	P	V		3.58	1	1	Yes	Total Petroleum Hydrocarbons (Aromatic Medium)	NA				8.0E+01	9.0E+01	6.3E+00	5.5E+00	5.5E+00	
1.1E+00	I	3.2E-04	I							5.9	1	0.8	Yes	Toxaphene	9001-35-2	7.1E-02	1.9E-02		1.5E-02				1.5E+02	3.0E+00
7.0E-03	X			7.5E-03	I					7.56	1	0.5	No	Traloxmethrin	68841-25-6					6.0E+00	9.8E+00		3.7E+00	
				3.0E-04	A		V			4.1	1	0.9	Yes	Tri-n-butyltin	688-73-3								1.6E+06	
				8.0E+01	X					0.25	1	1	Yes	Triacetin	102-76-1					1.6E+06	5.3E+08		1.6E+06	
				1.3E-02	I		V			4.6	1	0.9	Yes	Triallate	2303-17-5					2.6E+02	2.2E+02		1.2E+02	
				1.0E-02	I					1.1	1	1	Yes	Trisulfuron	82097-50-b					2.0E+02	6.0E+04		2.0E+02	
9.0E-03	P			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	
				1.0E-02	P					4	1	0.9	Yes	Tributyl Phosphate	126-73-8	8.7E+00	1.2E+01		5.1E+00	2.0E+02	3.3E+02		1.2E+02	
				3.0E-04	P						1	0	No	Tributyltin Compounds	NA				6.0E+00			6.0E+00		
				3.0E-04	I					4.05	1	1	Yes	Tributyltin Oxide	58-35-0					6.0E+00	9.5E+01		5.7E+00	
7.0E-02	I			3.0E+01	I	3.0E+01	H	V		3.16	1	1	Yes	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1					6.0E+05	1.9E+06	6.3E+04	5.5E+04	
				2.0E-02	I					1.33	1	1	Yes	Trichloroacetic Acid	76-03-9	1.1E+00	4.4E+01		1.1E+00	4.0E+02	1.8E+04		3.9E+02	6.0E+01
2.9E-02	H									-0.67	1	1	Yes	Trichloroaniline HCl, 2,4,6-	33663-50-2	2.7E+00	3.6E+03		2.7E+00				4.0E-01	
7.0E-03	X			3.0E-05	X					3.52	1	1	Yes	Trichloroaniline, 2,4,6-	634-93-5	1.1E+01	1.9E+01		7.0E+00	6.0E-01	1.2E+00		4.0E-01	
				8.0E-04	X		V			4.05	1	1	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	Yes	Trichlorobenzene, 1,2,4-	120-82-1	2.7E+00	1.9E+00		1.1E+00	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	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	Yes	Trichloroethane, 1,1,2-	79-00-5	1.4E+00	1.9E+01	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	Yes	Trichloroethylene	79-01-6	1.2E+00	7.2E+00	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	7.0E-01	H	V		2.53	1	1	Yes	Trichlorofluoromethane	75-69-4					6.0E+03	3.6E+04	1.5E+03	1.1E+03	
				1.0E-01	I					3.72	1	1	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	Yes	Trichlorophenol, 2,4,6-	88-06-2	7.1E+00	9.4E+00		4.0E+00	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-0																				

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Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHIL Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³ -y) ⁻¹	k e y	RfD _c (mg/kg-day)	k e y	RfC _c (mg/m ³)	k e y	v o l a t i l e	m u t a g e n	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)
7.7E-03	I			2.0E+00	P					-1.75	1	1	Yes	Triethylene Glycol	112-27-6					4.0E+04	1.8E+08		4.0E+04	
2.0E-02	P			7.5E-03	I			V		5.34	1	0.8	Yes	Trifluralin	1582-09-8	1.0E+01	3.3E+00		2.5E+00	1.5E+02	5.5E+01		4.0E+01	
				1.0E-02	P					-0.65	1	1	Yes	Trimethyl Phosphate	512-56-1	3.9E+00	2.7E+03		3.9E+00	2.0E+02	1.6E+05		2.0E+02	
				5.0E-03	P	V				3.66	1	1	Yes	Trimethylbenzene, 1,2,3-	526-73-8							1.0E+01	1.0E+01	
				7.0E-03	P	V				3.63	1	1	Yes	Trimethylbenzene, 1,2,4-	95-63-6							1.5E+01	1.5E+01	
				1.0E-02	X			V		3.42	1	1	Yes	Trimethylbenzene, 1,3,5-	108-67-8					2.0E+02	2.8E+02		1.2E+02	
3.0E-02	I			3.0E-02	I					1.18	1	1	Yes	Trinitrobenzene, 1,3,5-	99-35-4				2.5E+00	6.0E+02	4.7E+04		5.9E+02	
				5.0E-04	I					1.6	1	1	Yes	Trinitrotoluene, 2,4,6-	118-96-7	2.6E+00	1.0E+02			1.0E+01	4.5E+02		9.8E+00	
				2.0E-02	P					2.83	1	1	Yes	Triphenylphosphine Oxide	/91-28-6					4.0E+02	3.8E+03		3.6E+02	
2.3E+00	C	6.6E-04	C	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	Yes	Tris(1-chloro-2-propyl)phosphate	13674-84-5					2.0E+02	3.8E+03		1.9E+02	
								V		4.29	1	1	No	Tris(2,3-dibromopropyl)phosphate	126-72-7	3.4E-02		8.5E-03	6.8E-03					
2.0E-02	P			7.0E-03	P					1.44	1	1	Yes	Tris(2-chloroethyl)phosphate	115-96-8	3.9E+00	2.9E+02		3.8E+00	1.4E+02	1.2E+04		1.4E+02	
3.2E-03	P			1.0E-01	P					9.49	1	0	No	Tris(2-ethylhexyl)phosphate	78-42-2	2.4E+01			2.4E+01	2.0E+03			2.0E+03	
				3.0E-03	I	4.0E-05	A			1	1	1	Yes	Uranium (Soluble Salts)	NA					6.0E+01	1.4E+04		6.0E+01	3.0E+01
1.0E+00	C	2.9E-04	C					M		-0.15	1	1	Yes	Urethane	51-79-6	2.5E-02	5.9E+00		2.5E-02					
		8.3E-03	P	9.0E-03	I	7.0E-06	P			0.026	1	1	Yes	Vanadium Pentoxide	1314-62-1					1.8E+02	1.1E+03		1.5E+02	
				5.0E-03	S	1.0E-04	A			0.026	1	1	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	Yes	Verdigris	1929-77-7					2.0E+01	2.5E+01		1.1E+01	
				2.5E-02	I					3.1	1	0.9	Yes	Vinclozolin	50471-44-8					5.0E+02	3.7E+03		4.4E+02	
				1.0E+00	H	2.0E-01	I	V		0.73	1	1	Yes	Vinyl Acetate	108-05-4					2.0E+04	1.4E+06	4.2E+02	4.1E+02	
7.2E-01	I	4.4E-06	I	3.0E-03	I	1.0E-01	I	V	M	1.57	1	1	Yes	Vinyl Bromide	593-60-2			1.8E-01	1.8E-01	6.0E+01	8.9E+02	6.3E+00	6.3E+00	
				3.0E-04	I					1.62	1	1	Yes	Vinyl Chloride	75-01-4	2.1E-02	2.7E-01	3.4E-01	1.9E-02	6.0E+00	8.4E+01	2.1E+02	4.4E+01	2.0E+00
										2.7	1	1	Yes	Warfarin	81-81-2					6.0E+00	8.4E+01		5.6E+00	
				2.0E-01	S	1.0E-01	S	V		3.15	1	1	Yes	Xylene, p-	106-42-3					4.0E+03	7.6E+03	2.1E+02	1.9E+02	
				2.0E-01	S	1.0E-01	S	V		3.2	1	1	Yes	Xylene, m-	108-38-3					4.0E+03	7.1E+03	2.1E+02	1.9E+02	
				2.0E-01	S	1.0E-01	S	V		3.12	1	1	Yes	Xylene, o-	95-47-6					4.0E+03	8.0E+03	2.1E+02	1.9E+02	
				2.0E-01	I	1.0E-01	I	V		3.16	1	1	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	1	Yes	Zinc Phosphide	1314-84-7					6.0E+00	2.3E+03		6.0E+00	
				3.0E-01	I					1	1	1	Yes	Zinc and Compounds	7440-66-6					6.0E+03	2.3E+06		6.0E+03	
				5.0E-02	I					1.3	1	1	Yes	Zineb	12122-67-7					1.0E+03	9.7E+04		9.9E+02	
				8.0E-05	X					1	1	1	Yes	Zirconium	7440-67-7					1.6E+00	3.6E+02		1.6E+00	

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Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1				Protection of Groundwater SSL					
SFO (mg/kg-day) ¹	k _e	IUR (ug/m ³ -y) ¹	k _e	RD ₅₀ (mg/kg-day)	k _e	RF _c (mg/m ³ -y)	k _e	muta-gen	LOGP	GIABS	FA	IN EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HO=1 (ug/L)	Dermal SL Child HO=1 (ug/L)	Inhalation SL Child HO=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)
1.8E-02	C	5.1E-06	C	1.5E-01	I				-1.5	1	1	Yes	ALAR	1596-84-5	4.3E+00	1.3E+04		4.3E+00	3.0E+03	1.0E+07		3.0E+03		9.5E-04	
8.7E-03	I			4.0E-03	I				-0.85	1	1	Yes	Acephate	30560-19-1	9.0E+00	1.1E+04		8.9E+00	8.0E+01	1.1E+05		8.0E+01		2.0E-03	
		2.2E-06	I			9.0E-03	I	V	-0.34	1	1	Yes	Acetaldehyde	75-07-0			2.6E+00	2.6E+00			1.9E+01	1.9E+01		5.2E-04	
				2.0E-02	I				3.03	1	0.9	Yes	Acetochlor	34256-82-1					4.0E+02	2.9E+03		3.5E+02		2.8E-01	
				9.0E-01	I	3.1E+01	A	V	-0.24	1	1	Yes	Acetone	67-64-1					1.8E+04	4.4E+06	6.4E+04	1.4E+04		2.9E+00	
						2.0E-03	X	V	-0.03	1	1	Yes	Acetone Cyanohydrin	75-86-5								4.2E+00		8.4E-04	
						6.0E-02	I	V	-0.34	1	1	Yes	Acetonitrile	75-05-8								1.3E+02		2.6E-02	
3.8E+00	C	1.3E-03	C	1.0E-01	I				1.58	1	1	Yes	Acetophenone	98-06-2					2.0E+03	4.6E+04		1.9E+03		5.8E-01	
									3.12	1	1	Yes	Acetylaminofluorene, 2-	53-96-3	2.1E-02	6.4E-02		1.6E-02						7.2E-05	
				5.0E-04	I	2.0E-05	I	V	-0.01	1	1	Yes	Acrolein	107-02-8					1.0E+01	1.7E+03	4.2E-02	4.2E-02		8.4E-06	
5.0E-01	I	1.0E-04	I	2.0E-03	I	6.0E-03	I	M	-0.67	1	1	Yes	Acrylamide	79-06-1	5.0E-02	2.3E+01		5.0E-02	4.0E+01	2.1E+04		4.0E+01		1.1E-05	
				5.0E-01	I	1.0E-03	I	V	0.35	1	1	Yes	Acrylic Acid	79-10-7					1.0E+04	1.1E+06	2.1E+00	2.1E+00		4.2E-04	
5.4E-01	I	6.8E-05	I	4.0E-02	A	2.0E-03	I	V	0.25	1	1	Yes	Acrylonitrile	107-13-1	1.4E-01	1.4E+01	8.3E-02	5.2E-02	8.0E+02	8.8E+04	4.2E+00	4.1E+00		1.1E-05	
5.6E-02	C			1.0E-02	I	6.0E-03	P		-0.32	1	1	Yes	Adiponitrile	111-69-3											
									3.52	1	0.9	Yes	Alachlor	15972-60-8	1.4E+00	4.2E+00		1.0E+00	2.0E+02	6.9E+02		1.6E+02	2.0E+00	8.6E-04	1.7E-03
				1.0E-03	I				1.13	1	1	Yes	Aldicarb	116-06-3					2.0E+01	1.4E+03		2.0E+01	3.0E+00	4.9E-03	7.5E-04
				1.0E-03	I				-0.57	1	1	Yes	Aldicarb Sulfone	1646-88-4					2.0E+01	2.4E+04		2.0E+01	2.0E+00	4.4E-03	8.8E-04
									-0.78	1	1	Yes	Aldicarb sulfoxide	1646-87-3								4.0E+00	4.0E+00		
1.7E+01	I	4.9E-03	I	3.0E-05	I				6.5	1	1	No	Aldrin	309-00-2	4.6E-03		1.1E-03	9.2E-04	6.0E-01			6.0E-01		1.5E-04	
				2.5E-01	I				2.2	1	1	Yes	Allyl	74223-64-6					5.0E+03	2.4E+05		4.9E+03		1.9E+00	
				5.0E-03	I	1.0E-04	X	V	0.17	1	1	Yes	Allyl Alcohol	107-18-6					1.0E+02	1.3E+04	2.1E-01	2.1E-01		4.2E-05	
2.1E-02	C	6.0E-06	C	1.0E+00	P	5.0E-03	P		1.93	1	1	Yes	Allyl Chloride	107-05-1	3.7E+00	3.3E+01	9.4E-01	7.3E-01			2.1E+00	2.1E+00		2.3E-04	3.0E+04
				4.0E-04	I					1	1	Yes	Aluminum	7429-90-5					2.0E+04	4.5E+06		2.0E+04			
										1	1	Yes	Aluminum Phosphide	20859-73-8					8.0E+00	1.8E+03		8.0E+00			
				3.0E-04	I				2.31	1	1	Yes	Amdro	67485-29-4					6.0E+00	5.1E+02		5.9E+00		2.1E+03	
2.1E+01	C	6.0E-03	C	9.0E-03	I				2.98	1	1	Yes	Ametryn	834-12-8					1.8E+02	9.7E+02		1.5E+02		1.6E-01	1.5E-05
									2.86	1	1	Yes	Amnobeiphenyl, 4-	92-67-1	3.7E-03	1.5E-02		3.0E-03							
				8.0E-02	P				0.21	1	1	Yes	Amnophenol, m	591-27-5					1.6E+03	2.8E+05		1.6E+03		6.1E-01	
				2.0E-02	P				0.04	1	1	Yes	Amnophenol, p	123-30-8					4.0E+02	9.1E+04		4.0E+02		1.5E-01	
				2.5E-03	I				5.5	1	0.9	Yes	Amtraz	33089-61-1					5.0E+01	9.7E+00		8.2E+00		4.2E+00	
						1.0E-01	I	V	0.23	1	1	Yes	Ammonia	7664-41-7											
				2.0E-01	I					1	1	Yes	Ammonium Sulfamate	7773-06-0					4.0E+03	9.1E+05		4.0E+03			
						3.0E-03	X	V	0.89	1	1	Yes	Amyl Alcohol, tert-	75-85-4							6.3E+00	6.3E+00		1.3E-03	
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I		0.9	1	1	Yes	Aniline	62-53-3	1.4E+01	6.6E+02		1.3E+01	1.4E+02	7.7E+03		1.4E+02		4.6E-03	
4.0E-02	P			2.0E-03	X				3.39	1	0.9	Yes	Anthraquinone, 9,10-	84-85-1	1.9E+00	4.9E+00		1.4E+00	4.0E+01	1.1E+02		3.0E+01		1.4E-02	
				4.0E-04	I				0.15	1	1	Yes	Antimony (metallic)	7440-36-0					8.0E+00	2.7E+02		7.8E+00	6.0E+00	3.5E-01	2.7E-01
				5.0E-04	H				0.15	1	1	Yes	Antimony Pentoxide	1314-60-9					1.0E+01	3.4E+02		9.7E+00			
				9.0E-04	H				-7.28	0.15	1	No	Antimony Potassium Tartrate	11071-15-1					1.8E+01			1.8E+01			
				4.0E-04	H				0.15	1	1	Yes	Antimony Trisulfide	1332-81-8					8.0E+00	2.7E+02		7.8E+00			
						2.0E-04	I		0.15	1	1	Yes	Antimony Trisulfide	1309-84-4											
2.5E-02	I	7.1E-06	I	1.3E-02	I				3.1	1	0.9	Yes	Apollo	74115-24-5					2.6E+02	2.1E+03		2.3E+02		1.4E+01	
				5.0E-02	H				4.82	1	0.8	Yes	Aramite	140-57-8	3.1E+00	2.3E+00		1.3E+00	1.0E+03	8.2E+02		4.5E+02		1.5E-02	
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C			1	1	Yes	Arsenic, inorganic	7440-38-2	5.2E-02	9.3E+00		5.2E-02	6.0E+00	1.4E+03		6.0E+00	1.0E+01	1.5E-03	2.9E-01
				3.5E-06	C	5.0E-05	I			1	1	Yes	Arsine	7784-42-1					7.0E-02	1.6E+01		7.0E-02		1.9E+00	
				9.0E-03	I				4.28	1	0.9	Yes	Assure	76578-14-8					1.8E+02	3.8E+02		1.2E+02			
2.3E-01	C			5.0E-02	I				-0.27	1	1	Yes	Asulam	3337-71-1					1.0E+03	8.0E+05		1.0E+03	3.0E+00	2.6E-01	2.0E-03
8.8E-01	C	2.5E-04	C	3.5E-02	I				2.61	1	1	Yes	Atrazine	1912-24-9	3.4E-01	2.6E+00		3.0E-01	7.0E+02	6.2E+03		6.3E+02		2.0E-04	2.0E-03
									2.98	1	0.9	Yes	Auramine	492-80-8	8.9E-02	2.6E-01		6.6E-02						6.0E-04	
1.1E-01	I	3.1E-05	I	4.0E-04	I				4.48	1	1	No	Avermectin B1	65195-55-3					8.0E+00			8.0E+00		1.4E+01	
				1.0E+00	P	7.0E-06	P		3.82	1	1	Yes	Azobenzene	103-33-3	7.1E-01	7.0E-01	1.8E-01	1.2E-01				2.0E+04		9.2E-04	6.8E+00
									-1.7	1	1	Yes	Azodicarbonamide	123-77-3								2.0E+04			
5.0E-01	C	1.5E-01	C	2.0E-01	I	5.0E-04	H		0.07	1	1	Yes	Barium	7440-39-3					4.0E+03	6.4E+04		3.8E+03	2.0E+03	1.6E+02	8.2E+01
				2.0E-02	C	2.0E-04	C	M	0.025	1	1	Yes	Barium Chromate	10294-40-3	5.0E-02	2.3E-01		4.1E-02	4.0E+02	2.3E+03		3.4E+02			
				4.0E-03	I				1.52	1	1	Yes	Baygon	114-26-1					8.0E+01	3.6E+03		7.8E+01		2.5E-02	

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where nSL < 100X c SL; ** = where nSL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1				Protection of Groundwater SSL						
SFO (mg/kg-day) ¹	k _e	IUR (ug/m ³ -y)	k _e	RD ₅₀ (mg/kg-day)	k _e	RF _c (mg/m ³ -y)	k _e	muta-gen	LOGP	GIABS	FA	IN EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)	
1.1E+00	I	3.3E-04	I					V	1.29	1	1	Yes	Bis(2-chloroethyl)ether	111-44-4	7.1E-02	2.6E+00	1.7E-02	1.4E-02						3.6E-06		
2.2E+02	I	6.2E-02	I					V	0.57	1	1	Yes	Bis(chloromethyl)ether	542-88-1	3.5E-04	3.2E-02	9.1E-05	7.2E-05						1.7E-08		
				5.0E-02	I			V	3.32	1	1	Yes	Bisphenol A	80-05-7					1.0E+03	3.2E+03		7.7E+02		5.8E+01		
				2.0E-01	I	2.0E-02	H	V		1	1	Yes	Boron And Borates Only	7440-42-8					4.0E+03	9.1E+05		4.0E+03		1.3E+01		
				2.0E+00	P	2.0E-02	P	V	1.16	1	1	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	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			V		1	1	Yes	Bromate	15541-45-4	1.1E-01	2.0E+01		1.1E-01	8.0E+01	1.8E+04		8.0E+01	1.0E+01	8.5E-04	7.7E-02	
2.0E+00	X	6.0E-04	X					V	1.92	1	1	Yes	Bromo-2-chloroethane, 1-Bromobenzene	107-04-0	3.9E-02	5.5E-01	9.4E-03	7.4E-03						2.1E-06		
				8.0E-03	I	6.0E-02	I	V	2.99	1	1	Yes		108-96-1					1.6E+02	5.4E+02	1.3E+02	6.2E+01		4.2E-02		
								V	1.41	1	1	Yes	Bromochloromethane	74-97-5							8.3E+01		8.3E+01		2.1E-02	
6.2E-02	I	3.7E-05	C	2.0E-02	I			V	2	1	1	Yes	Bromodichloromethane	75-27-4	1.3E+00	1.8E+01	1.5E-01	1.3E-01	4.0E+02	6.4E+03		3.8E+02	8.0E+01(F)	3.7E-05	2.2E-02	
7.9E-03	I	1.1E-06	I	2.0E-02	I			V	2.4	1	1	Yes	Bromoforn	75-25-2	9.9E+00	1.4E+02	5.1E+00	3.3E+00	4.0E+02	6.2E+03		3.8E+02	8.0E+01(F)	8.7E-04	2.1E-02	
				1.4E-03	I	5.0E-03	I	V	1.19	1	1	Yes	Bromomethane	74-83-9					2.8E+01	2.0E+03	1.0E+01	7.5E+00		1.9E-03		
				5.0E-03	H			V	5.21	1	0.8	Yes	Bromophos	2104-96-3					1.0E+02	5.5E+01		3.5E+01		1.5E-01		
				2.0E-02	I			V	3.39	1	0.9	Yes	Bromoxynil	1689-84-5					4.0E+02	1.8E+03		3.3E+02		2.8E-01		
3.4E+00	C	3.0E-05	I	2.0E-02	I			V	5.4	1	0.8	Yes	Bromoxynil Octanoate	1689-99-2					4.0E+02	2.1E+02		1.4E+02		1.2E+00		
				1.0E-01	I	2.0E-03	I	V	1.99	1	1	Yes	Butadiene, 1,3-	106-99-0	2.3E-02	1.6E-01	1.9E-01	1.8E-02			4.2E+00	4.2E+00	2.0E+03	9.9E-06	4.1E-01	
1.9E-03	P			2.0E-01	I			V	0.88	1	1	Yes	Butanol, n-	71-36-3					2.0E+03	1.0E+05		2.0E+03		1.2E+00		
				2.0E-01	I			V	4.73	1	0.9	Yes	Butyl Benzyl Phthlate	85-68-7	4.1E+01	2.6E+01		1.6E+01	4.0E+03	2.9E+03		1.7E+03		2.3E-01		
				2.0E+00	P	3.0E+01	P	V	0.61	1	1	Yes	Butyl alcohol, sec-	78-92-2					4.0E+04	3.0E+06	6.3E+04	2.4E+04		5.0E+00		
				5.0E-02	I			V	4.15	1	1	Yes	Butylate	2008-41-5					1.0E+03	8.5E+02		4.6E+02		4.5E-01		
2.0E-04	C	5.7E-08	C					V	3.5	1	1	Yes	Butylated hydroxyanisole	25013-16-5	3.9E+02	6.2E+02		2.4E+02						4.5E-01		
3.6E-03	P			3.0E-01	P			V	5.1	1	1	Yes	Butylated hydroxytoluene	128-37-0	2.2E+01	3.8E+00		3.3E+00	6.0E+03	1.2E+03		1.0E+03		9.7E-02		
				5.0E-02	P			V	4.38	1	1	No	Butylbenzene, n-	104-51-8					1.0E+03			1.0E+03		3.2E+00		
				1.0E-01	X			V	4.57	1	1	No	Butylbenzene, sec-	135-98-8					2.0E+03			2.0E+03		5.9E+00		
				1.0E-01	X			V	4.11	1	1	Yes	Butylbenzene, tert-	98-06-6					2.0E+03		1.1E+03	6.9E+02		1.6E+00		
				2.0E-02	A			V	0.36	1	1	Yes	Cacodylic Acid	75-60-5					4.0E+02	6.7E+04		4.0E+02				
				1.8E-03	I	1.0E-03	I	A		0.025	1	Yes	Cadmium (Diet)	7440-43-9					1.0E+01	1.1E+02		9.2E+00	5.0E+00	6.9E-01	3.8E-01	
				1.8E-03	I	5.0E-04	I	A		0.05	1	Yes	Cadmium (Water)	7440-43-9					4.0E+02	2.3E+03		3.4E+02				
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M		0.025	1	Yes	Calcium Chromate	13765-19-0	5.0E-02	2.3E-01		4.1E-02								
				5.0E-01	I	2.2E-03	C	V	-0.19	1	1	Yes	Cephalochlorin	106-60-2					1.0E+04	9.0E+05		9.9E+03		2.5E+00		
1.5E-01	C	4.3E-05	C	2.0E-03	I			V	3.8	1	0.9	Yes	Captafol	2425-06-1	6.2E-01	1.7E+00		4.0E-01	4.0E+01	1.5E+02		3.2E+01		7.1E-04		
2.3E-03	C	6.6E-07	C	1.3E-01	I			V	2.8	1	1	Yes	Captan	133-06-2	3.4E+01	3.4E+02		3.1E+01	2.6E+03	3.0E+04		2.4E+03		2.2E-02		
				1.0E-01	I			V	2.36	1	1	Yes	Carbaryl	63-25-2					2.0E+03	2.4E+04		1.8E+03		1.7E+00		
				5.0E-03	I			V	2.32	1	1	Yes	Carbifuran	1563-66-2					1.0E+02	1.4E+03		9.4E+01	4.0E+01	3.7E-02	1.6E-02	
				1.0E-01	I	7.0E-01	I	V	1.94	1	1	Yes	Carbon Disulfide	75-15-0					2.0E+03	2.0E+04	1.5E+03	8.1E+02		2.4E-01		
7.0E-02	I	6.0E-06	I	4.0E-03	I	1.0E-01	I	V	2.83	1	1	Yes	Carbon Tetrachloride	56-23-5	1.1E+00	4.2E+00	9.4E-01	4.5E-01	8.0E+01	3.4E+02	2.1E+02	4.9E+01	5.0E+00	1.8E-04	1.9E-03	
				1.0E-02	I			V	3.81	1	0.8	Yes	Carbosulfan	56285-14-8					2.0E+02	6.9E+01		5.1E+01		1.2E+00		
				1.0E-01	I			V	2.14	1	1	Yes	Carboxin	5234-68-4					2.0E+03	4.1E+04		1.9E+03		1.0E+00		
								V		1	1	Yes	Ceric Oxide	1306-38-3										4.0E-01		
				1.0E-01	I			V	0.99	1	1	Yes	Chloral Hydrate	302-17-0					2.0E+03	1.5E+05		2.0E+03		7.0E-02		
				1.5E-02	I			V	1.9	1	1	Yes	Chlorarbiten	133-90-4					3.0E+02	7.4E+03		2.9E+02		1.5E-04		
4.0E-01	H							V	2.22	1	1	Yes	Chloralid	118-75-2	3.9E-01	3.4E+00		1.8E-01						1.5E-04		
3.5E-01	I	1.0E-04	I	5.0E-04	I	7.0E-04	I	V	6.26	1	0.7	No	Chlordane	12789-03-6	2.2E-01		5.6E-02	4.5E-02	1.0E+01		1.5E+00	1.3E+00	2.0E+00	3.0E-03	1.4E-01	
1.0E+01	I	4.6E-03	C	3.0E-04	I			V	5.41	1	0.8	Yes	Chlordecone (Kepone)	143-50-0	7.8E-03	6.2E-03		3.5E-03	6.0E+00	5.4E+00		2.9E+00		1.2E-04		
				7.0E-04	A			V	3.81	1	0.9	Yes	Chlorfenvinphos	470-90-6					1.4E+01	5.6E+01		1.1E+01		3.1E-02		
				2.0E-02	I			V	2.5	1	1	Yes	Chlorimuron, Ethyl-	90982-32-4					4.0E+02	1.5E+04		3.9E+02		1.3E-01		
				1.0E-01	I	1.5E-04	A	V	0.85	1	1	Yes	Chlorine	7782-50-5					2.0E+03	4.5E+05	3.0E-01	3.0E-01		1.4E-04		
				3.0E-02	I	2.0E-04	I	V		1	1	Yes	Chlorine Dioxide	10049-04-4					6.0E+02	1.4E+05	4.2E-01	4.2E-01				
				3.0E-02	I			V		1	1	Yes	Chlorite (Sodium Salt)	7758-19-2					6.0E+02	1.4E+05		6.0E+02	1.0E+03			
								V	2.05	1	1	Yes	Chloro-1,1-difluoroethane, 1-	75-68-3							1.0E+05	1.0E+05		5.2E+01		
				3.0E-04	I	2.0E-02	H	V	2.53	1	1	Yes	Chloro-1,3-butadiene, 2-	126-99-8			1.9E-02	1.9E-02	4.0E+02	1.8E+03	4.2E+01	3.7E+01		9.8E-06		
4.6E-01	H							V	-1.42	1	1	Yes	Chloro-2-methylaniline HCl, 4-	3165-93-3	1.7E-01	4.9E+02										

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncarcinogenic Child Hazard Index (HI) = 1				Protection of Groundwater SSL			
SFO (mg/kg-day) ¹	k _e	IUR (ug/m ³ -d) ¹	k _e	RD ₅₀ (mg/kg-day)	k _e	RF _c (mg/m ³ -d)	k _e	muta-gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)	
2.4E+02	C	6.9E-02	C	2.0E-01	I	1.0E-03	A		1.02	1	1	Yes	Chlorozotocin	54749-90-5	3.2E-04	7.1E-01		3.2E-04	4.0E+03	9.8E+03		2.8E+03		7.1E-08	2.6E+00	
				1.0E-03	A				4.96	1	0.8	Yes	Chlorpropham	101-21-3					2.0E+01	1.5E+01		8.4E+00		1.2E-01	1.2E-01	
				1.0E-02	H				4.31	1	0.9	Yes	Chlorpyrifos Methyl	5598-13-0					2.0E+02	2.9E+02		1.2E+02		5.4E-01	5.4E-01	
				5.0E-02	I				2	1	1	Yes	Chlorsulfuron	64902-72-3					1.0E+03	5.7E+04		9.9E+02		8.3E-01	8.3E-01	
				8.0E-04	H				5.8	1	0.8	Yes	Chlorthiophos	60238-56-4					1.6E+01	3.4E+00		2.8E+00		7.3E-02	7.3E-02	
5.0E-01	J	8.4E-02	S	1.5E+00	I	3.0E-03	I	1.0E-04	I	M			Chromium(III), Insoluble Salts	16065-83-1	5.0E-02	1.1E-01		3.5E-02	3.0E+04	8.9E+04		2.2E+04		4.0E+07	6.7E-04	
									0.013	1	1	Yes	Chromium(VI)	18540-29-9					6.0E+01	1.7E+02		4.4E+01	1.0E+02		1.8E+05	
									0.013	1	1	Yes	Chromium, Total	7440-47-3												
				9.0E-03	P	3.0E-04	P	6.0E-06	P				Cobalt	7440-48-4					6.0E+00	3.4E+03		6.0E+00		2.7E-01	2.7E-01	
				6.2E-04	I				4.0E-02	H			Coke Oven Emissions	8007-45-2									1.3E+03	2.8E+01	4.6E+01	
													Copper	7440-50-8					8.0E+02	1.8E+05		8.0E+02		2.8E+01	4.6E+01	
				5.0E-02	I	6.0E-01	C		1.96	1	1	Yes	Cresol, m-	108-39-4					1.0E+03	1.2E+04		9.3E+02		7.4E-01	7.4E-01	
				5.0E-02	I	6.0E-01	C		1.95	1	1	Yes	Cresol, o-	95-48-7					1.0E+03	1.2E+04		9.3E+02		7.5E-01	7.5E-01	
				1.0E-01	A	6.0E-01	C		1.94	1	1	Yes	Cresol, p-	106-44-5					2.0E+03	2.5E+04		1.9E+03		1.5E+00	1.5E+00	
				1.0E-01	A				3.1	1	1	Yes	Cresol, p-chloro-m-	59-60-7					2.0E+03	5.2E+03		1.4E+03		1.7E+00	1.7E+00	
				1.0E-01	A	6.0E-01	C		1.95	1	1	Yes	Cresols	1319-77-3					2.0E+03	2.4E+04		1.9E+03		1.5E+00	1.5E+00	
1.9E+00	H			1.0E-03	P				0.6	1	1	Yes	Crotonaldehyde, trans-	123-73-9	4.1E-02	2.6E+00		4.0E-02	2.0E+01	1.5E+03		2.0E+01		8.2E-06	8.2E-06	
				1.0E-01	I	4.0E-01	I	V	3.66	1	1	Yes	Cumene	98-82-8					2.0E+03	1.9E+03	8.3E+02	4.5E+02		7.4E-01	7.4E-01	
2.2E-01	C	6.3E-05	C						-3.16	1	1	No	Cupferron	135-20-6	3.5E-01			3.5E-01						6.1E-04	6.1E-04	
8.4E-01	H			2.0E-03	H				2.22	1	1	Yes	Cyanazine	21725-46-2	9.3E-02	1.5E+00		8.7E-02	4.0E+01	7.5E+02		3.8E+01		4.1E-05	4.1E-05	
				1.0E-03	I					1	1	Yes	Cyanides													
				5.0E-03	I					1	1	Yes	-Calcium Cyanide	592-01-8					2.0E+01	4.5E+03		2.0E+01				
				6.0E-04	I	8.0E-04	S	V		1	1	Yes	-Copper Cyanide	544-92-3					1.0E+02	2.3E+04		1.0E+02		2.0E+02	1.5E-02	
				1.0E-03	I				0.07	1	1	Yes	-Cyanide (CN-)	57-12-5					1.2E+01	2.7E+03	1.7E+00	1.5E+00		2.0E+02	2.0E+00	
				9.0E-02	I					1	1	Yes	-Cyanogen	460-19-5					2.0E+01	5.1E+03		2.0E+01				
				5.0E-02	I					1	1	Yes	-Cyanogen Bromide	506-68-3					1.8E+03	1.6E+06		1.8E+03				
				6.0E-04	I	8.0E-04	I	V		1	1	Yes	-Cyanogen Chloride	506-77-4					1.0E+03	5.8E+05		1.0E+03				
				2.0E-03	I				-0.25	1	1	Yes	-Hydrogen Cyanide	74-90-8					1.2E+01	2.7E+03	1.7E+00	1.5E+00		1.5E-02	1.5E-02	
				5.0E-03	I					0.04	1	Yes	-Potassium Cyanide	151-50-8					4.0E+01	4.5E+03		4.0E+01				
				1.0E-01	I					0.04	1	Yes	-Potassium Silver Cyanide	509-61-6					1.0E+02	4.5E+02		8.2E+01				
				1.0E-03	I					1	1	Yes	-Silver Cyanide	506-64-9					2.0E+03	1.8E+04		1.8E+03				
				2.0E-04	P					1	0	Yes	-Sodium Cyanide	143-33-9					2.0E+01	4.5E+03		2.0E+01	2.0E+02			
				2.0E-04	X					0.58	1	Yes	-Thiocyanates	NA					4.0E+00	9.1E+02		4.0E+00				
				5.0E-02	I					1	1	Yes	-Thiocyanic Acid	463-56-9					4.0E+00	9.1E+02		4.0E+00				
										1	1	Yes	-Zinc Cyanide	557-21-1					1.0E+03	3.8E+05		1.0E+03				
2.3E-02	H			6.0E+00	I	V			3.44	1	1	Yes	Cyclohexane	110-82-7							1.3E+04	1.3E+04		1.3E+01	1.3E+01	
				5.0E+00	I	7.0E-01	P	V	4.72	1	0.9	Yes	Cyclohexane, 1,2,3,4,5-pentabromo-8-chloro-	87-84-3	3.4E+00	8.0E+00		2.4E+00	1.0E+05	6.5E+06	1.5E+03	1.4E+03		1.4E-02	3.4E-01	
				2.0E-01	I				0.81	1	1	Yes	Cyclohexanone	108-94-1					1.0E+02	2.5E+02	2.1E+03	7.0E+01		4.6E-02	4.6E-02	
				5.0E-03	P	1.0E+00	X	V	2.86	1	1	Yes	Cyclohexene	110-83-8					4.0E+03	9.2E+04		3.8E+03		1.0E+00	1.0E+00	
				2.0E-01	I				1.49	1	1	Yes	Cyclohexylamine	108-91-8					1.0E+02			1.0E+02		6.9E+01	6.9E+01	
				5.0E-03	I				6.9	1	0.5	No	Cyhalothrin/karate	68085-85-8					1.0E+02			1.0E+02				
2.4E-01	I	6.9E-05	C	1.0E-02	I				6.6	1	0.7	No	Cypermethrin	52315-07-8					2.0E+02			2.0E+02		3.2E+01	3.2E+01	
3.4E-01	I	9.7E-05	C	7.5E-03	I				0.96	1	1	Yes	Cyromazine	66215-27-8					1.5E+02	1.2E+04		1.5E+02		3.8E-02	3.8E-02	
3.4E-01	I	9.7E-05	I	4.0E-04	A				6.02	1	0.8	Yes	DDD	72-54-8	3.2E-01	3.4E-02		3.1E-02						7.2E-03	7.2E-03	
				1.0E-02	I				6.51	1	0.8	No	DDE, p,p'-	72-55-9			5.8E-02		2.3E-01			4.6E-02		1.1E-02	1.1E-02	
				5.0E-04	I				6.91	1	0.7	No	DDT	50-29-3					2.3E-01			2.3E-01		7.7E-02	7.7E-02	
				1.0E-02	I				4.28	1	0.9	Yes	Dacthal	1861-32-1					2.0E+01	3.2E+02		1.2E+01		1.5E-01	1.5E-01	
7.0E-04	I			3.0E-02	I				0.78	1	1	Yes	Dalapon	75-99-0					6.0E+02	5.5E+04		6.0E+02	2.0E+02	1.2E-01	4.1E-02	
				7.0E-03	I				12.11	1	0	No	Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'-(BDE-209)	1163-19-5	1.1E+02			1.1E+02	1.4E+02			1.4E+02		6.2E+01	6.2E+01	
				4.0E-05	I				3.21	1	0.9	Yes	Demeton	8065-48-3					8.0E-01	4.1E+00		6.7E-01				
1.2E-03	I			6.0E-01	I				8.12	1	0	No	Di(2-ethylhexyl)adipate	103-23-1	6.5E+01			6.5E+01	1.2E+04			1.2E+04	4.0E+02	4.7E+00	2.9E+01	
6.1E-02	H			7.0E-02	X				4.38	1	1	Yes	Diallate	2303-16-4	1.3E+00	8.9E-01		5.2E-01						7.8E-04	7.8E-04	
				4.0E-04	A				3.81	1	0.9	Yes	Diazinon	333-41-5					1.4E+01	3.9E+01		1.0E+01		6.5E-02	6.5E-02	
8.0E-01	P	6.0E-03	P	1.0E-02	X				4.38	1	1	Yes	Dibenzothiophene	132-65-0					2.0E+02	9.6E+01		6.5E+01		1.2E+00	1.2E+00	
				2.0E-04	P	2.0E-04	I	V	2.96	1	1	Yes	Dibromo-3-chloropropane, 1,2-	96-12-8	3.1E-02	1.6E-01	3.4E-04	3.3E-04	4.0E+00	2.4E+01	4.2E-01	3.7E-01	2.0E-01	1.4E-07	8.6E-05	

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where nSL < 100X c SL; ** = where nSL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1				Protection of Groundwater SSL				
SFO (mg/kg-day) ¹	k _e	IUR (ug/m ³ -y) ¹	k _e	RD ₁₀ (mg/kg-day)	k _e	RF _c (mg/m ³ -y)	k _e	muta-gen	LOGP	GIABS	FA	IN EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)	
9.1E-02	I	2.6E-05	I	8.0E-03	X	7.0E-03	P	V	1.48	1	1	Yes	Dichloroethane, 1,2-Dichloroethylene, 1,1-	107-06-2 75-35-4	8.6E-01	1.8E+01	2.2E-01	1.7E-01	1.2E+02 1.0E+03	2.8E+03 8.5E+03	1.5E+01 2.8E+02	1.3E+01 2.8E+02	5.0E+00 7.0E+00	4.8E-05 1.0E-01	1.4E-03 2.5E-03	
				2.0E-03	I			V	1.86	1	1	Yes	Dichloroethylene, 1,2-cis-	156-59-2					4.0E+01	3.6E+02		3.6E+01	7.0E+01	1.1E-02	2.1E-02	
				2.0E-02	I			V	2.09	1	1	Yes	Dichloroethylene, 1,2-trans-	156-60-5					4.0E+02	3.6E+03		3.6E+02	1.0E+02	1.1E-01	3.1E-02	
				3.0E-03	I			V	3.06	1	1	Yes	Dichlorophenol, 2,4-	120-63-2					6.0E+01	1.9E+02		4.6E+01		5.4E-02		
				1.0E-02	I			V	2.81	1	1	Yes	Dichlorophenoxy Acetic Acid, 2,4-	94-75-7					2.0E+02	1.3E+03		1.7E+02	7.0E+01	4.5E-02	1.8E-02	
				8.0E-03	I			V	3.53	1	0.9	Yes	Dichlorophenoxybutyric Acid, 4-(2,4-	94-82-6					1.6E+02	4.8E+02		1.2E+02		1.1E-01		
3.6E-02	C	1.0E-05	C	9.0E-02	A	4.0E-03	I	V	1.96	1	1	Yes	Dichloropropane, 1,2-	78-87-5	2.2E+00	2.3E+01	5.6E-01	4.4E-01	1.8E+03	2.1E+04	8.3E+00	8.3E+00	5.0E+00	1.5E-04	1.7E-03	
				2.0E-02	P			V	2	1	1	Yes	Dichloropropane, 1,3-	142-28-9					4.0E+02	4.2E+03		3.7E+02		1.3E-01		
				3.0E-03	I			V	0.78	1	1	Yes	Dichloropropanol, 2,3-	616-23-9					6.0E+01	4.9E+03		5.9E+01		1.3E-02		
1.0E-01	I	4.0E-06	I	3.0E-02	I	2.0E-02	I	V	2.04	1	1	Yes	Dichloropropene, 1,3-	942-75-6	7.8E-01	7.5E+00	1.4E+00	4.7E-01	6.0E+02	6.5E+03	4.2E+01	3.9E+01		1.7E-04		
2.9E-01	I	8.3E-05	C	5.0E-04	I	5.0E-04	I	V	1.43	1	1	Yes	Dichlorvos	62-73-7	2.7E-01	1.3E+01		2.6E-01	1.0E+01	5.6E+02		9.9E+00		8.1E-05		
				8.0E-02	P	3.0E-04	X	V	3.51	1	1	Yes	Dicyclopentadiene	77-73-6					1.6E+03	3.5E+03	6.3E-01	6.3E-01		2.2E-03		
1.6E+01	I	4.6E-03	I	5.0E-05	I			V	5.4	1	0.8	Yes	Dieldrin	60-57-1	4.9E-03	2.6E-03		1.7E-03	1.0E+00	6.1E-01		3.8E-01		6.9E-05		
				3.0E-04	C			V			0		Diesel Engine Exhaust	NA												
				2.0E-03	P	2.0E-04	P		-1.43	1	1	Yes	Diethanolamine	111-42-2					4.0E+01	8.4E+04		4.0E+01		8.1E-03		
				3.0E-02	P	1.0E-04	P		0.56	1	1	Yes	Diethylene Glycol Monobutyl Ether	112-34-5					6.0E+02	8.6E+04		6.0E+02		1.3E-01		
				6.0E-02	P	3.0E-04	P		-0.54	1	1	Yes	Diethylene Glycol Monoethyl Ether	111-90-0					1.2E+03	7.8E+05		1.2E+03		2.4E-01		
				1.0E-03	P			V	0.05	1	1	Yes	Diethylformamide	617-84-5					2.0E+01	4.2E+03		2.0E+01		4.1E-03		
3.5E+02	C	1.0E-01	C					V	5.07	1	0.9	Yes	Diethylstilbestrol	56-53-1	2.2E-04	6.3E-05		4.9E-05						2.7E-05		
				8.0E-02	I			V	0.65	1	1	Yes	Difenzoquat	43222-48-6					1.6E+03	7.3E+05		1.6E+03				
				2.0E-02	I			V	3.88	1	0.9	Yes	Diffubenzuron	35367-38-5					4.0E+02	1.0E+03		2.9E+02		3.3E-01		
				4.0E+01	I	V		V	0.75	1	1	Yes	Diffuoroethane, 1,1-	75-37-6							8.3E+04	8.3E+04		2.8E+01		
4.4E-02	C	1.3E-05	C					V	3.38	1	1	Yes	Dihydroxofrole	94-58-6	1.8E+00	2.2E+00	4.3E-01	3.0E-01			1.5E+03			3.7E-04		
				7.0E-01	P	V		V	1.52	1	1	Yes	Diisopropyl Ether	108-20-3										3.7E-01		
				8.0E-02	I			V	1.03	1	1	Yes	Diisopropyl Methylphosphonate	1445-75-6					1.6E+03	1.3E+05		1.6E+03		4.5E-01		
				2.0E-02	I			V	-0.17	1	1	Yes	Dimethipin	55290-64-7					4.0E+02	2.4E+05		4.0E+02		8.8E-02		
				2.0E-04	I			V	0.78	1	1	Yes	Dimethoate	60-51-5					4.0E+00	6.4E+02		4.0E+00		9.0E-04		
1.6E+00	P							V	1.81	1	1	Yes	Dimethoxybenzidine, 3,3'-	119-90-4	4.8E-02	1.6E+00		4.7E-02						5.8E-05		
1.7E-03	P			6.0E-02	P			V	-0.61	1	1	Yes	Dimethyl methylphosphonate	756-79-6	4.6E+01	2.7E+04		4.6E+01				1.2E+03	8.1E+05	1.2E+03	9.7E-03	
4.6E+00	C	1.3E-03	C					V	4.58	1	1	Yes	Dimethylamino azobenzene [p-]	60-11-7	1.7E-02	6.9E-03		4.9E-03						2.1E-05		
5.8E-01	H							V	-1.51	1	1	Yes	Dimethylaniline HCl, 2,4-	21436-96-4	1.3E-01	4.0E+02		1.3E-01						1.2E-04		
2.0E-01	P			2.0E-03	X			V	1.68	1	1	Yes	Dimethylaniline, 2,4-	95-88-1	3.9E-01	6.8E+00		3.7E-01	4.0E+01	8.0E+02		3.8E+01		2.1E-04		
				2.0E-03	I			V	2.31	1	1	Yes	Dimethylaniline (N,N-	121-69-7					4.0E+01	3.0E+02		3.5E+01		1.3E-02		
1.1E+01	P							V	2.34	1	1	Yes	Dimethylbenzidine, 3,3'-	119-93-7	1.1E-01	8.2E-02		6.5E-03						4.3E-05		
				1.0E-01	P	3.0E-02	I	V	-1.01	1	1	Yes	Dimethylformamide	68-12-2					2.0E+03	1.8E+06	6.3E+01	6.1E+01		1.2E-02		
				1.0E-04	X	2.0E-06	X	V	-1.19	1	1	Yes	Dimethylhydrazine, 1,1-	57-14-2					2.0E+00	3.5E+03	4.2E-03	4.2E-03		9.3E-07		
5.5E+02	C	1.6E-01	C					V	-0.54	1	1	Yes	Dimethylhydrazine, 1,2-	540-73-8	1.4E-04	4.8E-02	3.5E-05	2.8E-05						6.5E-09		
				2.0E-02	I			V	2.3	1	1	Yes	Dimethylphenol, 2,4-	105-67-9					4.0E+02	3.1E+03		3.6E+02		4.2E-01		
				6.0E-04	I			V	2.36	1	1	Yes	Dimethylphenol, 2,6-	576-26-1					1.2E+01	8.5E+01		1.1E+01		1.3E-02		
				1.0E-03	I			V	2.23	1	1	Yes	Dimethylphenol, 3,4-	95-65-8					2.0E+01	1.7E+02		1.8E+01		2.1E-02		
4.5E-02	C	1.3E-05	C					V	2.58	1	1	Yes	Dimethylvinylchloride	513-37-1	1.7E+00	6.3E+00	4.3E-01	3.3E-01						2.0E-04		
				8.0E-05	X			V	2.13	1	1	Yes	Dinitro- <i>o</i> -cresol; 4,6-	534-52-1					1.6E+00	2.6E+01		1.5E+00		2.6E-03		
				2.0E-03	I			V	4.12	1	0.9	Yes	Dinitro- <i>o</i> -cyclohexyl Phenol, 4,6-	131-89-5					4.0E+01	5.4E+01		2.3E+01		7.7E-01		
				1.0E-04	P			V	1.69	1	1	Yes	Dinitrobenzene, 1,2-	528-29-0					2.0E+00	5.3E+01		1.9E+00		1.8E-03		
				1.0E-04	I			V	1.49	1	1	Yes	Dinitrobenzene, 1,3-	99-65-0					2.0E+00	7.2E+01		2.0E+00		1.8E-03		
				1.0E-04	P			V	1.46	1	1	Yes	Dinitrobenzene, 1,4-	100-25-4					2.0E+00	7.5E+01		2.0E+00		1.8E-03		
				2.0E-03	I			V	1.67	1	1	Yes	Dinitrophenol, 2,4-	51-28-5					4.0E+01	1.2E+03		3.9E+01		4.4E-02		
6.8E-01	I							V	2.18	1	1	Yes	Dinitrotoluene Mixture, 2,4,6-	NA	1.1E-01	1.4E+00		1.1E-01						1.5E-04		
3.1E-01	C	8.9E-05	C	2.0E-03	I			V	1.98	1	1	Yes	Dinitrotoluene, 2,4-	121-14-2	2.5E-01	4.1E+00		2.4E-01	4.0E+01	7.5E+02		3.8E+01		3.2E-04		
1.5E+00	P			3.0E-04	X			V	2.1	1	1	Yes	Dinitrotoluene, 2,6-	606-20-2	5.2E-02	7.1E-01		4.8E-02	6.0E+00	9.3E+01		5.7E+00		6.7E-05		
				2.0E-03	S			V	1.84	1	1	Yes	Dinitrotoluene, 2-Amino-4,6-	35572-78-2					4.0E+01	1.0E+03		3.9E+01		3.0E-02		
				2.0E-03	S			V	1.84	1	1	Yes	Dinitrotoluene, 4-Amino-2,6-	19408-51-0					4.0E+01	1.0E+03		3.9E+01		3.0E-02		
4.5E-01	X			9.0E-04	X			V	2.18	1	1	Yes														

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Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1				Protection of Groundwater SSL					
SFO (mg/kg-day) ¹	k _e	IUR (ug/m ³ -y) ¹	k _e	RD ₅₀ (mg/kg-day)	k _e	RF _c (mg/m ³ -y)	k _e	muta-gen	LOGP	GIABS	FA	IN EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)
9.9E-03	I	1.2E-06	I	6.0E-03	P	1.0E-03	I	V	0.45	1	1	Yes	Epichlorohydrin	106-89-8	7.9E+00	7.5E+02	4.7E+00	2.9E+00	1.2E+02	1.3E+04	2.1E+00	2.0E+00		4.5E-04	9.2E-03
				2.0E-02	I	V			0.86	1	1	Yes	Epoxybutane, 1,2-	106-88-7							4.2E+01				
				5.0E-03	I				-0.22	1	1	Yes	Ethephon	16672-87-0					1.0E+02	4.2E+04		1.0E+02		2.1E-02	
				5.0E-04	I				5.07	1	0.8	Yes	Ethion	563-12-2					1.0E+01	7.7E+00		4.3E+00		8.5E-03	
				1.0E-01	P	6.0E-02	P	V	0.59	1	1	Yes	Ethoxyethanol Acetate, 2-	111-15-9					2.0E+03	2.3E+05	1.3E+02	1.2E+02		2.5E-02	
				9.0E-02	P	2.0E-01	I	V	-0.32	1	1	Yes	Ethoxyethanol, 2-	110-80-5					1.8E+03	6.2E+05	4.2E+02	3.4E+02		6.8E-02	
4.8E-02	H			9.0E-01	I	7.0E-02	P	V	0.73	1	1	Yes	Ethyl Acetate	141-78-6	1.6E+00	4.3E+01		1.6E+00	1.8E+04	1.2E+06	1.5E+02	1.4E+02		3.1E-02	
				5.0E-03	P	8.0E-03	P	V	1.32	1	1	Yes	Ethyl Acrylate	140-68-5					1.0E+02	3.0E+03	1.7E+01	1.4E+01		3.5E-04	
				1.0E+01	I	V			1.43	1	1	Yes	Ethyl Chloride (Chloroethane)	75-00-3					4.0E+03	2.0E+05	2.1E+04	2.1E+04		5.9E+00	
				2.0E-01	I	V			0.89	1	1	Yes	Ethyl Ether	60-29-7					1.8E+03	2.3E+04	6.3E+02	4.6E+02		8.8E-01	
				9.0E-02	H	3.0E-01	P	V	1.94	1	1	Yes	Ethyl Methacrylate	97-63-2					1.8E+03	2.3E+04	6.3E+02	4.6E+02		1.1E-01	
				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		2.8E-03	
1.1E-02	C	2.5E-06	C	1.0E-01	I	1.0E+00	I	V	3.15	1	1	Yes	Ethylbenzene	100-41-4	7.1E+00	1.2E+01	2.2E+00	1.5E+00	2.0E+03	3.8E+03	2.1E+03	8.1E+02	7.0E+02	1.7E-03	7.9E-01
				7.0E-02	P				-0.94	1	1	Yes	Ethylene Cyanohydrin	109-78-4					1.4E+03	1.1E+06		1.4E+03		2.8E-01	
				9.0E-02	P				-2.04	1	1	No	Ethylene Diamine	107-15-3					1.8E+03			1.8E+03		4.2E-01	
				2.0E+00	I	4.0E-01	C	V	-1.36	1	1	Yes	Ethylene Glycol	107-21-1					4.0E+04	5.7E+07		4.0E+04		8.1E+00	
				1.0E-01	I	1.8E+00	I		0.83	1	1	Yes	Ethylene Glycol Monobutyl Ether	111-76-2					2.0E+03	1.4E+05		2.0E+03		4.1E-01	
3.1E-01	C	8.8E-05	C			3.0E-02	C	V	-0.3	1	1	Yes	Ethylene Oxide	75-21-8	2.5E-01	5.2E+01	6.4E-02	5.1E-02			6.3E+01	6.3E+01		1.1E-05	
4.5E-02	C	1.3E-05	C	8.0E-05	I				-0.66	1	1	Yes	Ethylene Thiourea	96-45-7	1.7E+00	9.7E+02		1.7E+00	1.6E+00	1.0E+03		1.6E+00		3.6E-04	
6.5E+01	C	1.9E-02	C						-0.28	1	1	Yes	Ethyleneimine	151-56-4	1.2E-03	2.4E-01	3.0E-04	2.4E-04						5.2E-08	
				3.0E+00	I				2.19	1	1	Yes	Ethylphthalyl Ethyl Glycolate	84-72-0					6.0E+04	1.5E+06		5.8E+04		1.3E+02	
				8.0E-03	I				2.55	1	1	Yes	Express	101200-48-0					1.6E+02	5.0E+03		1.6E+02		6.1E-02	
				2.5E-04	I				3.23	1	0.9	Yes	Fenamiphos	22224-92-6					5.0E+00	3.4E+01		4.4E+00		4.3E-03	
				2.5E-02	I				5.7	1	0.8	Yes	Fenpropathrin	39515-41-8					5.0E+02	7.3E+01		6.4E+01		2.9E+00	
				1.3E-02	I				2.42	1	1	Yes	Fluometuron	2164-17-2					2.6E+02	3.4E+03		2.4E+02		1.9E-01	
				4.0E-02	C	1.3E-02	C			1	1	Yes	Fluoride	16984-48-8					8.0E+02	1.8E+05		8.0E+02		1.2E+02	
				6.0E-02	I	1.3E-02	C			1	1	Yes	Fluorine (Soluble Fluoride)	7782-41-4					1.2E+03	2.7E+05		1.2E+03	4.0E+03	1.8E+02	6.0E+02
				8.0E-02	I				3.16	1	0.9	Yes	Fluridone	59756-60-4					1.6E+03	1.4E+04		1.4E+03		1.6E+02	
				2.0E-02	I				3.34	1	0.9	Yes	Flurprimidol	56425-91-3					4.0E+02	2.4E+03		3.4E+02		1.6E+00	
				6.0E-02	I				3.7	1	0.9	Yes	Flutolanil	86332-96-5					1.2E+03	4.5E+03		9.5E+02		5.0E+00	
				1.0E-02	I				6.81	1	0.6	No	Fluvalinate	69409-94-5					2.0E+02			2.0E+02		2.9E+02	
3.5E-03	I			1.0E-01	I				2.85	1	1	Yes	Folpet	133-07-3	2.2E+01	2.0E+02		2.0E+01	2.0E+03	2.1E+04		1.8E+03		4.7E-03	
1.9E-01	I			2.0E-03	I				2.9	1	1	Yes	Fomesafen	72178-02-0	4.1E-01	8.7E+00		3.9E-01	4.0E+01	6.3E+01	2.0E+01	2.4E+01		1.3E-03	
				1.3E-05	I	9.8E-03	A	V	0.35	1	1	Yes	Formaldehyde	50-00-0			4.3E-01	4.3E-01	4.0E+03	3.2E+05	2.0E+01	2.0E+01		4.7E-02	8.7E-05
				9.0E-01	P	3.0E-04	X	V	-0.54	1	1	Yes	Formic Acid	64-18-8					1.8E+04	6.3E+06	6.3E-01	6.3E-01		1.3E-04	
				3.0E+00	I				-2.4	1	1	No	FosetylAL	39348-24-8					6.0E+04			6.0E+04			
				1.0E-03	X				4.12	1	1	Yes	Furans						2.0E+01	1.3E+01		7.9E+00		1.5E-01	
				1.0E-03	I				1.34	1	1	Yes	~Furan	110-00-9					2.0E+01	4.8E+02		1.9E+01		7.3E-03	
				9.0E-01	I	2.0E+00	I	V	0.46	1	1	Yes	~Tetrahydrofuran	109-99-9					1.8E+04	1.7E+06	4.2E+03	3.4E+03		7.5E-01	
3.8E+00	H			3.0E-03	I	5.0E-02	H	V	-0.04	1	1	Yes	Furazolidone	67-45-8	2.1E-02	9.8E+00		2.0E-02	6.0E+01	7.1E+03	1.0E+02	3.8E+01		3.9E-05	
				1.5E+00	C	4.3E-04	C		0.41	1	1	Yes	Furfural	98-01-1					2.0E+01	1.8E+00		5.0E-02		8.1E-03	
				3.0E-02	I	8.6E-06	C		1.8	1	1	Yes	Furium	531-82-8	5.2E-02	1.8E+00		5.0E-02	6.0E+01	7.1E+03	1.0E+02	3.8E+01		6.8E-05	
				4.0E-04	I				4.38	1	0.9	Yes	Furmecyclo	60568-05-0	2.6E+00	1.9E+00	1.1E+00		6.0E+01	7.1E+03	1.0E+02	3.8E+01		1.2E-03	
				8.0E-05	C				-5.34	1	1	No	Glutamate, Ammonium	77182-82-2					8.0E+00			8.0E+00		1.8E-03	
				1.0E-02	X				-0.18	1	1	Yes	Glutaraldehyde	111-30-8					8.0E+00			8.0E+00		1.8E-03	
				1.0E-01	I				-0.12	1	1	Yes	Glycidyl	765-34-4					8.0E+00	1.8E+03	2.1E+00	1.7E+00		3.3E-04	
				3.0E-03	I				-3.4	1	1	No	Glyphosate	1071-83-6					2.0E+03			2.0E+03	7.0E+02	8.8E+00	3.1E+00
				1.0E-02	X				4.73	1	0.8	Yes	Goal	42874-03-3					6.0E+01	6.6E+01		3.2E+01		2.5E+00	
				2.0E-02	P				-1.63	1	1	Yes	Guanidine	113-00-8					2.0E+02	4.2E+05		2.0E+02		4.5E-02	
				3.0E-03	A	1.0E-02	A		-1.7	1	1	Yes	Guanidine Chloride	50-01-1					4.0E+02	1.0E+09		4.0E+02		1.7E-02	
				5.0E-05	I				2.75	1	1	Yes	Gutnion	86-90-0					6.0E+01	8.3E+02		5.6E+01		1.7E-02	
				1.3E-02	I				4.07	1	0.9	Yes	Haloxypol, Methyl	69806-40-2					1.0E+00	3.1E+00		7.6E-01		8.4E-03	
4.5E+00	I	1.3E-03	I	5.0E-04	I				1.56	1	1	Yes	Harmony	79277-27-3					2.6E+02	3.5E+04		2.6E+02		7.8E-02	
				1.3E-05	I				6.1	1	0.8	Yes	Heptachlor	76											

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncarcinogenic Child Hazard Index (HI) = 1				Protection of Groundwater SSL		
SFO (mg/kg-day)	k _e	IUR (ug/m ³ -y)	k _e	RD ₅₀ (mg/kg-day)	k _e	RF _c (mg/m ³ -y)	k _e	muta-gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)
				5.0E-03		3.0E-02	I	V	1.38	1	1	Yes	Hexanone, 2-	591-78-6					1.0E+02	2.7E+03	6.3E+01	3.8E+01		8.8E-03	
				3.3E-02			I		1.85	1	1	Yes	Hexazinone	51235-04-2					6.6E+02	2.4E+04		6.4E+02		3.0E-01	
3.0E+00	I	4.9E-03	I			3.0E-05	P	V	-2.07	1	1	Yes	Hydrazine	302-01-2	2.6E-02	1.1E+02	1.1E-03	1.1E-03			6.3E-02	6.3E-02			
3.0E+00	I	4.9E-03	I							1	1	Yes	Hydrazine Sulfate	10034-93-2	2.6E-02	4.7E+00		2.6E-02							
				4.0E-02		2.0E-02	I	V	0.23	1	1	Yes	Hydrogen Chloride	7647-01-0					8.0E+02	1.8E+05	4.2E+01	4.2E+01			
				1.4E-02			C	V	0.23	1	1	Yes	Hydrogen Fluoride	7664-39-3							2.9E+01	2.8E+01			
				2.0E-03			I	V	0.23	1	1	Yes	Hydrogen Sulfide	7783-06-4							4.2E+00	4.2E+00			
6.0E-02	P			4.0E-02	P				0.59	1	1	Yes	Hydroquinone	123-31-9	1.3E+00	1.1E+02		1.3E+00	8.0E+02	7.9E+04		7.9E+02		8.7E-04	
				1.3E-02	I				3.82	1	0.9	Yes	Imazali	35554-44-0					2.6E+02	6.8E+02		1.9E+02		3.2E+00	
				2.5E-01	I				1.86	1	1	Yes	Imazaquin	81335-37-7					5.0E+03	2.6E+05		4.9E+03		2.5E+01	
				1.0E-02	A					1	1	Yes	Iodine	7553-56-2					2.0E+02	4.5E+04		2.0E+02		1.2E+01	
				4.0E-02	I				3	1	0.9	Yes	Iprodione	36734-19-7					8.0E+02	9.1E+03		7.4E+02		2.3E-01	
				7.0E-01	P					1	1	Yes	Iron	7439-89-6					1.4E+04	3.2E+06		1.4E+04		3.5E+02	
				3.0E-01	I		V		0.76	1	1	Yes	Isobutyl Alcohol	78-83-1					6.0E+03	3.6E+05		5.9E+03		1.2E+00	
				2.0E-01	I	2.0E+00	C		1.7	1	1	Yes	Isophorone	78-59-1	8.2E+01	1.6E+03		7.8E+01	4.0E+03	8.6E+04		3.8E+03		2.6E-02	
				1.5E-02	I		V		5.8	1	0.8	Yes	Isopropalin	33820-53-0					3.0E+02	4.6E+01		4.0E+01		9.2E-01	
				2.0E+00	P	2.0E-01	P	V	0.05	1	1	Yes	Isopropanol	67-63-0					4.0E+04	6.5E+06	4.2E+02	4.1E+02		8.4E-02	
				1.0E-01	I				0.27	1	1	Yes	Isopropyl Methyl Phosphonic Acid	1832-54-8					2.0E+03	3.9E+05		2.0E+03		4.3E-01	
				5.0E-02	I				3.94	1	0.9	Yes	Isoxaben	82558-50-7					1.0E+03	2.7E+03		7.3E+02		2.0E+00	
						3.0E-01	A	V	8	1	0	No	JP-7	NA							6.3E+02	6.3E+02		1.2E+00	
				7.5E-02	I				3.43	1	0.9	Yes	Kerb	23950-58-5					1.5E+03	5.5E+03		1.2E+03		1.2E+00	
				2.0E-03	I				4.81	1	0.9	Yes	Lactofen	77501-63-4					4.0E+01	6.7E+01		2.5E+01		1.2E+00	
													Lead Compounds												
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M	0.025	1	0.8	Yes	~Lead Chromate	7758-97-6	5.0E-02	2.3E-01		4.1E-02	4.0E+02	2.3E+03		3.4E+02			
8.5E-03	C	1.2E-05	C									Yes	~Lead Phosphate	7446-27-7	9.2E+00	1.6E+03		9.1E+00							
2.8E-01	C	8.0E-05	C						-0.08	1	1	Yes	~Lead acetate	301-04-2	2.8E-01	2.7E+02		2.8E-01				1.5E+01	1.5E+01	1.4E+01	
8.5E-03	C	1.2E-05	C									No	~Lead subacetate	1335-32-6	9.2E+00			9.2E+00							
				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		4.7E-06	
				2.0E-03	I				3.2	1	0.9	Yes	Linuron	330-55-2					4.0E+01	2.0E+02		3.3E+01		2.9E-02	
				2.0E-03	P					1	1	Yes	Lithium	7439-93-2					4.0E+01	9.1E+03		4.0E+01		1.2E+01	
				2.0E-01	I				2.18	1	1	Yes	Londax	83055-99-6					4.0E+03	2.4E+05		3.9E+03		1.0E+00	
				5.0E-04	I				3.25	1	1	Yes	MCPA	94-74-6					1.0E+01	3.0E+01		7.5E+00		2.0E-03	
				1.0E-02	I				3.5	1	0.9	Yes	MCPB	94-81-5					2.0E+02	5.5E+02		1.5E+02		5.8E-02	
				1.0E-03	I				3.13	1	1	Yes	MCPP	93-65-2					2.0E+01	7.1E+01		1.6E+01		4.7E-03	
				2.0E-02	I				2.36	1	1	Yes	Malathion	121-75-5					4.0E+02	1.1E+04		3.9E+02		1.0E-01	
				1.0E-01	I	7.0E-04	C		1.62	1	1	Yes	Maleic Anhydride	108-31-6					2.0E+03	3.8E+04		1.9E+03		3.9E-01	
				5.0E-01	I				-0.84	1	1	Yes	Maleic Hydrate	123-33-1					1.0E+04	8.9E+06		1.0E+04		2.1E+00	
				1.0E-04	P				-0.6	1	1	Yes	Malorohitrite	109-77-3					2.0E+00	9.1E+02		2.0E+00		4.1E-04	
				3.0E-02	H				1.33	1	0.9	Yes	Mancozeb	8018-01-7					6.0E+02	4.9E+03		5.4E+02		2.0E-01	
				5.0E-03	I				0.62	1	1	Yes	Maneb	12427-38-2					1.0E+02	4.4E+03		9.8E+01		1.4E-01	
				1.4E-01	I	5.0E-05	I			1	1	Yes	Manganese (Diet)	7439-96-5											
				2.4E-02	S	5.0E-05	I		0.04	1	1	Yes	Manganese (Non-diet)	7439-96-5					4.8E+02	4.4E+03		4.3E+02		2.8E+01	
				9.0E-05	H				1.04	1	1	Yes	Mepiquat Chloride	950-10-7					1.8E+00	2.5E+02		1.8E+00		2.6E-03	
				3.0E-02	I				-2.82	1	1	No	Mercury Compounds	24307-26-4					6.0E+02			6.0E+02		2.0E-01	
				3.0E-04	I	3.0E-04	S		-0.22	0.07	1	Yes	~Mercuric Chloride (and other Mercury Salts)	7487-84-7					6.0E+00	9.5E+01		5.7E+00	2.0E+00	3.3E-02	1.0E-01
						3.0E-04	I	V	0.62	1	1	Yes	~Mercury (elemental)	7439-97-6							6.3E-01	6.3E-01	2.0E+00		
				1.0E-04	I					1	1	Yes	~Methyl Mercury	22967-92-6					2.0E+00	4.5E+02		2.0E+00			
				8.0E-05	I				0.71	1	1	Yes	~Phenylmercuric Acetate	62-38-4					1.6E+00	5.7E+02		1.6E+00		5.0E-04	
				3.0E-05	I		V		7.67	1	0.3	No	Merphos	150-50-5					6.0E-01			6.0E-01		5.9E-02	
				3.0E-05	I				5.7	1	0.9	Yes	Merphos Oxide	78-48-8					6.0E-01	9.9E-02		8.5E-02		4.2E-04	
				6.0E-02	I				1.65	1	1	Yes	Metalaxyl	67837-19-1					1.2E+03	6.4E+04		1.2E+03		3.3E-01	
				1.0E-04	I	3.0E-02	P	V	0.68	1	1	Yes	Methacrylonitrile	126-98-7					2.0E+00	1.3E+02	6.3E+01	1.9E+00		4.4E-04	
				5.0E-05	I				-0.8	1	1	Yes	Methamidophos	10265-92-6					1.0E+00	1.0E+03		1.0E+00		2.1E-04	
				2.0E+00	I	2.0E+01	I	V	-0.77	1	1	Yes	Methanol	67-56-1					4.0E+04	1.8E+07	4.2E+04	2.0E+04		4.1E+00	
				1.0E-03	I				2.2	1	1	Yes	Methidathion	950-37-8					2.0E+01	5.8E+02		1.9E+01		4.7E-03	
				2.5E-02	I				0.6	1	1	Yes	Methylol	16752-77-5					5.0E+02	6.8E+04		5.0E+02		1.1E-01	
4.9E-02	C	1.4E-05	C						1.47	1	1	Yes	Methoxy-5-nitroaniline, 2-	99-59-2	1.6E+00	5.2E+01		1.5E+00	1.0E+02	5.9E+01		3.7E+01	4.0E+01	5.3E-04	2.0E+00
				5.0E-03	I																				

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1				Protection of Groundwater SSL						
SFO (mg/kg-day) ¹	k	IUR (ug/m ³ -y)	RfD (mg/kg-day)	k	RfC (mg/m ³ -y)	k	y	muta-gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)	
1.8E-03	C	2.6E-07			3.0E+01		V		0.94	1	1	Yes	Methyl tert-Butyl Ether (MTBE)	1634-04-4	4.3E+01	1.9E+03	2.2E+01	1.4E+01			6.3E+03	6.3E+03			3.2E-03	
9.0E-03	P		3.0E-04	X					-2.06	1	1	Yes	Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2	8.7E+00	1.4E+02		8.1E+00	6.0E+00	5.9E+04		6.0E+00		3.6E-03		
8.3E+00	C	2.4E-03	2.0E-02	X					1.87	1	1	Yes	Methyl-5-Nitroaniline, 2-	99-55-8					4.0E+02	7.3E+03		3.8E+02		4.5E-03		
1.3E-01	C	3.7E-05							-0.92	1	1	Yes	Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	9.4E-03	1.0E+01		9.4E-03						3.2E-06		
			1.0E-02	A					-2.06	1	1	Yes	Methylaniline Hydrochloride, 2-	636-21-5	6.0E-01	3.7E+03		6.0E-01						2.6E-04		
									-1.18	1	1	Yes	Methylarsonic acid	124-58-3					2.0E+02	3.6E+05		2.0E+02				
1.0E-01	X		2.0E-04	X						1	0	No	Methylbenzene, 1,4-diamine monohydrochloride, 2-	74612-12-7	7.8E-01			7.8E-01	4.0E+00			4.0E+00				
2.2E+01	C	6.3E-03							6.42	1	0.8	No	Methylbenzene-1,4-diamine sulfate, 2-	615-50-9					6.0E+00			6.0E+00			2.2E-03	
2.0E-03	I	1.0E-08	6.0E-03	I	6.0E-01	V	M	1.25	1	1	1	Yes	Methylene Chloride	75-09-2	1.3E+01	3.4E+02	2.0E+02	1.1E+01	1.2E+02	3.7E+03	1.3E+03	1.1E+02	5.0E+00	2.9E-03	1.3E-03	
1.0E-01	P	4.3E-04	2.0E-03	P					3.91	1	0.9	Yes	Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	2.5E-01	4.2E-01		1.6E-01	4.0E+01	7.5E+01		2.6E+01		1.8E-03		
4.6E-02	I	1.3E-05							4.37	1	1	Yes	Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	1.7E+00	6.4E-01		4.6E-01						2.6E-03		
1.6E+00	C	4.6E-04			2.0E-02	C		1.59	1	1	1	Yes	Methylenebisbenzamide, 4,4'-	101-77-9	4.9E-02	1.6E+00		4.7E-02							2.1E-04	
			7.0E-02	H	6.0E-04	I	V	5.22	1	0.9	Yes	Methylenediphenyl Diisocyanate	101-68-8													
									3.48	1	1	Yes	Methylstyrene, Alpha-	98-83-9					1.4E+03	1.7E+03		7.8E+02			1.3E+00	
			1.5E-01	I					5.13	1	1	Yes	Metolachlor	51218-45-2					3.0E+03	2.6E+04		2.7E+03			3.2E+00	
			2.5E-02	I					1.7	1	1	Yes	Metribuzin	21087-64-9					5.0E+02	1.8E+04		4.9E+02			1.5E-01	
			3.0E+00	P					6.1	1	1	No	Mineral oils	8012-95-1					6.0E+04	2.6E+04		6.0E+04			2.4E+03	
1.8E+01	C	5.1E-03	2.0E-04	I		V		6.89	1	0.5	No	Mirex	2385-85-5	4.3E-03		1.1E-03	8.8E-04	4.0E+00				4.0E+00			6.3E-04	
			2.0E-03	I					3.21	1	1	Yes	Molinate	2212-67-1				4.0E+01	1.2E+02			3.0E+01			1.7E-02	
			5.0E-03	I						1	1	Yes	Molybdenum	7439-98-7				1.0E+02	2.3E+04		1.0E+02			2.0E+00		
			1.0E-01	I						1	1	Yes	Monochloramine	10599-90-3				2.0E+03	4.5E+05			2.0E+03	4.0E+03			
			2.0E-03	P					1.66	1	1	Yes	Monomethylaniline	100-61-8				4.0E+01	7.5E+02			3.8E+01			1.4E-02	
			3.0E-04	X					4.04	1	0.9	Yes	N,N-Diphenyl-1,4-benzenediamine	74-31-7				6.0E+00	8.8E+00			3.6E+00			3.7E-01	
			2.0E-03	I		V		1.38	1	1	1	Yes	Naled	300-76-5				4.0E+01	6.8E+03		2.1E+02	4.0E+01			1.8E-02	
1.8E+00	C	0.0E+00	3.0E-02	X	1.0E-01	P	V	2.28	1	0	No	Naphtha, High Flash Aromatic (HFAN)	64742-95-6	4.3E-02	3.5E-01		3.9E-02	6.0E+02				1.5E+02			2.0E-04	
			1.0E-01	I					3.36	1	0.9	Yes	Napropamide	15299-99-7				2.0E+03	8.9E+03			1.6E+03			1.1E+01	
			2.6E-04	C	1.1E-02	C	1.4E-05	C		1	1	Yes	Nickel Acetate	373-02-4				2.2E+02	2.6E+05			2.2E+02				
			2.6E-04	C	1.1E-02	C	1.4E-05	C		1	1	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		0	Yes	Nickel Carbonyl	13463-39-3			2.2E-02	2.2E-02	2.2E+02		2.9E-02	2.9E-02				
			2.6E-04	C	1.1E-02	C	1.4E-05	C		0.04	1	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	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	Yes	Nickel Refinery Dust	NA				2.2E+02	1.0E+04			2.2E+02			3.2E+01	
			2.6E-04	C	2.0E-02	C	1.4E-05	A		0.04	1	Yes	Nickel Soluble Salts	7440-02-0				4.0E+02	1.8E+04			3.9E+02			2.6E+01	
1.7E+00	C	4.8E-04	1.1E-02	C	1.4E-05	C		0.04	1	1	Yes	Nickel Sulfide	12035-72-2	4.6E-02	1.6E+00		4.5E-02	2.2E+02	1.0E+04			2.2E+02				
			2.6E-04	C	1.1E-02	C	1.4E-05	C		1	0	Yes	Nickelocene	1271-28-9				2.2E+02				2.2E+02				
			1.6E+00	I						1	1	Yes	Nitrate	14797-55-0				3.2E+04	7.3E+06			3.2E+04	1.0E+04			
										1	0	Yes	Nitrate + Nitrite (as N)	NA								1.0E+04				
			1.0E-01	I						1	1	Yes	Nitrite	14797-55-0				2.0E+03	4.5E+05			2.0E+03				
			1.0E-02	X	5.0E-05	X		1.85	1	1	1	Yes	Nitroaniline, 2-	88-74-4				2.0E+02	3.4E+03			1.9E+02			8.0E-02	
2.0E-02	P		4.0E-03	P	6.0E-03	P		1.39	1	1	1	Yes	Nitroaniline, 4-	100-01-8	3.9E+00	1.2E+02		8.0E+01	2.8E+03			7.8E+01			1.6E-03	
			2.0E-03	I	9.0E-03	I	V	1.85	1	1	1	Yes	Nitrobenzene	98-95-3			1.4E-01	1.4E-01	4.0E+01	6.2E+02	1.9E+01	1.3E+01			9.2E-05	
			3.0E+03	P				-4.56	1	1	1	No	Nitrocellulose	9004-70-0				6.0E+07				6.0E+07			1.3E+04	
			7.0E-02	H				-0.47	1	1	1	Yes	Nitrofurantoin	67-20-9				1.4E+03	1.6E+06			1.4E+03			6.1E-01	
1.3E+00	C	3.7E-04						0.23	1	1	1	Yes	Nitrofurazone	59-87-0	6.0E-02	1.6E+01		6.0E-02							5.4E-05	
1.7E-02	P		1.0E-04	P				1.62	1	1	1	Yes	Nitroglycine	55-63-0	4.6E+00	1.8E+02		4.5E+00	2.0E+00	8.7E+01		2.0E+00			8.5E-04	
			1.0E-01	I				-0.89	1	1	1	Yes	Nitroguanidine	656-88-7					2.0E+03	1.8E+06		2.0E+03			4.8E-01	
			8.8E-06	P	5.0E-03	P	V	-0.35	1	1	1	Yes	Nitroethane	75-52-5			6.4E-01	6.4E-01			1.0E+01	1.0E+01			1.4E-04	
			2.7E-03	H	2.0E-02	I	V	0.93	1	1	1	Yes	Nitropropane, 2-	79-46-9			2.1E-03	2.1E-03			4.2E+01	4.2E+01			5.5E-07	
2.7E+01	C	7.7E-03						0.23	1	1	1	Yes	Nitroso-N-ethylurea, N-	759-73-9	9.3E-04	1.5E-01		9.2E-04							2.2E-07	
1.2E+02	C	3.4E-02						-0.03	1	1	1	Yes	Nitroso-N-methylurea, N-	884-93-5	2.1E-04	4.5E-02		2.1E-04							4.6E-08	
5.4E+00	I	1.6E-03						2.63	1	1	1	Yes	Nitroso-di-N-butylamine, N-	924-18-3	1.4E-02	7.6E-02	3.5E-03	2.7E-03							5.5E-06	
7.0E+00	I	2.0E-03						1.36	1	1	1	Yes	Nitroso-di-N-propylamine, N-	621-64-7	1.1E-02	3.4E-01		1.1E-02							8.1E-06	
2.8E+00	I	8.0E-04						-1.28	1	1	1	Yes	Nitrosodiethanolamine, N-	1116-54-7	2.8E-02	7.8E+01		2.8E-02							5.6E-06	
1.5E+02	I	4.3E-02						0.48	1	1	1	Yes	Nitrosodiethylamine, N-	55-18-5	1.7E-04	1.6E-02										

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncarcinogenic Child Hazard Index (HI) = 1		Protection of Groundwater SSL					
SFO (mg/kg-day) ¹	k _e (y)	IUR (ug/m ³ -y)	k _e (mg/kg-day)	RD ₅₀ (mg/m ³)	k _e (mg/m ³)	RF _c (y)	k _e (y)	muta-gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)	
				5.0E-03	I				4.8	1	0.8	Yes	Oxadiazon	19666-30-9					1.0E+02	9.0E+01		4.7E+01		4.8E-01		
				2.5E-02	I				0.47	1	1	Yes	Oxamyl	23135-22-0					5.0E+02	5.0E+02		5.0E+02	2.0E+02	1.1E-01	4.4E-02	
				1.3E-02	I				3.2	1	0.9	Yes	Paclitaxel	76738-62-0					2.6E+02	1.7E+03		2.3E+02		4.6E-01		
				4.5E-03	I				4.5	1	1	No	Paraquat Dichloride	1910-42-5					9.0E+01			9.0E+01		1.2E+00		
				6.0E-03	H				3.83	1	0.9	Yes	Parathion	56-38-2					1.2E+02	3.0E+02		8.6E+01		4.3E-01		
				5.0E-02	H		V		3.83	1	1	Yes	Pebulate	1114-71-2					1.0E+03	1.3E+03		5.6E+02		4.5E-01		
				4.0E-02	I				5.18	1	0.9	Yes	Pendimethalin	40487-42-1					8.0E+02	2.3E+02		1.8E+02		2.1E+00		
				2.0E-03	I				6.84	1	0.6	No	Pentabromodiphenyl Ether	32534-81-9					4.0E+01			4.0E+01		1.8E+00		
				1.0E-04	I				7.66	1	0.6	No	Pentabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-99)	60348-60-9					2.0E+00			2.0E+00		8.7E-02		
				8.0E-04	I		V		5.17	1	0.9	Yes	Pentachlorobenzene	608-93-5					1.6E+01	3.9E+00		3.2E+00		2.4E-02		
9.0E-02	P								3.22	1	1	Yes	Pentachloroethane	76-01-7	8.7E-01	2.4E+00		6.4E-01						3.1E-04		
2.6E-01	H								4.64	1	0.9	Yes	Pentachloronitrobenzene	82-68-8	3.0E-01	1.9E-01		1.2E-01						1.4E-03		
4.0E-01	I	5.1E-06	C						5.12	1	0.9	Yes	Pentachlorophenol	87-86-5	1.9E-01	5.0E-02		4.0E-02					1.0E+00	4.0E-04	1.0E-02	
4.0E-03	X								2.38	1	1	Yes	Pentaerythritol tetranitrate (PETN)	78-11-5	1.9E+01	4.1E+02		1.9E+01				2.9E+01		2.8E-02		
									3.39	1	1	Yes	Pentane, n-	109-66-0							2.1E+03		2.1E+03		1.0E+01	
													Perchlorates													
				7.0E-04	I					1	1	Yes	~Ammonium Perchlorate	7790-98-9					1.4E+01	3.2E+03		1.4E+01				
				7.0E-04	I					1	1	Yes	~Lithium Perchlorate	7791-03-9					1.4E+01	3.2E+03		1.4E+01				
				7.0E-04	I					1	1	Yes	~Perchlorate and Perchlorate Salts	14797-73-0					1.4E+01	3.2E+03		1.4E+01	1.5E+01(F)			
				7.0E-04	I					1	1	Yes	~Potassium Perchlorate	7778-74-7					1.4E+01	1.6E+03		1.4E+01				
				7.0E-04	I					1	1	Yes	~Sodium Perchlorate	7601-89-0					1.4E+01	3.2E+03		1.4E+01				
				2.0E-02	P		V		1.8173	1	1	Yes	Perfluorobutane Sulfonate	375-73-5					4.0E+02	8.3E+03		3.8E+02		2.1E-01		
2.2E-03	C	6.3E-07	C						6.5	1	0.6	No	Permethrin	52645-53-1	3.5E+01	1.1E+03		3.4E+01				1.0E+03		2.4E+02		
									1.58	1	1	Yes	Phenacetin	62-44-2											9.7E-03	
				2.5E-01	I				3.59	1	0.9	Yes	Phenmedipham	13684-63-4					5.0E+03	1.9E+04		4.0E+03		2.1E+01		
				3.0E-01	I	2.0E-01	C		1.46	1	1	Yes	Phenol	108-95-2					6.0E+03	1.4E+05		5.8E+03		3.3E+00		
				5.0E-04	X				4.15	1	1	Yes	Phenothiazine	92-84-2					1.0E+01	7.5E+00		4.3E+00		1.4E-02		
				6.0E-03	I				-0.33	1	1	Yes	Phenylenediamine, m-	108-45-2					1.2E+02	4.8E+04		1.2E+02		3.2E-02		
4.7E-02	H								0.15	1	1	Yes	Phenylenediamine, o-	95-54-5	1.7E+00	2.8E+02		1.6E+00						4.4E-04		
				1.9E-01	H				-0.3	1	1	Yes	Phenylenediamine, p-	106-50-3					3.8E+03	1.4E+06		3.8E+03		1.0E+00		
1.9E-03	H								3.09	1	1	Yes	Phenylphenol, 2-	90-43-7	4.0E+01	1.1E+02		3.0E+01				1.2E+01		4.0E-01		
				2.0E-04	H				3.56	1	0.9	Yes	Phorate	298-02-2								3.0E+00		3.4E-03		
									-0.71	1	1	Yes	Phosgene	75-44-5							2.1E+03		2.1E+03		1.0E+01	
				2.0E-02	I				2.78	1	1	Yes	Phosmet	732-11-6					4.0E+02	5.3E+03		3.7E+02		8.2E-02		
				4.9E+01	P					1	1	Yes	Phosphates, Inorganic													
				4.9E+01	P					1	0	Yes	~Aluminum metaphosphate	13776-88-0					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	0	Yes	~Ammonium polyphosphate	68333-79-9					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Calcium pyrophosphate	7790-76-3					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Diammonium phosphate	7783-28-0					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Dicalcium phosphate	757-83-9					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Dimagnesium phosphate	7782-75-4					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Dipotassium phosphate	7758-11-4					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Disodium phosphate	7558-79-4					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Monobutylammonium phosphate	13530-50-2					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Monoammonium phosphate	7722-76-1					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Monocalcium phosphate	7758-23-8					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Monomagnesium phosphate	7757-86-0					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Monopotassium phosphate	7778-77-0					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Monosodium phosphate	7558-80-7					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Polyphosphoric acid	8017-16-1					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	0.9	Yes	~Potassium triphosphate	13845-36-8					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Sodium acid pyrophosphate	7758-16-9					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Sodium aluminum phosphate (acidic)	7785-88-8					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	0	Yes	~Sodium aluminum phosphate (anhydrous)	10279-59-1					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	0.8	Yes	~Sodium aluminum phosphate (tetrahydrate)	10305-76-7					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	0.9	Yes	~Sodium hexametaphosphate	10124-56-8					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Sodium polyphosphate	68915-31-1					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Sodium trimetaphosphate	7785-84-4					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Sodium tripolyphosphate	7758-29-4					9.7E+05	2.2E+08		9.7E+05				
				4.9E+01	P					1	1	Yes	~Tetrapotassium phosphate													

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncarcinogenic Child Hazard Index (HI) = 1		Protection of Groundwater SSL				
SFO (mg/kg-day) ¹	k _e	IUR (ug/m ³ -y)	k _e	RD ₅₀ (mg/kg-day)	k _e	RF _c (mg/m ³ -y)	k _e	muta-gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)
1.0E-01	I		I						4.5	1	0.9	Yes	~Dibutyl Phthalate	84-74-2					2.0E+03	1.6E+03		9.0E+02		2.3E+00	
8.0E-01	I		I						2.42	1	1	Yes	~Diethyl Phthalate	84-66-2					1.6E+04	2.0E+05		1.5E+04		6.1E+00	
1.0E-01	I		V						2.25	1	1	Yes	~Dimethylterephthalate	120-61-6					2.0E+03	2.7E+04		1.9E+03		4.9E-01	
1.0E-02	P								8.1	1	0	No	~Octyl Phthalate, di-N-	117-84-0					2.0E+02			2.0E+02		5.7E+01	
1.0E+00	H								2	1	1	Yes	~Phthalic Acid, P-	100-21-0					2.0E+04	3.3E+05		1.9E+04		6.8E+00	
2.0E+00	I	2.0E-02	C						1.6	1	1	Yes	~Phthalic Anhydride	85-44-9					4.0E+04	1.1E+06		3.9E+04		8.5E+00	
7.0E-02	I								1.9	1	1	Yes	Picloram	1918-02-1					1.4E+03	4.3E+04		1.4E+03	5.0E+02	3.8E-01	1.4E-01
1.0E-04	X								0.93	1	1	Yes	Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3					2.0E+00	2.1E+02		2.0E+00		1.3E-03	
1.0E-02	I								4.2	1	0.9	Yes	Phosphorus, Methyl	2932-93-7					2.0E+02	3.1E+02		1.2E+02		1.2E-01	
3.0E+01	C	8.6E-03	C	7.0E-06	H					1	0	No	Polybrominated Biphenyls	59536-65-1	2.6E-03			2.6E-03	1.4E-01			1.4E-01			
7.0E-02	S	2.0E-05	S	7.0E-05	I				5.69	1	0.9	No	Polychlorinated Biphenyls (PCBs)	12674-11-2	1.1E+00		2.8E-01	2.2E-01	1.4E+00			1.4E+00		2.1E-02	
2.0E+00	S	5.7E-04	S						4.65	1	1	Yes	~Aroclor 1221	11104-28-2	3.9E-02	1.1E-02	9.8E-03	4.6E-03						7.9E-05	
2.0E+00	S	5.7E-04	S						4.4	1	1	Yes	~Aroclor 1232	11141-16-5	3.9E-02	1.1E-02	9.8E-03	4.6E-03						7.9E-05	
2.0E+00	S	5.7E-04	S						6.34	1	0.7	No	~Aroclor 1242	53469-21-9	3.9E-02		9.8E-03	7.8E-03						1.2E-03	
2.0E+00	S	5.7E-04	S						6.2	1	0.7	No	~Aroclor 1248	12672-29-6	3.9E-02		9.8E-03	7.8E-03				4.0E-01		2.1E-03	
2.0E+00	S	5.7E-04	S	2.0E-05	I				6.5	1	0.5	No	~Aroclor 1254	11097-69-1	3.9E-02		9.8E-03	7.8E-03				4.0E-01		5.5E-03	
2.0E+00	S	5.7E-04	S						7.55	1	0	No	~Aroclor 1260	11096-82-5	3.9E-02		9.8E-03	7.8E-03						2.0E+00	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	6.34	1	0.7	No	~Aroclor 5460	11126-42-4					1.2E+01			1.2E+01		2.0E+00	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	8.27	1	0	No	~Heptachlorobiphenyl, 2,3,3',4,4',5,5' (PCB 189)	39635-31-9	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		2.8E-03	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.5	1	0	No	~Hexachlorobiphenyl, 2,3',4,4',5,5' (PCB 167)	52663-72-6	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		1.7E-03	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.6	1	0	No	~Hexachlorobiphenyl, 2,3,3',4,4',5,5' (PCB 157)	69782-90-7	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		1.7E-03	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.6	1	0	No	~Hexachlorobiphenyl, 2,3,3',4,4',5 (PCB 156)	38380-08-4	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		1.7E-03	
3.9E+03	E	1.1E+00	E	2.3E-08	E	1.3E-06	E	V	7.41	1	0.1	No	~Hexachlorobiphenyl, 3,3',4,4',5,5' (PCB 169)	32774-16-6	2.0E-05		4.9E-06	4.0E-06	4.7E-04		2.8E-03	4.0E-04		1.7E-06	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	6.98	1	0.4	No	~Pentachlorobiphenyl, 2',3,4,4',5 (PCB 123)	65510-44-3	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		1.0E-03	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.12	1	0.3	No	~Pentachlorobiphenyl, 2,3',4,4',5 (PCB 118)	31508-00-6	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		1.0E-03	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	6.79	1	0.5	No	~Pentachlorobiphenyl, 2,3,3',4,4' (PCB 105)	32598-14-4	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		1.0E-03	
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	6.98	1	0.4	No	~Pentachlorobiphenyl, 2,3,4,4',5 (PCB 114)	74472-37-0	2.0E-02		4.9E-03	4.0E-03	4.7E-01		2.8E+00	4.0E-01		1.0E-03	
1.3E+04	E	3.8E+00	E	7.0E-09	E	4.0E-07	E	V	6.98	1	0.4	No	~Pentachlorobiphenyl, 3,3',4,4',5 (PCB 126)	57465-28-8	6.0E-06		1.5E-06	1.2E-06	1.4E-04		8.3E-04	1.2E-04		3.0E-07	
2.0E+00	I	5.7E-04	I						7.1	1	0.7	No	~Polychlorinated Biphenyls (high risk)	1336-36-3									5.0E-01	6.8E-03	7.8E-02
4.0E-01	I	1.0E-04	I						7.1	1	0.7	No	~Polychlorinated Biphenyls (low risk)	1336-36-3	1.9E-01		5.6E-02	4.4E-02						6.8E-03	7.8E-02
7.0E-02	I	2.0E-05	I						7.1	1	0.7	No	~Polychlorinated Biphenyls (lowest risk)	1336-36-3										9.4E-04	
1.3E+01	E	3.8E-03	E	7.0E-06	E	4.0E-04	E	V	6.63	1	0.6	No	~Tetrachlorobiphenyl, 3,3',4,4' (PCB 77)	32598-13-3	6.0E-03		4.9E-03	6.0E-03	1.4E-01		1.4E-01			6.2E-05	
3.9E+01	E	1.1E-02	E	2.3E-06	E	1.3E-04	E	V	6.34	1	0.7	No	~Tetrachlorobiphenyl, 3,4,4',5 (PCB 81)	70362-50-4	2.0E-03		4.9E-04	4.0E-04	4.7E-02		2.8E-01	4.0E-02			
				6.0E-04	I				10.46	1	0	No	Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9											
													Polynuclear Aromatic Hydrocarbons (PAHs)												
				6.0E-02	I				3.92	1	1	Yes	~Acenaphthene	83-32-9					1.2E+03	9.6E+02		5.3E+02		5.5E+00	
7.3E-01	E	1.1E-04	C	3.0E-01	I				4.45	1	1	Yes	~Anthracene	120-12-7					6.0E+03	2.5E+03		1.8E+03		5.8E+01	
									5.76	1	1	No	~Benz(a)anthracene	56-55-3	3.4E-02		1.8E-02	1.2E-02						4.3E-03	
1.2E+00	C	1.1E-04	C						6.11	1	0.9	No	~Benzo(b)fluoranthene	205-82-3	6.5E-02			6.5E-02					2.0E-01	7.8E-02	2.4E-01
7.3E+00	I	1.1E-03	C						6.13	1	1	No	~Benzo(a)pyrene	50-32-8	3.4E-03			3.4E-03						4.0E-03	
7.3E-01	E	1.1E-04	C						5.78	1	1	No	~Benzo(k)fluoranthene	205-99-2	3.4E-02			3.4E-02						4.1E-02	
7.3E-02	E	1.1E-04	C						6.11	1	0.9	No	~Benzo(k)fluoranthene	207-08-9	3.4E-01			3.4E-01						4.0E-01	
7.3E-03	E	1.1E-05	C	8.0E-02	I				3.9	1	1	Yes	~Chloronaphthalene, 1,8-	91-58-7					1.6E+03	1.4E+03		7.5E+02		3.9E+00	
7.3E-03	E	1.1E-05	C						5.81	1	1	No	~Chrysene	218-01-9	3.4E+00			3.4E+00						1.2E+00	
7.3E+00	E	1.2E-03	C						6.75	1	0.6	No	~Dibenz(a,h)anthracene	53-70-3	3.4E-03			3.4E-03						1.3E-02	
1.2E+01	C	1.1E-03	C						7.71	1	0.3	No	~Dibenzo(a,e)pyrene	192-65-4	6.5E-03			6.5E-03						8.4E-02	
2.5E+02	C	7.1E-02	C						5.8	1	0.9	No	~Dimethylbenz(a)anthracene, 7,12-	57-97-6	1.0E-04			1.0E-04						9.9E-05	
				4.0E-02	I				5.16	1	1	No	~Fluoranthene	206-44-0					8.0E+02			8.0E+02		8.9E+01	
				4.0E-02	I				4.18	1	1	Yes	~Fluorene	86-73-7					8.0E+02	4.6E+02		2.9E+02		5.4E+00	
7.3E-01	E	1.1E-04	C						6.7	1	0.6	No	~Indeno(1,2,3-cd)pyrene	193-39-5	3.4E-02			3.4E-02						1.3E-01	
2.9E-02	P			7.0E-02	A				3.87	1	1	Yes	~Methylnaphthalene, 1-	90-12-0	2.7E+00	1.9E+00		1.1E+00						5.8E-03	
				4.0E-03	I				3.86	1	1	Yes	~Methylnaphthalene, 2-	81-57-6					8.0E+01	6.5E+01		3.6E			

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where nSL < 100X cSL; ** = where nSL < 10X cSL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHILD Hazard Index (HI) = 1			Protection of Groundwater SSL					
SFO (mg/kg-day)	k _e (y)	IUR (ug/m ³ -y)	k _e (mg/kg-day)	RfD _c (mg/m ³ -y)	k _e (mg/m ³ -y)	RfC _c (mg/m ³ -y)	k _e (y)	muta-gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)
2.0E-02	I								2.93	1		Yes	Propazine	139-40-2					4.0E+02	2.4E+03		3.4E+02		3.1E-01	
2.0E-02	I								2.6	1		Yes	Propionamide	122-42-9					4.0E+02	2.8E+03		3.5E+02		2.2E-01	
1.3E-02	I								3.72	1	0.9	Yes	Propiconazole	60207-90-1					2.6E+02	1.1E+03		2.1E+02		6.9E-01	
									0.59	1		Yes	Propionaldehyde	123-38-6							1.7E+01		3.4E-03		
1.0E-01	X	1.0E+00	X	V					3.69	1	1	Yes	Propyl benzene	103-65-1					2.0E+03	1.8E+03	2.1E+03	6.6E+02		1.2E+00	
									1.77	1	1	Yes	Propylene	115-07-1							6.3E+03		6.0E+00		
2.0E+01	P								-0.92	1	1	Yes	Propylene Glycol	57-55-6					4.0E+05	3.2E+08		4.0E+05		8.1E+01	
									1.59	1	1	Yes	Propylene Glycol Dinitrate	6423-43-4											
7.0E-01	H								0.002	1	1	Yes	Propylene Glycol Monoethyl Ether	1569-02-4					1.4E+04	3.3E+06		1.4E+04		2.8E+00	
7.0E-01	H	2.0E+00	I	V					-0.49	1	1	Yes	Propylene Glycol Monomethyl Ether	107-98-2					1.4E+04	3.9E+06	4.2E+03	3.2E+03		6.5E-01	
2.4E-01	I	3.7E-06	I						0.03	1	1	Yes	Propylene Oxide	75-56-9	3.2E-01	4.5E+01	1.5E+00				6.3E+01		5.6E-05		
									2.6	1	1	Yes	Pursuit	81335-77-5					5.0E+03	7.2E+04		4.7E+03		4.1E+00	
									6.2	1	0.7	No	Pyridin	51630-58-1					5.0E+02			5.0E+02		3.2E+02	
									0.65	1	1	Yes	Pyridine	110-86-1					2.0E+01	1.5E+03		2.0E+01		6.8E-03	
									4.44	1	0.9	Yes	Quinalphos	13593-03-8					1.0E+01	1.0E+01		5.1E+00		4.3E-02	
3.0E+00	I								2.03	1	1	Yes	Quinoline	91-22-5	2.6E-02	2.8E-01		2.4E-02						7.8E-05	
												Yes	Refractory Ceramic Fibers	NA											
									6.14	1	0.7	Yes	Resmethrin	10453-86-8					6.0E+02	7.6E+01		6.7E+01		4.2E+01	
									4.88	1	0.8	Yes	Ronnel	299-84-3					1.0E+03	6.8E+02		4.1E+02		3.7E+00	
2.2E-01	C	6.3E-05	C						4.1	1	0.9	Yes	Rotenone	83-79-4					8.0E+01	2.6E+02		6.1E+01		3.2E+01	
									3.45	1	1	Yes	Saflrole	94-59-7	1.1E-01	5.9E-01		9.5E-02						5.9E-05	
									5.57	1	0.8	Yes	Savey	78587-05-0					5.0E+02	1.4E+02		1.1E+02		5.0E-01	
												Yes	Selenious Acid	7783-00-8					1.0E+02	2.3E+04		1.0E+02			
												Yes	Selenium	7782-49-2					1.0E+02	2.3E+04		1.0E+02	5.0E+01	5.2E-01	2.6E-01
												Yes	Selenium Sulfide	7446-34-6					1.0E+02	2.3E+04		1.0E+02			
									4.38	1	0.9	Yes	Sethoxydim	74051-80-2					1.8E+03	2.4E+03		1.0E+03		9.3E+00	
												Yes	Silica (crystalline, respirable)	7631-86-9											
										0.04	1	Yes	Silver	7440-22-4											
1.2E-01	H								2.18	1	1	Yes	Simazine	122-34-9	6.5E-01	8.9E+00		6.1E-01					4.0E+00	3.0E-04	2.0E-03
									0.37	1	1	Yes	Sodium Acifluorfen	62476-59-9					2.6E+02	2.1E+05		2.6E+02		2.1E+00	
												Yes	Sodium Azide	26628-22-8					8.0E+01	1.8E+04		8.0E+01			
5.0E-01	C	1.5E-01	C						0.025	1	1	Yes	Sodium Dichromate	10588-01-9	5.0E-02	2.3E-01		4.1E-02							
2.7E-01	H								-1.431	1	1	Yes	Sodium Diethyldithiocarbamate	148-18-5	2.9E-01	8.2E+02		2.9E-01							
												Yes	Sodium Fluoride	7681-49-4					1.0E+03	2.3E+05		1.0E+03			
												No	Sodium Fluoroacetate	62-74-8					4.0E-01			4.0E-01		8.1E-05	
												Yes	Sodium Metavanadate	13718-26-8					2.0E+01	4.5E+03		2.0E+01			
2.4E-02	H								3.53	1	0.9	Yes	Stirofos (Tetrachlorovinphos)	961-11-5	3.2E+00	1.8E+01		2.8E+00						8.1E-03	
5.0E-01	C	1.5E-01	C						0.025	1	1	Yes	Strontium Chromate	7789-06-2	5.0E-02	2.3E-01		4.1E-02							
												Yes	Strontium, Stable	7440-24-6					1.2E+04	2.7E+06		1.2E+04		4.2E+02	
									1.93	1	1	Yes	Strychnine	57-24-9					6.0E+00	3.2E+02		5.9E+00		6.5E-02	
2.0E-01	I	1.0E+00	I	V					2.95	1	1	Yes	Styrene	100-42-5					4.0E+03	1.0E+04	2.1E+03	1.2E+03	1.0E+02	1.3E+00	1.1E-01
3.0E-03	P								3.1	1	1	Yes	Styrene-Acrylonitrile (SAN) Trimer	NA					6.0E+01	2.4E+02		4.8E+01			
1.0E-03	P	2.0E-03	X						-0.77	1	1	Yes	Sulfolane	126-33-0					2.0E+01	1.7E+04		2.0E+01		4.4E-03	
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		6.5E-02	
												Yes	Sulfur Trioxide	7446-11-9							2.1E+00		2.1E+00		
												Yes	Sulfuric Acid	7664-93-9											
2.5E-02	I								2.94	1	0.9	Yes	Systhane	88671-89-0					5.0E+02	4.8E+03		4.5E+02		5.6E+00	
3.0E-02	H								3.3	1	0.9	Yes	TCMTB	21664-17-0					6.0E+02	2.4E+03		4.8E+02		3.3E+00	
7.0E-02	I								1.79	1		Yes	Tebuthiuron	34014-18-1					1.4E+03	4.7E+04		1.4E+03		3.9E-01	
2.0E-02	H								5.96	1	0.7	No	Temephos	3383-96-8					4.0E+02			4.0E+02		7.6E+01	
1.3E-02	I								1.89	1		Yes	Terbacil	5902-51-2					2.6E+02	7.0E+03		2.5E+02		7.5E-02	
2.5E-05	H								4.48	1	0.9	Yes	Terbufos	13071-79-9					5.0E-01	4.5E-01		2.4E-01		5.2E-04	
1.0E-03	I								3.74	1	0.9	Yes	Terbutryn	889-50-0					2.0E+01	4.1E+01		1.3E+01		1.9E-02	
1.0E-04	I								6.77	1	0.6	No	Tetrabromodiphenylether, 2,2',4,4'-(BDE-47)	5436-43-1					2.0E+00			2.0E+00		5.4E-02	
3.0E-04	I								4.64	1		Yes	Tetrachlorobenzene, 1,2,4,5-	95-94-3					6.0E+00	2.4E+00		1.7E+00		7.9E-03	
2.6E-02	I	7.4E-06	I						2.93	1		Yes	Tetrachloroethane, 1,1,1,2-	630-28-6	3.0E+00	1.0E+01	7.6E-01	5.7E-01	6.0E+02	2.4E+03		4.8E+02		2.2E-04	
2.0E-01	I	5.8E-05	C						2.39	1		Yes	Tetrachloroethane, 1,1,2,2-	79-34-5	3.9E-01	3.1E+00	9.7E-02	7.6E-02	4.0E+02	3.6E+03		3.6E+02		3.0E-05	
2.1E-03	I	2.6E-07	I						3.4	1		Yes	Tetrachloroethylene	127-18-4	3.7E+01	6.3E+01	2.2E-01	1.1E+01	1.2E+02	2.3E+02	8.3E+01	4.1E+01	5.0E+00	5.1E-03	2.3E-03
2.0E+01	H								4.45	1	0.9	Yes	Tetrachlorophenol												

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1				Protection of Groundwater SSL			
SFO (mg/kg-day) ¹	k _e (y)	IUR (ug/m ³ -y)	RfD _c (mg/kg-day)	k _e (y)	RfD _c (mg/m ³ -y)	k _e (y)	muta-gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)	
6.0E-01	H		1.0E-04	A	V				1	1	Yes	Tin	7440-31-5					1.2E+04	2.7E+06	2.1E-01	1.2E+04		3.0E+03		
			1.0E-04	A	V				1	1	Yes	Titanium Tetrachloride	7550-45-0							2.1E-01	2.1E-01				
1.8E-01	X		8.0E-02	I	5.0E+00	I	V	2.73	1	1	Yes	Toluene	108-88-3	4.3E-01	7.9E+01		4.3E-01	1.6E+03	5.3E+03	1.0E+04	1.1E+03	1.0E+03	7.6E-01	6.9E-01	
3.0E-02	P		2.0E-04	X				0.16	1	1	Yes	Toluene-2,5-diamine	95-70-5	2.6E+00	6.5E+01		2.5E+00	4.0E+00	8.3E+02	2.3E+03	4.0E+00		1.3E-04		
			4.0E-03	X				1.39	1	1	Yes	Toluidine, p-	106-49-0					8.0E+01	2.3E+03	7.7E+01	7.7E+01		1.1E-03		
			3.0E+00	P			V	6.1	1	1	No	Total Petroleum Hydrocarbons (Aliphatic High)	NA					6.0E+04			6.0E+04		2.4E+03		
			6.0E-01	P	V			3.9	1	1	Yes	Total Petroleum Hydrocarbons (Aliphatic Low)	NA							1.3E+03	1.3E+03		8.8E+00		
			1.0E-02	X	1.0E-01	P	V	5.65	1	1	No	Total Petroleum Hydrocarbons (Aliphatic Medium)	NA					2.0E+02		2.1E+02	1.0E+02		1.5E+00		
			4.0E-02	P				5.16	1	1	No	Total Petroleum Hydrocarbons (Aromatic High)	NA					8.0E+02			8.0E+02		8.9E+01		
			4.0E-03	P	3.0E-02	P	V	2.13	1	1	Yes	Total Petroleum Hydrocarbons (Aromatic Low)	NA					8.0E+01	6.0E+02	6.3E+01	3.3E+01		1.7E-02		
			4.0E-03	P	3.0E-03	P	V	3.58	1	1	Yes	Total Petroleum Hydrocarbons (Aromatic Medium)	NA					8.0E+01	9.0E+01	6.3E+00	5.5E+00		2.3E-02		
1.1E+00	I	3.2E-04						5.9	1	0.8	Yes	Toxaphene	8001-35-2	7.1E-02	1.9E-02		1.5E-02	1.5E+02	6.0E+00	1.5E+02	3.7E+00	3.0E+00	2.4E-03	4.6E-01	
			7.5E-03	I				7.56	1	0.5	No	Tralometrin	66841-25-6					6.0E+00	9.8E+00		1.5E+02		5.8E+01		
			3.0E-04	A			V	4.1	1	0.9	Yes	Tri-n-butyltin	688-73-3										8.2E-02		
			8.0E+01	X				0.25	1	1	Yes	Triacetin	102-76-1					1.6E+06	5.3E+08		1.6E+06		4.5E+02		
			1.3E-02	I			V	4.6	1	0.9	Yes	Triallate	2303-17-5					2.6E+02	2.2E+02		1.2E+02		2.6E-01		
			1.0E-02	I				1.1	1	1	Yes	Triasulfuron	82097-50-5					2.0E+02	6.0E+04		2.0E+02		2.1E-01		
			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		6.4E-02		
9.0E-03	P		1.0E-02	P				4	1	0.9	Yes	Tributyl Phosphate	126-73-8	8.7E+00	1.2E+01		5.1E+00	2.0E+02	3.3E+02		1.2E+02		2.5E-02		
			3.0E-04	P				1	0	0	No	Tributyltin Compounds	NA					6.0E+00			6.0E+00				
			3.0E-04	I				4.05	1	1	Yes	Tributyltin Oxide	56-35-9					6.0E+00	9.5E+01		5.7E+00		2.9E+02		
			3.0E+01	I	3.0E+01	H	V	3.16	1	1	Yes	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1				1.1E+00	6.0E+05	1.9E+06	6.3E+04	5.5E+04	6.0E+01	1.4E+02	1.2E-02	
7.0E-02	I		2.0E-02	I				1.33	1	1	Yes	Trichloroacetic Acid	76-03-9	1.1E+00	4.4E+01		1.1E+00	4.0E+02	1.8E+04		3.9E+02		2.2E-04		
2.9E-02	H							-0.67	1	1	Yes	Trichloroaniline HCl, 2,4,6-	33663-50-2		2.7E+00	3.6E+03		2.7E+00					7.4E-03		
7.0E-03	X		3.0E-05	X				3.52	1	1	Yes	Trichloroaniline, 2,4,6-	634-93-5	1.1E+01	1.9E+01		7.0E+00	6.0E-01	1.2E+00		4.0E-01		3.6E-03		
			8.0E-04	X			V	4.05	1	1	Yes	Trichlorobenzene, 1,2,3-	87-61-6					1.6E+01	1.3E+01		7.0E+00		2.1E-02		
2.9E-02	P		1.0E-02	I	2.0E-03	P	V	4.02	1	1	Yes	Trichlorobenzene, 1,2,4-	120-82-1	2.7E+00	1.9E+00		1.1E+00	2.0E+02	1.6E+02	4.2E+00	4.0E+00	7.0E+01	3.3E-03	2.0E-01	
			2.0E+00	I	5.0E+00	I	V	2.49	1	1	Yes	Trichloroethane, 1,1,1-	71-55-6	4.0E+04	2.5E+05		1.0E+04	4.0E+04	2.5E+05	1.0E+04	8.0E+03	2.0E+02	2.8E+00	7.0E-02	
5.7E-02	I	1.6E-05	4.0E-03	I	2.0E-04	X	V	1.89	1	1	Yes	Trichloroethane, 1,1,2-	79-00-5	1.4E+00	1.9E+01	3.5E-01	2.8E-01	8.0E+01	1.3E+03	4.2E-01	4.1E-01	5.0E+00	8.9E-05	1.6E-03	
4.6E-02	I	4.1E-06	5.0E-04	I	2.0E-03	I	V	2.42	1	1	Yes	Trichloroethylene	79-01-6	1.2E+00	7.2E+00	9.6E-01	4.9E-01	1.0E+01	6.9E+01	4.2E+00	2.8E+00	5.0E+00	1.8E-04	1.8E-03	
			3.0E-01	I	7.0E-01	H	V	2.53	1	1	Yes	Trichlorofluoromethane	75-69-4					6.0E+03	3.6E+04	1.5E+03	1.1E+03		7.3E-01		
			1.0E-01	I				3.72	1	1	Yes	Trichlorophenol, 2,4,5-	95-95-4					2.0E+03	2.9E+03		2.5E+03		4.4E+00		
1.1E-02	I	3.1E-06	1.0E-03	P				3.69	1	1	Yes	Trichlorophenol, 2,4,6-	88-06-2	7.1E+00	9.4E+00		4.0E+00	2.0E+01	3.0E+01		1.2E+01		1.5E-02		
			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		6.8E-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	6.1E-02	2.8E-02	
3.0E+01	I		5.0E-03	I			V	2.43	1	1	Yes	Trichloropropane, 1,1,2-	599-77-6				7.5E-04	1.0E+02	7.5E+02		8.8E+01		3.5E-02		
			4.0E-03	I	3.0E-04	I	V	2.27	1	1	Yes	Trichloropropane, 1,2,3-	96-18-4	8.4E-04	7.1E-03			8.0E+01	7.7E+02	6.3E-01	6.2E-01		3.2E-07		
			3.0E-03	X	3.0E-04	P	V	2.78	1	1	Yes	Trichloropropene, 1,2,3-	96-19-5					6.0E+01	2.6E+02	6.3E-01	6.2E-01		3.1E-04		
			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		1.5E+01		
			3.0E-03	I				5.18	1	0.8	Yes	Tri-diphenylphosphine oxide	58138-08-2					6.0E+01	2.6E+01		1.8E+01		1.3E-01		
			7.0E-03	I	V			1.45	1	1	Yes	Triethylamine	121-44-8							1.5E+01		1.5E+01		4.4E-03	
			2.0E+00	P				-1.75	1	1	Yes	Triethylene Glycol	112-27-6					4.0E+04	1.8E+08		4.0E+04		8.8E+00		
7.7E-03	I		7.5E-03	I			V	5.34	1	0.8	Yes	Triethyltin chloride	1562-09-8	1.0E+01	1.3E+00		2.5E+00	1.5E+02	5.5E+01		4.0E+01		8.2E-02		
2.0E-02	P		1.0E-02	P				-0.65	1	1	Yes	Trimethyl Phosphate	512-56-1	3.9E+00	2.7E+03		3.9E+00	2.0E+02	1.6E+05		2.0E+02		8.6E-04		
			5.0E-03	P	V			3.66	1	1	Yes	Trimethylbenzene, 1,2,3-	526-73-8							1.0E+01		1.0E+01		1.5E-02	
			7.0E-03	P	V			3.63	1	1	Yes	Trimethylbenzene, 1,2,4-	85-63-6							1.5E+01		1.5E+01		2.1E-02	
			1.0E-02	X			V	3.42	1	1	Yes	Trimethylbenzene, 1,3,5-	109-67-8					2.0E+02	2.8E+02		1.2E+02		1.7E-01		
3.0E-02	I		3.0E-02	I				1.18	1	1	Yes	Trinitrobenzene, 1,3,5-	89-35-4					6.0E+02	4.7E+04		5.9E+02		2.1E+00		
			5.0E-04	I				1.6	1	1	Yes	Trinitrotoluene, 2,4,6-	118-96-7	2.6E+00	1.0E+02		2.5E+00	1.0E+01	4.5E+02		9.8E+00		1.5E-02		
			2.0E-02	P				2.83	1	1	Yes	Triphenylphosphine Oxide	791-28-6					4.0E+02	3.8E+03		3.6E+02		1.5E+00		
			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		8.0E+00		
2.3E+00	C	6.6E-04	1.0E-02	X			V	2.59	1	1	No	Tris(1-chloro-2-propyl)phosphate	13674-84-6					2.0E+02	3.8E+03		1.9E+02		6.5E-01		
2.0E-02	P		7.0E-03	P				1.44	1	1	Yes	Tris(2,3-dibromopropyl)phosphate	128-72-7	3.4E-02		8.5E-03	6.8E-03	3.4E-02				1.9E+02		1.3E-04	
3.2E-03	P</																								

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer CHLD Hazard Index (HI) = 1			Protection of Groundwater SSL				
SFO (mg/kg-day) ¹	ke y	IUR (ug/m ³) ¹	ke y	RD ₀ (mg/kg-day)	ke y	RF _C (mg/m ³)	ke y	RF _S (mg/m ³)	ke y	muta- gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (ug/L)	Dermal SL TR=1.0E-6 (ug/L)	Inhalation SL TR=1.0E-6 (ug/L)	Carcinogenic SL TR=1.0E-6 (ug/L)	Ingestion SL Child HQ=1 (ug/L)	Dermal SL Child HQ=1 (ug/L)	Inhalation SL Child HQ=1 (ug/L)	Noncarcinogenic SL Child HI=1 (ug/L)	MCL (ug/L)	Risk-based SSL (mg/kg)	MCL-based SSL (mg/kg)
				3.0E-01		1						1	1	Yes	Zinc and Compounds	7440-66-6					6.0E+03	2.3E+06		6.0E+03		3.7E+02	
				5.0E-02		1				1.3		1	1	Yes	Zinc	12122-67-7					1.0E+03	9.7E+04		9.9E+02		2.9E+00	
				8.0E-05		X						1	1	Yes	Zirconium	7440-67-7					1.6E+00	3.6E+02		1.6E+00		4.8E+00	

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Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	ke y	IUR (ug/m ³) ⁻¹	ke y	RfD _o (mg/kg-day)	ke y	RfC _i (mg/m ³)	ke y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)
1.8E-02	C	5.1E-06	C	1.5E-01	I				1.0E+00	1.0E-01		1.4E+09		ALAR	1596-84-5	1.8E+02	4.3E+02	3.3E+06	1.3E+02	1.8E+05	4.1E+05		1.2E+05
8.7E-03	I	2.2E-06	I	4.0E-03	I	9.0E-03	I	V	1.0E+00	1.0E-01		1.4E+09	8.7E+03	Acephate	30560-19-1	3.8E+02	8.9E+02		2.6E+02	4.7E+03	1.1E+04		3.3E+03
														Acetaldehyde	75-07-0			4.9E+01			3.4E+02		3.4E+02
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Acetochlor	34256-82-1					2.3E+04	5.5E+04		1.6E+04
				9.0E-01	I	3.1E+01	A	V	1.0E+00			1.1E+05	1.4E+04	Acetone	67-64-1					1.1E+06		1.8E+06	6.7E+05
						2.0E-03	X	V	1.0E+00			1.1E+05	1.4E+09	Acetone Cyanohydrin	75-86-5						2.1E+02		2.1E+02
				6.0E-02	I				1.0E+00		1.3E+05	1.4E+09	1.3E+04	Acetonitrile	75-05-8							3.4E+03	3.4E+03
3.8E+00	C	1.3E-03	C	1.0E-01	I				1.0E+00	1.0E-01		1.4E+09		Acetophenone	98-86-2					1.2E+05			1.2E+05
									1.0E+00	1.0E-01		1.4E+09		Acetylaminofluorene, 2-	53-96-3	8.6E-01	2.0E+00	1.3E+04	6.0E-01				
				5.0E-04	I	2.0E-05	I	V	1.0E+00		2.3E+04	1.4E+09	6.9E+03	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	1.0E+00	1.0E-01		1.4E+09		Acrylamide	79-06-1	6.5E+00	1.5E+01	1.7E+05	4.6E+00	2.3E+03	5.5E+03	3.6E+07	1.6E+03
				5.0E-01	I	1.0E-03	I	V	1.0E+00			1.1E+05	1.4E+09	Acrylic Acid	79-10-7					5.8E+05		4.2E+02	4.2E+02
5.4E-01	I	6.8E-05	I	4.0E-02	A	2.0E-03	I	V	1.0E+00		1.1E+04	1.4E+09	7.7E+03	Acrylonitrile	107-13-1	6.1E+00		1.4E+00	1.1E+00	4.7E+04		6.7E+01	6.7E+01
5.6E-02	C			1.0E-02	I	6.0E-03	P		1.0E+00	1.0E-01		1.4E+09		Adiponitrile	111-69-3					1.2E+04	2.8E+04		3.6E+07
									1.0E+00	1.0E-01		1.4E+09		Alachlor	15972-60-8	5.8E+01	1.4E+02		4.1E+01				8.2E+03
				1.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Aldicarb	116-06-3					1.2E+03	2.8E+03		8.2E+02
				1.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Aldicarb Sulfone	1646-88-4					1.2E+03	2.8E+03		8.2E+02
									1.0E+00	1.0E-01		1.4E+09		Aldicarb sulfoxide	1646-87-3								
1.7E+01	I	4.9E-03	I	3.0E-05	I				1.0E+00			1.4E+09	1.7E+06	Aldrin	509-00-2	1.9E-01		4.3E+00	1.8E-01	3.5E+01			3.5E+01
				2.5E-01	I				1.0E+00	1.0E-01		1.4E+09		Allyl	74223-64-6					2.9E+05	6.9E+05		2.1E+05
				5.0E-03	I	1.0E-04	X	V	1.0E+00		1.1E+05	1.4E+09	3.4E+04	Allyl Alcohol	107-18-6					5.8E+03		1.5E+01	1.5E+01
2.1E-02	C	6.0E-06	C	1.0E+00	P	5.0E-03	P		1.0E+00		1.4E+03	1.4E+09	1.8E+03	Allyl Chloride	107-05-1	1.6E+02		3.2E+00	3.2E+00			6.9E+00	6.9E+00
				4.0E-04	I				1.0E+00			1.4E+09		Aluminum	7429-90-5					1.2E+06		3.0E+07	1.1E+06
									1.0E+00			1.4E+09		Aluminum Phosphide	20859-73-8					4.7E+02			4.7E+02
2.1E+01	C	6.0E-03	C	3.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Amdri	67485-29-4	1.6E-01	3.7E-01	2.8E+03	1.1E-01	3.5E+02	8.3E+02		2.5E+02
				9.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Ametryn	834-12-8					1.1E+04	2.5E+04		7.4E+03
									1.0E+00	1.0E-01		1.4E+09		Aminodiphenyl, 4-	92-67-1								
				8.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Aminophenol, m-	591-27-5					9.3E+04	2.2E+05		6.6E+04
				2.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Aminophenol, p-	123-30-8					2.3E+04	5.5E+04		1.6E+04
				2.5E-03	I				1.0E+00	1.0E-01		1.4E+09		Amtraz	33089-61-1					2.9E+03	6.9E+03		2.1E+03
				2.0E-01	I	1.0E-01	I	V	1.0E+00			1.4E+09		Ammonia	7804-41-7					2.3E+05			2.3E+05
						3.0E-03	X	V	1.0E+00		1.4E+04	1.4E+09	2.6E+04	Ammonium Sulfamate	7773-06-0								3.4E+02
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I		1.0E+00	1.0E-01		1.4E+09		Amyl Alcohol, tert.	75-85-4	5.7E+02	1.4E+03	1.0E+07	4.0E+02	8.2E+03	1.9E+04	6.0E+06	5.7E+03
4.0E-02	P			2.0E-03	X				1.0E+00	1.0E-01		1.4E+09		Aniline	62-53-3	8.2E+01	1.9E+02		5.7E+01	2.3E+03	5.5E+03		1.6E+03
				4.0E-04	I				1.5E-01			1.4E+09		Antraquinone, 9,10-	84-85-1					4.7E+02			4.7E+02
									1.5E-01			1.4E+09		Antimony (metallic)	7440-36-0								
				5.0E-04	H				1.5E-01			1.4E+09		Antimony Pentoxide	1314-60-9					5.8E+02			5.8E+02
				9.0E-04	H				1.5E-01			1.4E+09		Antimony Potassium Tartrate	11071-15-1					1.1E+03			1.1E+03
				4.0E-04	H				1.5E-01			1.4E+09		Antimony Tetroxide	1332-81-6					4.7E+02			4.7E+02
				2.0E-04	I				1.5E-01			1.4E+09		Antimony Trioxide	1309-84-4							1.2E+06	1.2E+06
2.5E-02	I	7.1E-06	I	1.3E-02	I				1.0E+00	1.0E-01		1.4E+09		Apollo	74115-24-5	1.3E+02	3.1E+02	2.3E+06	9.2E+01	1.5E+04	3.6E+04		1.1E+04
				5.0E-02	H				1.0E+00	1.0E-01		1.4E+09		Aramite	140-67-8					5.8E+04	1.4E+05		4.1E+04
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C		1.0E+00	3.0E-02		1.4E+09		Arsenic, Inorganic	7440-38-2	3.6E+00	1.7E+01	3.9E+03	3.0E+00	5.8E+02	2.8E+03	8.9E+04	4.8E+02
				3.5E-06	C	5.0E-05	I		1.0E+00			1.4E+09		Arsine	7784-42-1					4.1E+00		3.0E+05	4.1E+00
				9.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Assure	76578-14-8					1.1E+04	2.5E+04		7.4E+03
				5.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Asulam	3337-71-1					5.8E+04	1.4E+05		4.1E+04
2.3E-01	C			3.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Atrazine	1912-24-9	1.4E+01	3.4E+01	6.7E+04	1.0E+01	4.1E+04	9.7E+04		2.9E+04
8.8E-01	C	2.5E-04	C						1.0E+00	1.0E-01		1.4E+09		Auramine	492-80-8	3.7E+00	8.8E+00		2.6E+00				
1.1E-01	I	3.1E-05	I	4.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Avermectin B1	65195-55-3	3.0E+01		2.1E+02	2.6E+01	4.7E+02	1.1E+03		3.3E+02
				1.0E+00	P	7.0E-06	P		1.0E+00	1.0E-01		1.4E+09	5.2E+05	Azobenzene	103-33-3					1.2E+06	2.8E+06	4.2E+04	4.0E+04
									1.0E+00	1.0E-01		1.4E+09		Azodicarbonamide	123-77-3								
5.0E-01	C	1.5E-01	C	2.0E-01	I	5.0E-04	H		7.0E-02			1.4E+09		Barium	7440-39-3	6.5E+00		1.1E+02	6.2E+00	2.3E+05	3.0E+06		2.2E+05
				2.0E-02	C	2.0E-04	C	M	2.5E-02			1.4E+09		Barium Chromate	10294-40-3					2.3E+04	1.2E+06		2.3E+04
				4.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Baygon	114-26-1					4.7E+03	1.1E+04		3.3E+03
				3.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Bayleton	43121-43-3					3.5E+04	8.3E+04		2.5E+04
				2.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Baythroid	68359-37-5					2.9E+04	6.9E+04		2.1E+04

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Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	ke IUR (ug/m ³ -y) ⁻¹	ke (mg/kg-day)	RfD _o (mg/kg-day)	ke RfC _i (mg/m ³ -y)	ke V o c mutagen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
2.7E-01	X		2.0E-03	H	3.0E-05	I		1.0E+00	1.4E+09	1.9E+04	Chloroacetaldehyde, 2-Chloroacetic Acid Chloroacetophenone, 2-	107-20-0 79-11-8 532-27-4	1.2E+01			1.2E+01	2.3E+03	5.5E+03		1.8E+05	1.6E+03 1.8E+05
2.0E-01	P		4.0E-03	I	5.0E-02	P V		1.0E+00	1.4E+09	1.4E+09	Chloroaniline, p-Chlorobenzene Chlorobenzilate	106-47-8 108-90-7 510-15-6	1.6E+01	3.9E+01		1.1E+01	4.7E+03	1.1E+04		1.4E+03	3.3E+03 1.3E+03 1.6E+04
1.1E-01	C	3.1E-05	2.0E-02	I				1.0E+00	1.4E+09	1.4E+09	Chlorobenzoic Acid, p-Chlorobenzotrifluoride, 4-Chlorobutane, 1-Chlorodifluoromethane Chloroethanol, 2-Chloroform	74-11-3 98-56-6 109-69-3 75-45-6 107-07-3 67-66-3	3.0E+01	7.0E+01	5.4E+05	2.1E+01	3.5E+04	8.3E+04		8.9E+03	2.5E+04 2.5E+03 4.7E+04
3.1E-02	C	2.3E-05	1.0E-02	I	9.8E-02	A V		1.0E+00	1.4E+09	2.6E+03	Chloromethane Chloromethyl Methyl Ether Chloronitrobenzene, o-Chlorotoluene, p-	74-87-3 107-30-2 88-73-3 1897-45-6 95-49-8 106-43-4	1.1E+02		1.4E+00	1.4E+00	2.3E+04	1.2E+04		1.1E+03	2.1E+05 2.1E+05 2.3E+04 1.0E+03
2.4E+00	C	6.9E-04	3.0E-03	P	1.0E-05	X		1.0E+00	1.4E+09	5.3E+03	Chloromethane Chloromethyl Methyl Ether Chloronitrobenzene, o-Chlorotoluene, p-	74-87-3 107-30-2 88-73-3 1897-45-6 95-49-8 106-43-4	1.4E+00	1.1E+01	2.6E+01	9.5E-02	8.9E-02	7.7E+00			3.5E+03 8.3E+03 6.0E+04 2.4E+03
6.3E-03	P		1.0E-03	P	6.0E-04	P		1.0E+00	1.4E+09	1.4E+09	Chloronitrobenzene, p-Chlorophenol, 2-Chloropicrin	100-00-5 95-57-8 76-06-2	5.2E+02	1.2E+03		3.6E+02	1.2E+03	2.8E+03		3.6E+06	8.2E+02 5.8E+03 8.2E+00
3.1E-03	C	8.9E-07	1.5E-02	I				1.0E+00	1.4E+09	8.1E+03	Chlorothalonil Chlorotoluene, o-Chlorotoluene, p-	1897-45-6 95-49-8 106-43-4	1.1E+03	2.5E+03	1.9E+07	7.4E+02	1.8E+04	4.1E+04			1.2E+04 2.3E+04 2.3E+04
2.4E+02	C	6.9E-02	2.0E-01	I				1.0E+00	1.4E+09	1.4E+09	Chlorzoxolon Chlorpropylamine Chlorpyrifos	54749-90-5 101-21-3 2921-88-2	1.4E-02	3.2E-02	2.4E+02	9.6E-03	2.3E+05	5.5E+05			1.6E+05 8.2E+02
5.0E-01	J	8.4E-02	3.0E-03	I	1.0E-04	I	M	2.5E-02	1.4E+09	1.4E+09	Chlorpyrifos Methyl Chlorosulfuron Chlorthiophos Chromium(III) Insoluble Salts Chromium(VI) Chromium, Total	5598-13-0 64902-72-3 60238-56-4 16895-83-1 18540-29-9 7440-47-3	6.5E+00		2.0E+02	6.3E+00	1.2E+04	2.8E+04		6.0E+05	8.2E+03 4.1E+04 6.6E+02 1.8E+06 3.5E+03
9.0E-03	P	6.2E-04	3.0E-04	P	6.0E-06	P		1.0E+00	1.4E+09	1.4E+09	Cobalt Coke Oven Emissions Copper	440-48-4 8007-45-2 7440-50-8			1.9E+03	1.9E+03	3.5E+02			3.6E+04	3.5E+02 4.7E+04
5.0E-02	I		6.0E-01	C				1.0E+00	1.4E+09	1.4E+09	Cresol, m-Cresol, o-Cresol, p-Cresol	108-39-4 95-48-7 106-44-5					5.8E+04	1.4E+05		3.6E+09	4.1E+04 4.1E+04 8.2E+04
1.9E+00	H		1.0E-01	A	6.0E-01	C		1.0E+00	1.4E+09	1.4E+09	Cresol, p-chloro-m-Cresolis Crotohaldehyde, trans-Cumene	59-50-7 1319-77-3 123-73-9 98-82-8	1.7E+00		1.7E+00		1.2E+05	2.8E+05		3.6E+09	8.2E+04 8.2E+04 1.2E+03 9.9E+03
2.2E-01	C	6.3E-05	2.0E-03	H				1.0E+00	1.4E+09	1.4E+09	Cupferron Cyanazine	135-20-6 21725-46-2	1.5E+01	3.5E+01	2.6E+05	1.0E+01	2.3E+03	5.5E+03			1.6E+03
			1.0E-03	I				1.0E+00	1.4E+09	1.4E+09	Cyanides						1.2E+03				1.2E+03
			5.0E-03	I				1.0E+00	1.4E+09	1.4E+09	-Calcium Cyanide -Copper Cyanide	592-01-8 544-92-3					5.8E+03				5.8E+03
			6.0E-04	I	8.0E-04	S V		1.0E+00	1.4E+09	3.5E+03	-Cyanide (CN ₂) -Cyanogen -Cyanogen Bromide	57-12-5 460-19-5 506-68-3					7.0E+02		1.2E+01		1.2E+01 1.2E+03 1.1E+05
			5.0E-02	I				1.0E+00	1.4E+09	1.4E+09	-Cyanogen Chloride -Hydrogen Cyanide -Potassium Cyanide	506-77-4 74-90-8 151-50-8					5.8E+04		1.8E+02		5.8E+04 1.5E+02 2.3E+03
			5.0E-03	I				4.0E-02	1.4E+09	1.4E+09	-Potassium Silver Cyanide -Silver Cyanide -Sodium Cyanide	506-61-6 506-64-9 143-33-9					5.8E+03				5.8E+03 1.2E+05 1.2E+03
			2.0E-04	P				1.0E+00	1.4E+09	1.4E+09	-Thiocyanates -Thiocyanic Acid -Zinc Cyanide	NA 463-56-9 557-21-1					2.3E+02				2.3E+02 2.3E+02 5.8E+04
2.3E-02	H		6.0E+00	I V				1.0E+00	1.4E+09	1.0E+03	Cyclohexane Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-Cyclohexanone	110-82-7 87-84-3 108-94-1	1.4E+02	3.4E+02		1.0E+02	5.8E+06		2.7E+04		2.7E+04 1.3E+05 1.3E+05
			5.0E-03	P	1.0E+00	X V		1.0E+00	2.8E+02	1.5E+03	Cyclohexene Cyclohexylamine Cyhalothrin/karate	110-83-8 108-91-8 68085-85-8					5.8E+03		6.4E+03		3.1E+03 2.3E+05 5.8E+03
			2.0E-01	I				1.0E+00	2.9E+05	1.4E+09	Cyhalothrin/karate	68085-85-8					5.8E+03	1.4E+04			2.3E+05 4.1E+03
			1.0E-02	I				1.0E+00	1.4E+09	1.4E+09	Cypermethrin	52315-07-8					1.2E+04	2.8E+04			8.2E+03

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Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	ke y ⁻¹	IUR (ug/m ³ -y) ⁻¹	ke y ⁻¹	RfD _o (mg/kg-day)	ke y ⁻¹	RfC _o (mg/m ³ -y)	ke y ⁻¹	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
2.4E-01	I	6.9E-05	C	7.5E-03	I				1.0E+00	1.0E-01		1.4E+09		Cyromazine	66215-27-8									
3.4E-01	I	9.7E-05	C		I			V				1.0E+00	2.1E+06	DDD	72-54-8	1.4E+01	3.2E+01	2.4E+05	9.6E+00	8.8E+03	2.1E+04		6.2E+03	
3.4E-01	I	9.7E-05	I	5.0E-04	I				1.0E+00	3.0E-02		1.4E+09		DDE, p,p'-	72-55-9	9.6E+00	2.7E+02	9.3E+00	9.3E+00	5.8E+02	4.6E+03		5.2E+02	
				1.0E-02	I				1.0E+00	1.0E-01		1.4E+09		DDT	50-29-3	9.6E+00	7.6E+01	1.7E+05	8.5E+00	1.2E+04	2.8E+04		8.2E+03	
				3.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Dacthal	1861-32-1									
7.0E-04	I			7.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Dalapon	75-99-0	4.7E+03	1.1E+04		3.3E+03	3.5E+04	8.3E+04		2.5E+04	
				4.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5					8.2E+03	1.9E+04		5.7E+03	
									1.0E+00	1.0E-01		1.4E+09		Demeton	8065-48-3					4.7E+01	1.1E+02		3.3E+01	
1.2E-03	I			6.0E-01	I				1.0E+00	1.0E-01		1.4E+09		Di(2-ethylhexyl)adipate	103-23-1	2.7E+03	6.4E+03		1.9E+03	7.0E+05	1.7E+06		4.9E+05	
6.1E-02	H			7.0E-04	A				1.0E+00	1.0E-01		1.4E+09		Diallate	2303-16-4	5.4E+01	1.3E+02		3.8E+01	8.2E+02	1.9E+03		5.7E+02	
					X			V	1.0E+00			5.2E+05		Diazinon	333-41-5					1.2E+04			1.2E+04	
8.0E-01	P	6.0E-03	P	1.0E-02	X			V	1.0E+00			9.8E+02	5.2E+05	Dibenzothiophene	132-65-0	4.1E+00		6.5E-02	6.4E-02	2.3E+02		2.8E+01	1.2E+04	
				2.0E-04	P	2.0E-04	I	V	1.0E+00			1.6E+02	1.4E+09	Dibromo-3-chloropropane, 1,2-	96-12-8					3.5E+02			2.5E+01	
				4.0E-04	X			V	1.0E+00			1.6E+02	1.4E+09	Dibromobenzene, 1,3-	108-36-1					4.7E+02			4.7E+02	
8.4E-02	I	2.7E-05	C	1.0E-02	I			V	1.0E+00			8.0E+02	1.4E+09	Dibromobenzene, 1,4-	106-37-6	3.9E+01		3.6E+00	3.3E+00	1.2E+04			1.2E+04	
2.0E+00	I	6.0E-04	I	2.0E-02	I			V	1.0E+00			8.0E+02	1.4E+09	Dibromochloromethane	124-48-1	1.6E+00		1.8E-01	1.6E-01	2.3E+04			2.3E+04	
				9.0E-03	I	9.0E-03	I	V	1.0E+00			1.3E+03	1.4E+09	Dibromoethane, 1,2-	106-93-4					1.1E+04		3.4E+02	3.3E+02	
				1.0E-02	H	4.0E-03	X	V	1.0E+00			2.8E+03	1.4E+09	Dibromomethane (Methylene Bromide)	74-95-3					1.2E+04		9.9E+01	9.8E+01	
				3.0E-04	P				1.0E+00	1.0E-01		1.4E+09		Dibutyltin Compounds	NA					3.5E+02	8.3E+02		2.5E+02	
				3.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Dicamba	1918-00-9					3.5E+04	8.3E+04		2.5E+04	
4.2E-03	P							V	1.0E+00			5.2E+02	1.4E+09	Dichloro-2-butene, 1,4-	764-41-0			3.6E-02	3.6E-02					
4.2E-03	P							V	1.0E+00			5.2E+02	1.4E+09	Dichloro-2-butene, cis-1,4-	1476-11-5			3.2E-02	3.2E-02					
4.2E-03	P							V	1.0E+00			7.6E+02	1.4E+09	Dichloro-2-butene, trans-1,4-	110-57-6			3.2E-02	3.2E-02					
5.0E-02	I			4.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Dichloroacetic Acid	79-43-6	6.5E+01	1.5E+02		4.6E+01	4.7E+03	1.1E+04		3.3E+03	
				9.0E-02	I	2.0E-01	H	V	1.0E+00			3.8E+02	1.4E+09	Dichlorobenzene, 1,2-	95-50-1					1.1E+05			1.0E+04	
5.4E-03	C	1.1E-05	C	7.0E-02	A	8.0E-01	I	V	1.0E+00			1.4E+09	1.0E+04	Dichlorobenzene, 1,4-	106-46-7	6.1E+02		1.2E+01	1.1E+01	8.2E+04		3.7E+04	2.5E+04	
4.5E-01	I	3.4E-04	C						1.0E+00	1.0E-01		1.4E+09		Dichlorobenzidine, 3,3'	91-94-1	7.3E+00	1.7E+01	4.9E+04	5.1E+00	2.3E+05			7.4E+03	
				9.0E-03	X				1.0E+00	1.0E-01		1.4E+09		Dichlorobenzophenone, 4,4'	90-98-2					1.1E+04	2.5E+04		3.7E+02	
				2.0E-01	I	1.0E-01	X	V	1.0E+00			8.5E+02	1.4E+09	Dichlorodifluoromethane	75-71-8					2.3E+05			3.7E+02	
5.7E-03	C	1.6E-06	C	2.0E-01	P			V	1.0E+00			1.7E+03	1.4E+09	Dichloroethane, 1,1-	75-34-3	5.7E+02		1.6E+01	1.6E+01	2.3E+05			2.3E+05	
9.1E-02	I	2.6E-05	I	6.0E-03	X	7.0E-03	P	V	1.0E+00			3.0E+03	1.4E+09	Dichloroethane, 1,2-	107-06-2	3.6E+01		2.2E+00	2.0E+00	7.0E+03		1.4E+02	1.4E+02	
				5.0E-02	I	2.0E-01	I	V	1.0E+00			1.2E+03	1.4E+09	Dichloroethylene, 1,1-	75-35-4					5.8E+04		1.0E+03	1.0E+03	
				2.0E-03	I			V	1.0E+00			2.4E+03	1.4E+09	Dichloroethylene, 1,2-cis-	158-59-2					2.3E+03			2.3E+03	
				2.0E-02	I			V	1.0E+00			1.9E+03	1.4E+09	Dichloroethylene, 1,2-trans-	156-60-5					2.3E+04			2.3E+04	
				3.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Dichlorophenol, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
				1.0E-02	I				1.0E+00	5.0E-02		1.4E+09		Dichlorophenoxy Acetic Acid, 2,4-	94-75-7					1.2E+04	5.5E+04		9.6E+03	
				8.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Dichlorophenoxybutyric Acid, 4-(2,4-	94-82-6					9.3E+03	2.2E+04		6.6E+03	
3.6E-02	C	1.0E-05	C	9.0E-02	A	4.0E-03	I	V	1.0E+00			1.4E+03	1.4E+09	Dichloropropene, 1,2-	78-87-5	9.1E+01		4.6E+00	4.4E+00	1.1E+05		6.6E+01	6.6E+01	
				2.0E-02	P			V	1.0E+00			1.5E+03	1.4E+09	Dichloropropane, 1,3-	142-28-9					2.3E+04			2.3E+04	
				3.0E-03	I				1.0E+00	1.0E-01		1.4E+09		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	1.0E+00			1.8E+03	1.4E+09	Dichloropropene, 1,3-	542-75-8	3.3E+01		1.1E+01	8.2E+00	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		1.0E+00	1.0E-01		1.4E+09		Dichlorvos	62-73-7	1.1E+01	2.7E+01	2.0E+05	7.9E+00	5.8E+02	1.4E+03	3.0E+06	4.1E+02	
				8.0E-02	P	3.0E-04	X	V	1.0E+00			1.4E+09	4.1E+03	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				1.0E+00	1.0E-01		1.4E+09		Dieldrin	60-57-1	2.0E-01	4.8E-01	3.6E+03	1.4E-01	5.8E+01	1.4E+02		4.1E+01	
				3.0E-04	C			I	1.0E+00	1.0E-01		1.4E+09		Diesel Engine Exhaust	NA									
				2.0E-03	P	2.0E-04	P		1.0E+00	1.0E-01		1.4E+09		Diethanolamine	111-42-2					2.3E+03	5.5E+03	1.2E+06	1.6E+03	
				3.0E-02	P	1.0E-04	P		1.0E+00	1.0E-01		1.4E+09		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		1.0E+00	1.0E-01		1.4E+09		Diethylene Glycol Monoethyl Ether	111-90-0					7.0E+04	1.7E+05	1.8E+06	4.8E+04	
3.5E+02	C	1.0E-01	C	1.0E-03	P			V	1.0E+00			1.1E+05	1.4E+09	Diethylformamide	617-84-5					1.2E+03			1.2E+03	
				1.0E+00	1.0E-01				1.0E+00	1.0E-01		1.4E+09		Diethylstilbestrol	56-53-1	9.3E-03	2.2E-02	1.7E+02	6.6E-03					
				8.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Difenzoquat	43222-48-6					9.3E+04	2.2E+05		6.6E+04	
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Diflubenzuron	35367-38-5					2.3E+04	5.5E+04		1.6E+04	
				4.0E+01	I	V			1.0E+00			1.4E+03	1.4E+09	Difluoroethane, 1,1-	75-37-6							2.0E+05	2.0E+05	
4.4E-02	C	1.3E-05	C					V	1.0E+00			1.4E+09	1.5E+03	Dihydroisoflural	94-58-6	7.4E+01		1.4E+00	1.4E+00</					

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	ke y	IUR (ug/m ³ -y) ⁻¹	ke y	RfD _o (mg/kg-day)	ke y	RfC _i (mg/m ³ -y)	ke y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)
1.1E+01	P								1.0E+00	1.0E-01		1.4E+09		Dimethylbenzidine, 3,3'-	119-93-7	3.0E-01	7.0E-01		2.1E-01				
				1.0E-01	P	3.0E-02	I	V	1.0E+00		1.1E+05	1.4E+09	1.3E+05	Dimethylformamide	68-12-2					1.2E+05		1.7E+04	1.5E+04
5.5E+02	C	1.6E-01	C	1.0E-04	X	2.0E-06	X	V	1.0E+00		1.7E+05	1.4E+09	1.6E+05	Dimethylhydrazine, 1,1-	57-14-7					1.2E+02		1.4E+00	1.4E+00
									1.0E+00		1.9E+05	1.4E+09	1.7E+05	Dimethylhydrazine, 1,2-	540-73-8	5.9E-03		1.3E-02	4.1E-03				
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Dimethylphenol, 2,4-	105-67-9					2.3E+04	5.5E+04		1.6E+04
				6.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Dimethylphenol, 2,6-	576-26-1					7.0E+02	1.7E+03		4.9E+02
				1.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Dimethylphenol, 3,4-	95-65-8					1.2E+03	2.8E+03		8.2E+02
4.5E-02	C	1.3E-05	C					V	1.0E+00		1.1E+03	1.4E+09	1.0E+03	Dimethylvinylchloride	513-37-1		9.5E-01	9.4E-01					
				8.0E-05	X				1.0E+00	1.0E-01		1.4E+09		Dinitro-o-cresol, 4,6-	534-52-1					9.3E+01	2.2E+02		6.6E+01
				2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5					2.3E+03	5.5E+03		1.6E+03
				1.0E-04	P				1.0E+00	1.0E-01		1.4E+09		Dinitrobenzene, 1,2-	528-29-0					1.2E+02	2.8E+02		8.2E+01
				1.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Dinitrobenzene, 1,3-	99-65-0					1.2E+02	2.8E+02		8.2E+01
				1.0E-04	P				1.0E+00	1.0E-01		1.4E+09		Dinitrobenzene, 1,4-	100-25-4					1.2E+02	2.8E+02		8.2E+01
				2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Dinitrophenol, 2,4-	51-28-5					2.3E+03	5.5E+03		1.6E+03
6.8E-01	I								1.0E+00	1.0E-01		1.4E+09		Dinitrotoluene Mixture, 2,4/2,6-	NA	4.8E+00	1.1E+01		3.4E+00				
3.1E-01	C	8.9E-05	C	2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Dinitrotoluene, 2,4-	121-14-2	1.1E+01	2.4E+01	1.9E+05	7.4E+00	2.3E+03	5.4E+03		1.6E+03
1.5E+00	P			3.0E-04	X				1.0E+00	9.9E-02		1.4E+09		Dinitrotoluene, 2,6-	606-20-2	2.2E+00	5.2E+00		1.5E+00	3.5E+02	8.4E+02		2.5E+02
				2.0E-03	S				1.0E+00	6.0E-03		1.4E+09		Dinitrotoluene, 2-Amino-4,6-	35572-78-2					2.3E+03	9.2E+04		2.3E+03
				2.0E-03	S				1.0E+00	9.0E-03		1.4E+09		Dinitrotoluene, 4-Amino-2,6-	19406-51-0					2.3E+03	6.1E+04		2.3E+03
4.5E-01	X			9.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Dinitrotoluene, Technical grade	25321-14-6		7.3E+00	1.7E+01	5.1E+00	1.1E+03	2.5E+03		7.4E+02
				1.0E-03	I				1.0E+00	1.0E-01		1.4E+09		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	V			1.0E+00		1.2E+05	1.4E+09	4.0E+04	Dioxane, 1,4-	123-91-1	3.3E+01		9.7E+01	2.4E+01	3.5E+04		5.2E+03	4.5E+03
6.2E+03	I	1.3E+00	I						1.0E+00	3.0E-02		1.4E+09		Dioxins	NA	5.3E-04	4.2E-03	1.3E+01	4.7E-04				
1.3E+05	C	3.8E+01	C	7.0E-10	I	4.0E-08	C	V	1.0E+00	3.0E-02		1.4E+09	2.0E+06	-Hexachlorodibenzo-p-dioxin, Mixture -1,2,3,7,8-	1746-01-8	2.5E-05	2.0E-04	6.3E-04	2.2E-05	8.2E-04	6.4E-03	3.4E-01	7.2E-04
				3.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Diphenamid	957-51-7					3.5E+04	8.3E+04		2.5E+04
				8.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Diphenyl Sulfone	127-63-9					9.3E+02	2.2E+03		6.6E+02
				2.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Diphenylamine	122-39-4					2.9E+04	6.9E+04		2.1E+04
8.0E-01	I	2.2E-04	I						1.0E+00	1.0E-01		1.4E+09		Diphenylhydrazine, 1,2-	122-86-7	4.1E+00	9.7E+00	7.6E+04	2.9E+00				
				2.2E-03	I				1.0E+00	1.0E-01		1.4E+09		Diquat	85-00-7					2.6E+03	6.1E+03		1.8E+03
7.1E+00	C	1.4E-01	C						1.0E+00	1.0E-01		1.4E+09		Direct Black 38	1937-37-1	4.6E-01	1.1E+00	1.2E+02	3.2E-01				
7.4E+00	C	1.4E-01	C						1.0E+00	1.0E-01		1.4E+09		Direct Blue 6	2602-46-2	4.4E-01	1.0E+00	1.2E+02	3.1E-01				
6.7E+00	C	1.4E-01	C						1.0E+00	1.0E-01		1.4E+09		Direct Brown 95	16071-86-6	4.9E-01	1.2E+00	1.2E+02	3.4E-01				
				4.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Disulfoton	298-04-4					4.7E+01	1.1E+02		3.3E+01
				1.0E-02	I			V	1.0E+00		1.4E+09	4.5E+04		Dithane, 1,4-	505-29-3					1.2E+04			1.2E+04
				2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Diuron	330-54-1					2.3E+03	5.5E+03		1.6E+03
				4.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Iodine	20439-10-3					4.7E+03	1.1E+04		3.3E+03
				2.5E-02	I			V	1.0E+00		1.4E+09	1.2E+05		EPTC	759-94-4					2.9E+04			2.9E+04
				6.0E-03	I			V	1.0E+00		1.4E+09	4.1E+05		Endosulfan	115-29-7					7.0E+03			7.0E+03
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Endothal	145-73-3					2.3E+04	5.5E+04		1.6E+04
				3.0E-04	I				1.0E+00	1.0E-01		1.4E+09		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	1.0E+00		1.1E+04	1.4E+09	1.9E+04	Epichlorohydrin	106-89-8	3.3E+02		1.9E+02	1.2E+02	7.0E+03			8.3E+01
				2.0E-02	I	V			1.0E+00		1.5E+04	1.4E+09	7.7E+03	1,2-Epoxybutane, 1,2-	106-88-7							8.3E+01	6.7E+02
				5.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Ethephon	16672-87-0					5.8E+03	1.4E+04		4.1E+03
				5.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Ethion	563-12-2					5.8E+02	1.4E+03		4.1E+02
				1.0E-01	P	6.0E-02	P	V	1.0E+00		3.1E+04	1.4E+09	6.2E+04	Ethoxyethanol Acetate, 2-	111-15-9					1.2E+05		1.6E+04	1.4E+04
				9.0E-02	P	2.0E-01	I	V	1.0E+00		1.1E+05	1.4E+09	9.8E+04	Ethoxyethanol, 2-	110-80-5					1.1E+05		8.6E+04	4.7E+04
				9.0E-01	I	7.0E-02	P	V	1.0E+00		1.1E+04	1.4E+09	8.6E+03	Ethyl Acetate	141-78-6					1.1E+06		2.6E+03	2.6E+03
4.8E-02	H			5.0E-03	P	8.0E-03	P	V	1.0E+00		2.5E+03	1.4E+09	6.3E+03	Ethyl Acrylate	140-88-5	6.8E+01			6.8E+01	5.8E+03			2.1E+02
									1.0E+01	I	V			Ethyl Chloride (Chloroethane)	75-00-3								5.7E+04
				2.0E-01	I			V	1.0E+00		1.0E+04	1.4E+09	3.1E+03	Ethyl Ether	60-29-7					2.3E+05			2.3E+05
				9.0E-02	H	3.0E-01	P	V	1.0E+00		1.1E+03	1.4E+09	5.8E+03	Ethyl Methacrylate	97-63-2					1.1E+05		7.6E+03	7.1E+03
				1.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Ethyl-p-nitrophenyl Phosphonate	2104-64-5					1.2E+01	2.8E+01		8.2E+00
1.1E-02	C	2.5E-06	C	1.0E-01	I	1.0E+00	I	V	1.0E+00		4.8E+02	1.4E+09	5.7E+03	Ethylbenzene	100-41-4	3.0E+02		2.8E+01	2.5E+01	1.2E+05	2.8E+01		2.0E+04
				7.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Ethylene Cyanohydrin	109-78-4					8.2E+04	1.9E+05		5.7E+04
				9.0E-02	P			V	1.0E+00		1.9E+05	1.4E+09	1.8E+05	Ethylene Diamine	107-15-3					1.1E+05			1.1E+05
				2.0E+00																			

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	ke y	IUR (ug/m ³ -y) ⁻¹	ke y	RfD _o (mg/kg-day)	ke y	RfC _i (ug/m ³ -y)	ke y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
				2.5E-02	I				1.0E+00	1.0E-01				Fenpropathrin	39515-41-8					2.9E+04	6.9E+04		2.1E+04	
				1.3E-02	I				1.0E+00	1.0E-01				Fluometuron	2164-17-2					1.5E+04	3.6E+04		1.1E+04	
				4.0E-02	C	1.3E-02	C		1.0E+00					Fluoride	16984-48-8					4.7E+04		7.7E+07	4.7E+04	
				6.0E-02	I	1.3E-02	C		1.0E+00					Fluorine (Soluble Fluoride)	7782-41-4					7.0E+04		7.7E+07	7.0E+04	
				8.0E-02	I				1.0E+00	1.0E-01				Fluridone	59756-60-4					9.3E+04	2.2E+05		6.6E+04	
				2.0E-02	I				1.0E+00	1.0E-01				Flurprimidol	56425-91-3					2.3E+04	5.5E+04		1.6E+04	
				6.0E-02	I				1.0E+00	1.0E-01				Flutolanil	66332-96-5					7.0E+04	1.7E+05		4.9E+04	
				1.0E-02	I				1.0E+00	1.0E-01				Fluvalinate	69409-94-5					1.2E+04	2.8E+04		8.2E+03	
				1.0E-01	I				1.0E+00	1.0E-01				Folpet	133-07-3					1.2E+05	2.8E+05		8.2E+04	
3.5E-03	I															9.3E+02	2.2E+03		6.6E+02					
				1.9E-01	I				1.0E+00	1.0E-01				Fomesafen	72178-02-0				1.2E+01					
						2.0E-03	I		1.0E+00	1.0E-01				Fonofos	944-22-9		1.7E+01	4.1E+01			2.3E+03	5.5E+03		1.6E+03
						1.3E-05	I		1.0E+00		4.2E+04	1.4E+09	7.8E+04	Formaldehyde	50-00-0			7.3E+01	7.3E+01	2.3E+05	2.3E+05	3.3E+03	3.3E+03	
				9.0E-01	P	3.0E-04	X V		1.0E+00		1.1E+05	1.4E+09	9.3E+04	Formic Acid	64-18-6					1.1E+06		1.2E+02	1.2E+02	
				3.0E+00	I				1.0E+00	1.0E-01				Fosetyl-AL	39148-24-8					3.5E+06	8.3E+06		2.5E+06	
														Furans										
				1.0E-03	X		V		1.0E+00	3.0E-02		1.4E+09	2.0E+05	Dibenzofuran	132-64-9					1.2E+03	9.2E+03		1.0E+03	
				1.0E-03	I		V		1.0E+00	3.0E-02	6.2E+03	1.4E+09	2.6E+03	Furan	110-00-9					1.2E+03	9.2E+03		1.0E+03	
				9.0E-01	I	2.0E+00	I V		1.0E+00	3.0E-02	1.7E+05	1.4E+09	1.2E+04	Tetrahydrofuran	109-99-9					1.1E+06	8.3E+06	1.1E+05	9.6E+04	
3.8E+00	H			3.0E-03	I	5.0E-02	H V		1.0E+00	1.0E-01		1.4E+09		Furazolidone	67-45-8	8.6E-01	2.0E+00		6.0E-01					
1.5E+00	C	4.3E-04	C						1.0E+00		1.0E+04	1.4E+09	4.9E+04	Furfural	98-01-1					3.5E+03		1.1E+04	2.6E+03	
				3.0E-02	I	8.6E-06	C		1.0E+00	1.0E-01		1.4E+09		Furium	531-82-8	2.2E+00	5.2E+00	3.9E+04	1.5E+00					
									1.0E+00	1.0E-01		1.4E+09		Furmecyclox	60568-05-0	1.1E+02	2.6E+02	1.9E+06	7.7E+01					
				4.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Glufosinate, Ammonium	77182-82-2					4.7E+02	1.1E+03		3.3E+02	
						8.0E-05	C		1.0E+00	1.0E-01		1.4E+09		Glutaraldehyde	111-30-8					2.3E+05	2.3E+05	4.8E+05	4.8E+05	
				4.0E-04	I	1.0E-03	H V		1.0E+00		1.1E+05	1.4E+09	7.3E+04	Glycidyl	765-34-4					4.7E+02		3.2E+02	1.9E+02	
				1.0E-01	I				1.0E+00	1.0E-01		1.4E+09		Glyphosate	1071-83-6					1.2E+05	2.8E+05		8.2E+04	
				3.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Goal	42874-03-3					3.5E+03	8.3E+03		2.5E+03	
				1.0E-02	X		V		1.0E+00			1.4E+09	1.5E+05	Guandine	143-00-8					1.2E+04			1.2E+04	
				2.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Guandine Chloride	50-01-1					2.3E+04	5.5E+04		1.6E+04	
				3.0E-03	A	1.0E-02	A		1.0E+00	1.0E-01		1.4E+09		Guthion	86-50-0					3.5E+03	8.3E+03	6.0E+07	2.5E+03	
				5.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Haloxyp, Methyl	69806-40-2					5.8E+01	1.4E+02		4.1E+01	
				1.3E-02	I				1.0E+00	1.0E-01		1.4E+09		Harmony	79277-27-3					1.5E+04	3.6E+04		1.1E+04	
4.5E+00	I	1.3E-03	I	5.0E-04	I		V		1.0E+00			4.8E+05		Heptachlor	76-44-8	7.3E-01	4.5E+00	6.3E-01		5.8E+02			5.8E+02	
				1.3E-05	I				1.0E+00		1.4E+09	8.4E+05		Heptachlor Epoxide	1024-57-3					1.5E+01			1.5E+01	
				2.0E-03	I		V		1.0E+00		1.4E+09	3.8E+05		Hexabromobenzene	87-82-1	3.6E-01	4.0E+00	3.3E-01		2.3E+03			2.3E+03	
				2.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BDE-153)	88631-49-2					2.3E+02	5.5E+02		1.6E+02	
1.6E+00	I	4.6E-04	I	8.0E-04	I		V		1.0E+00		1.7E+01	1.4E+09	1.1E+04	Hexachlorobenzene	118-74-1	2.0E+00	1.8E+00	9.6E-01		9.3E+02			9.3E+02	
7.8E-02	I	2.2E-05	I	1.0E-03	P		V		1.0E+00			1.4E+09		Hexachlorobutadiene	87-68-3	4.2E+01	6.0E+00	5.3E+00		1.2E+03			1.2E+03	
6.3E+00	I	1.8E-03	I	8.0E-03	A				1.0E+00	1.0E-01		1.4E+09		Hexachlorocyclohexane, Alpha	319-84-6	5.2E+01	1.2E+00	9.3E+03	3.6E-01	9.3E+03	2.2E+01		6.6E+03	
1.8E+00	I	5.3E-04	I						1.0E+00	1.0E-01		1.4E+09		Hexachlorocyclohexane, Beta-	319-85-7	1.8E+00	4.3E+00	3.1E+04	1.3E+00					
1.1E+00	C	3.1E-04	C	3.0E-04	I				1.0E+00	4.0E-02		1.4E+09		Hexachlorocyclohexane, Gamma-(Lindane)	58-89-9	3.0E+00	1.8E+01	5.4E+04	2.5E+00	3.5E+02	2.1E+03		3.0E+02	
1.8E+00	I	5.1E-04	I						1.0E+00	1.0E-01		1.4E+09		Hexachlorocyclohexane, Technical	608-73-1	1.8E+00	4.3E+00	3.3E+04	1.3E+00					
				6.0E-03	I	2.0E-04	I V		1.0E+00		1.6E+01	1.4E+09	8.5E+03	Hexachlorocyclopentadiene	77-47-4					7.0E+03		7.5E+00	7.5E+00	
				7.0E-04	I	3.0E-02	I V		1.0E+00		1.4E+09	8.0E+03		Hexachloroethane	67-72-1	8.2E+01		8.9E+00	8.0E+00	8.2E+02		1.1E+03	4.6E+02	
				3.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Hexachlorophene	70-30-4					3.5E+02	8.3E+02		2.5E+02	
1.1E-01	I			3.0E-03	I				1.0E+00	1.5E-02		1.4E+09		Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	3.0E+01	4.7E+02		2.8E+01	3.5E+03	5.5E+04		3.3E+03	
				4.0E-04	P	1.0E-05	I V		1.0E+00		5.2E+03	1.4E+09	3.0E+05	Hexamethylene Diisocyanate, 1,6-	822-06-0					4.7E+02	1.1E+03		1.3E+01	
									1.0E+00	1.0E-01		1.4E+09		Hexamethylphosphoramide	680-31-9								3.3E+02	
				6.0E-02	H	7.0E-01	I V		1.0E+00		1.4E+02	1.4E+09	8.3E+02	Hexane, N-	110-54-3					7.0E+04		2.5E+03	2.5E+03	
				2.0E+00	P				1.0E+00	1.0E-01		1.4E+09		Hexanedioic Acid	124-04-9					2.3E+06	5.5E+06		1.6E+06	
				5.0E-03	I	3.0E-02	I V		1.0E+00		3.3E+03	1.4E+09	1.3E+04	Hexanone, 2-	591-78-6					5.8E+03		1.7E+03	1.3E+03	
				3.3E-02	I				1.0E+00	1.0E-01		1.4E+09		Hexazinone	51235-04-2					3.9E+04	9.1E+04		2.7E+04	
3.0E+00	I	4.9E-03	I			3.0E-05	P V		1.0E+00			1.4E+09		Hydrazine	302-01-2	1.1E+00	3.4E+03	1.1E+00				1.8E+05	1.8E+05	
3.0E+00	I	4.9E-03	I						1.0E+00			1.4E+09		Hydrazine Sulfate	10034-93-2	1.1E+00	3.4E+03		1.1E+00					
						2.0E-02	I V		1.0E+00			1.4E+09		Hydrogen Chloride	7647-01-0							1.2E+08	1.2E+08	
				4.0E																				

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Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	ke y	IUR (ug/m ³) ⁻¹	ke y	RfD _o (mg/kg-day)	ke y	RfC _i (mg/m ³)	ke y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
9.5E-04	I			2.0E-01 1.5E-02	I	2.0E+00 C			1.0E+00 1.0E+00	1.0E-01		1.4E+09 4.2E+05		Isophorone Isopropanol	78-59-1 33820-53-0	3.4E+03	8.1E+03		2.4E+03	2.3E+05 1.8E+04	5.5E+05	1.2E+10	1.6E+05 1.8E+04	
				2.0E+00 1.0E-01 5.0E-02	P	2.0E-01 P V			1.0E+00 1.0E+00 1.0E-01	1.1E+05	1.4E+09 2.8E+04			Isopropanol Isopropyl Methyl Phosphonic Acid Isoxaben	67-63-0 1832-54-8 82558-50-7					2.3E+06 1.2E+05 5.8E+04		2.4E+04	2.4E+04 8.2E+04 4.1E+04	
						3.0E-01 A V			1.0E+00 1.0E+00 1.0E-01		1.4E+09 1.4E+09 1.4E+09			JP-7 Kerb Lactofen	NA 23950-58-5 77501-63-4							1.8E+09	1.8E+09 6.2E+04 1.6E+03	
				7.5E-02 2.0E-03	I				1.0E+00 1.0E+00	1.0E-01		1.4E+09 1.4E+09		Lead Compounds --Lead Chromate --Lead Phosphate	7758-97-6 7446-27-7	6.5E+00 3.8E+02		1.1E+02 1.4E+06	6.2E+00 3.8E+02	2.3E+04		1.2E+06	2.3E+04	
5.0E-01 8.5E-03	C C	1.5E-01 1.2E-05	C C	2.0E-02 2.0E-02	C	2.0E-04 C	M		2.5E-02 1.0E+00		1.4E+09 1.4E+09			--Lead acetate --Lead and Compounds --Lead subacetate	301-04-2 7439-92-1 1335-32-6	1.2E+01 3.8E+02	2.8E+01 9.1E+02	2.1E+05 1.4E+06	8.2E+00 2.7E+02				8.0E+02	
				1.0E-07 2.0E-03 2.0E-03	I P		V		1.0E+00 1.0E+00 1.0E+00	2.4E+00	1.4E+09 1.4E+09 1.4E+09	1.9E+03		--Tetraethyl Lead Linuron Lithium	78-00-2 330-55-2 7439-93-2					1.2E-01 2.3E+03 2.3E+03	5.5E+03		1.6E+03 2.3E+03	
				2.0E-01 5.0E-04 1.0E-02	I I I				1.0E+00 1.0E+00 1.0E+00	1.0E-01	1.4E+09 1.4E+09 1.4E+09			Londax MCPA MCPB	83055-99-6 94-74-6 94-81-5					2.3E+05 5.8E+02 1.2E+04	5.5E+05 1.4E+03 2.8E+04		1.6E+05 4.1E+02 8.2E+03	
				1.0E-03 2.0E-02 1.0E-01	I I I				1.0E+00 1.0E+00 1.0E+00	1.0E-01	1.4E+09 1.4E+09 1.4E+09			MCPB Malathion Malic Anhydride	93-65-2 121-75-5 198-31-6					1.2E+03 2.3E+04 1.2E+05	2.8E+03 5.5E+04 2.8E+05	4.2E+06	8.2E+02 1.6E+04 8.0E+04	
				5.0E-01 1.0E-04 3.0E-02	I P H				1.0E+00 1.0E+00 1.0E+00	1.0E-01	1.4E+09 1.4E+09 1.4E+09			Maleic Hydrazide Malononitrile Mancozeb	123-33-1 109-77-3 8018-01-7					5.8E+05 1.2E+02 3.5E+04	1.4E+06 2.8E+02 8.3E+04		4.1E+05 8.2E+01 2.5E+04	
				5.0E-03 1.4E-01 2.4E-02	I I S				1.0E+00 1.0E+00 4.0E-02	1.0E-01	1.4E+09 1.4E+09 1.4E+09			Maneb Manganese (Diet) Manganese (Ndn-diet)	12427-38-2 7439-96-5 7439-96-5					5.8E+03 2.8E+04	1.4E+04	3.0E+05	4.1E+03 2.6E+04	
				9.0E-05 3.0E-02	H I				1.0E+00 1.0E+00	1.0E-01	1.4E+09 1.4E+09			Mephosolan Mepiquat Chloride Mercury Compounds	950-10-7 24307-26-4					1.1E+02 3.5E+04	2.5E+02 8.3E+04		7.4E+01 2.5E+04	
				3.0E-04 1.0E-04	I I	3.0E-04 S 3.0E-04 I V			7.0E-02 1.0E+00 1.0E+00		1.4E+09 1.4E+09 1.4E+09	3.0E+04		--Mercuric Chloride (and other Mercury salts) --Mercury (elemental) --Methyl Mercury	7487-94-7 7439-97-6 22967-92-6					3.5E+02 1.2E+02		1.8E+06 4.0E+01	3.5E+02 4.0E+01 1.2E+02	
				8.0E-05 3.0E-05 3.0E-05	I I I		V		1.0E+00 1.0E+00 1.0E+00	1.0E-01	1.4E+09 1.4E+09 1.4E+09	1.9E+06		--Phenylmercuric Acetate Merphos Merphos Oxide	62-38-4 150-50-5 78-48-8					9.3E+01 3.5E+01 3.5E+01	2.2E+02 8.3E+01		6.6E+01 3.5E+01 2.5E+01	
				6.0E-02 1.0E-04 5.0E-05	I I I				1.0E+00 1.0E+00 1.0E+00	1.0E-01	1.4E+09 1.4E+09 1.4E+09	6.8E+03		Metaxyl Methacrylonitrile Methamidophos	57837-19-1 128-98-7 10265-92-6					7.0E+04 1.2E+02 5.8E+01	1.7E+05 8.9E+02		4.9E+04 1.0E+02 4.1E+01	
				2.0E+00 1.0E-03 2.5E-02	I I I	2.0E+01 I V			1.0E+00 1.0E+00 1.0E+00	1.1E+05	1.4E+09 1.4E+09 1.4E+09	2.9E+04		Methanol Methidathion Methomyl	67-56-1 950-37-8 16752-77-5					2.3E+06 1.2E+03 2.9E+04	2.5E+06 2.8E+03 6.9E+04		1.2E+06 8.2E+02 2.1E+04	
4.9E-02	C	1.4E-05	C	5.0E-03 8.0E-03	I P	1.0E-03 P V			1.0E+00 1.0E+00 1.2E+05	1.0E-01	1.4E+09 1.4E+09 1.4E+09	1.2E+05		Methoxy-5-nitroaniline, 2- Methoxychlor Methoxyethanol Acetate, 2- Methoxyethanol, 2- Methyl Acetate Methyl Acrylate	99-59-2 72-43-5 110-49-6 109-86-4 79-20-9 96-33-3	6.7E+01	1.6E+02	1.2E+06	4.7E+01	5.8E+03 9.3E+03	1.4E+04	5.4E+02	4.1E+03 5.1E+02	
				5.0E-03 1.0E+00 3.0E-02	P X H	2.0E-02 I V 2.0E-02 P V			1.0E+00 1.0E+00 1.0E+00	1.1E+05 2.9E+04 6.8E+03	1.4E+09 1.4E+09 1.4E+09	1.0E+05 8.1E+03 7.0E+03		Methyl Ethyl Ketone (2-Butanone) Methyl Hydrazine Methyl Isobutyl Ketone (4-methyl-2-pentanone)	78-93-3 60-34-4 108-10-1			1.9E+00	1.9E+00			7.0E+05 1.2E+03 9.3E+04	2.7E+05 1.4E+01 1.4E+05	1.9E+05 1.4E+01 5.6E+04
				1.4E+00 2.5E-04	I I	7.0E-01 I V			1.0E+00 1.0E+00	1.0E-01	1.4E+09 1.4E+09	6.3E+03		Methyl Isocyanate Methyl Methacrylate Methyl Parathion	624-83-9 80-62-6 298-00-0					1.6E+06 2.9E+02	6.9E+02		1.9E+04 2.1E+02	
				6.0E-02 9.9E-02	X C	4.0E-02 H V			1.0E+00 1.0E+00	1.0E-01	1.4E+09 1.4E+09	1.2E+04		Methyl Phosphonic Acid Methyl Styrene (Mixed Isomers) Methyl methanesulfonate	993-13-5 25013-15-4 66-27-3				2.3E+01	7.0E+04 7.0E+03	1.7E+05	2.1E+03	4.9E+04 1.6E+03	
1.8E-03	C	2.6E-07	C	3.0E+00 3.0E-04	I X	3.0E+00 I V			1.0E+00 1.0E+00	1.0E-01	8.9E+03 1.4E+09	4.9E+03		Methyl tert-Butyl Ether (MTBE) Methyl-1,4-benzenediamine dihydrochloride, 2- Methyl-5-Nitroaniline, 2- Methyl-N-nitro-N-nitrosoguanidine, N- Methylaniline Hydrochloride, 2-	1634-04-4 615-45-2 99-55-8 70-25-7 636-21-5	1.8E+03		2.3E+02	2.1E+02	3.5E+02 2.3E+04	8.3E+02 5.5E+04	6.4E+04	6.4E+04 2.5E+02 1.6E+04	

Regional Screening Level (RSL) Composite Worker Soil Table (TR=1E-6, HQ=1) June 2015

Key: I = IRIS; P = PPRTR; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1								
SFO (mg/kg-day) ⁻¹	ke y ⁻¹	IUR (ug/m ³ -y) ⁻¹	ke y ⁻¹	RfD _o (mg/kg-day)	ke y ⁻¹	RfC _o (mg/m ³ -y)	ke y ⁻¹	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
				1.0E-02	A				1.0E+00	1.0E-01		1.4E+09		Methylarsonic acid	124-58-3					1.2E+04	2.8E+04		8.2E+03	
1.0E-01	X			2.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Methylbenzene, 1,4-diamine monohydrochloride, 2-	74612-12-7					2.3E+02	5.5E+02		1.6E+02	
2.2E+01	C	6.3E-03	C	3.0E-04	X			M	1.0E+00	1.0E-01		1.4E+09		Methylbenzene, 1,4-diamine sulfate, 2-	615-50-9	3.3E+01	7.7E+01		2.3E+01	3.5E+02	8.3E+02		2.5E+02	
									1.0E+00	1.0E-01		1.4E+09		Methylcholanthrene, 3-	56-49-5	1.5E-01	3.5E-01	2.6E+03	1.0E-01					
2.0E-03	I	1.0E-08	I	6.0E-03	I	6.0E-01	I	V	M	1.0E+00		3.3E+03	1.4E+09	2.2E+03	Methylene Chloride	75-09-2	1.6E+03	2.7E+03	1.0E+03	7.0E+03		5.8E+03	3.2E+03	
1.0E-01	P	4.3E-04	C	2.0E-03	P			M	1.0E+00	1.0E-01		1.4E+09		Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	3.3E+01	7.7E+01	3.9E+04	2.3E+01	2.3E+03	5.5E+03		1.6E+03	
4.6E-02	I	1.3E-05	C						1.0E+00	1.0E-01		1.4E+09		Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	7.1E+01	1.7E+02	1.3E+06	5.0E+01					
1.6E+00	C	4.6E-04	C			2.0E-02	C		1.0E+00	1.0E-01		1.4E+09		Methylenedisobenzeneamine, 4,4'-	101-77-9	2.0E+00	4.8E+00	3.6E+04	1.4E+00			1.2E+08	1.2E+08	
				7.0E-02	H	6.0E-04	I		1.0E+00	1.0E-01		1.4E+09		Methylenediphenyl Diisocyanate	101-68-8							3.6E+06	3.6E+06	
								V	1.0E+00		5.0E+02	1.4E+09	1.3E+04	Methylstyrene, Alpha-	98-83-9					8.2E+04			8.2E+04	
				1.5E-01	I				1.0E+00	1.0E-01		1.4E+09		Metolachlor	51218-45-2					1.8E+05	4.1E+05		1.2E+05	
				2.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Metribuzin	21087-64-9					2.9E+04	6.9E+04		2.1E+04	
				3.0E+00	P			V	1.0E+00		3.4E-01	1.4E+09	1.4E+03	Mineral oils	8012-95-1					3.5E+06			3.5E+06	
1.8E+01	C	5.1E-03	C	2.0E-04	I			V	1.0E+00			1.4E+09	8.6E+05	Mirex	2385-85-5	1.8E-01		2.1E+00	1.7E-01	2.3E+02			2.3E+02	
				2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Molinate	2212-67-1					2.3E+03	5.5E+03		1.6E+03	
				5.0E-03	I				1.0E+00			1.4E+09		Molybdenum	7439-98-7					5.8E+03			5.8E+03	
				1.0E-01	I				1.0E+00			1.4E+09		Monochloramine	10599-90-3					1.2E+05			1.2E+05	
				2.0E-03	P				1.0E+00	1.0E-01		1.4E+09		Monomethylaniline	100-61-8					2.3E+03	5.5E+03		1.6E+03	
				3.0E-04	X				1.0E+00	1.0E-01		1.4E+09		N,N-Diphenyl-1,4-benzenediamine	74-31-7					3.5E+02	8.3E+02		2.5E+02	
				2.0E-03	I			V	1.0E+00			1.4E+09	5.7E+04	Naled	300-76-5					2.3E+03			2.3E+03	
1.8E+00	C	0.0E+00	C	3.0E-02	X	1.0E-01	P	V	1.0E+00			1.4E+09		Naphtha, High Flash Aromatic (HFAN)	64742-95-6	1.8E+00	4.3E+00		1.3E+00	3.5E+04		6.0E+08		3.5E+04
				1.0E-01	I				1.0E+00	1.0E-01		1.4E+09		Naphthylamine, 2-	91-59-8									
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1.0E+00	1.0E-01	1.4E+09		Napropamide	15299-99-7					1.2E+05	2.8E+05		8.2E+04	
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1.0E+00	1.0E-01	1.4E+09		Nickel Acetate	373-02-4		6.4E+04	6.4E+04		1.3E+04	3.0E+04	8.3E+04	8.1E+03	
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1.0E+00	1.0E-01	1.4E+09		Nickel Carbonate	3333-87-3		6.4E+04	6.4E+04		1.3E+04	3.0E+04	8.3E+04	8.1E+03	
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1.0E+00		1.4E+09		Nickel Carbonyl	13463-39-3		6.4E+04	6.4E+04		1.3E+04		8.3E+04	1.1E+04	
				2.6E-04	C	1.1E-02	C	1.4E-05	C	4.0E-02		1.4E+09		Nickel Hydroxide	12054-48-7		6.4E+04	6.4E+04		1.3E+04		8.3E+04	1.1E+04	
				2.6E-04	C	1.1E-02	C	2.0E-05	C	4.0E-02		1.4E+09		Nickel Oxide	1313-99-1		6.4E+04	6.4E+04		1.3E+04		1.2E+05	1.2E+04	
				2.4E-04	I	1.1E-02	C	1.4E-05	C	4.0E-02		1.4E+09		Nickel Refinery Dust	NA		6.9E+04	6.9E+04		1.3E+04		8.3E+04	1.1E+04	
				2.6E-04	C	2.0E-02	I	9.0E-05	A	4.0E-02		1.4E+09		Nickel Soluble Salts	7440-02-0		6.4E+04	6.4E+04		2.3E+04		4.5E+05	2.2E+04	
1.7E+00	C	4.8E-04	I	1.1E-02	C	1.4E-05	C		4.0E-02		1.4E+09		Nickel Sulfide	12036-72-2	1.9E+00	3.5E+04	1.9E+00		1.3E+04		8.3E+04	1.1E+04		
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1.0E+00	1.0E-01	1.4E+09		Nickelocene	1271-28-9		6.4E+04	6.4E+04		1.3E+04	3.0E+04	8.3E+04	8.1E+03	
				1.6E+00	I				1.0E+00		1.4E+09		Nitrate	14797-55-8					1.9E+06			1.9E+06		
									1.0E+00		1.4E+09		Nitrate + Nitrite (as N)	NA										
				1.0E-01	I				1.0E+00		1.4E+09		Nitrite	14797-65-0					1.2E+05			1.2E+05		
				1.0E-02	X	5.0E-05	X		1.0E+00	1.0E-01		1.4E+09		Nitroamine, 2-	88-74-4					1.2E+04	2.8E+04	3.0E+05	8.0E+03	
				4.0E-03	P	6.0E-03	P		1.0E+00	1.0E-01		1.4E+09		Nitroamine, 4-	100-01-6	1.6E+02	3.9E+02		1.1E+02	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	1.0E+00	3.1E+03	1.4E+09	7.3E+04	Nitrobenzene	98-95-3		2.2E+01	2.2E+01	2.3E+03		2.9E+03	1.3E+03	
				3.0E+03	P				1.0E+00	1.0E-01		1.4E+09		Nitrocellulose	9004-70-0					3.5E+09	8.3E+09		2.5E+09	
				7.0E-02	H				1.0E+00	1.0E-01		1.4E+09		Nitrofurantoin	67-20-9					8.2E+04	1.9E+05		5.7E+04	
1.3E+00	C	3.7E-04	C						1.0E+00	1.0E-01		1.4E+09		Nitrofurazone	59-87-0	2.5E+00	5.9E+00	4.5E+04	1.8E+00	1.2E+02	2.8E+02		8.2E+01	
1.7E-02	P			1.0E-04	P				1.0E+00	1.0E-01		1.4E+09		Nitroglycerin	55-63-0	1.9E+02	4.5E+02		1.4E+02	1.2E+02	2.8E+02		8.2E+01	
				1.0E-01	I				1.0E+00	1.0E-01		1.4E+09		Nitroguanidine	556-88-7					1.2E+05	2.8E+05		8.2E+04	
				8.8E-06	P	5.0E-03	P	V	1.0E+00	1.8E+04	1.4E+09	1.7E+04		Nitromethane	75-52-5		2.4E+01	2.4E+01		3.7E+02			3.7E+02	
2.7E+01	C	7.7E-03	C	2.7E-03	H	2.0E-02	I	V	1.0E+00	4.9E+03	1.4E+09	1.3E+04		Nitropropane, 2	79-46-9		6.0E-02	6.0E-02				1.2E+03	1.2E+03	
				2.7E-03	H				1.0E+00	1.0E-01		1.4E+09		Nitroso-N-ethylurea, N-	759-73-9	1.2E-01	2.9E-01	2.2E+03	8.5E-02					
1.2E+02	C	3.4E-02	C					M	1.0E+00	1.0E-01		1.4E+09		Nitroso-N-methylurea, N-	684-93-5	2.7E-02	6.4E-02	4.9E+02	1.9E-02					
5.4E+00	I	1.6E-03	I					V	1.0E+00			1.4E+09	2.4E+05	Nitroso-di-N-butylamine, N-	924-16-3	6.1E-01		1.9E+00	4.6E-01					
7.0E+00	I	2.0E-03	C						1.0E+00	1.0E-01		1.4E+09		Nitroso-di-N-propylamine, N-	621-64-7	4.7E-01	1.1E+00	8.3E+03	3.3E-01					
2.8E+00	I	8.0E-04	C						1.0E+00	1.0E-01		1.4E+09		Nitrosodiphenylamine, N-	1116-54-7	1.2E+00	2.8E+00	2.1E+04	8.2E-01					
1.5E+02	I	4.3E-02	I					M	1.0E+00	1.0E-01		1.4E+09		Nitrosodimethylamine, N-	55-18-5	2.2E-02	5.2E-02	3.9E+02	1.5E-02					
5.1E+01	I	1.4E-02	I	8.0E-06	P	4.0E-05	X	V	M	1.0E+00	2.4E+05	1.4E+09	8.2E+04	Nitrosodimethylamine, N-	62-75-9	6.4E-02	7.2E-02	3.4E-02		9.3E+00		1.4E+01	5.7E+00	
4.9E-03	I	2.6E-06	C						1.0E+00	1.0E-01		1.4E+09		Nitrosodiphenylamine, N-	86-30-6	6.7E+02	1.6E+03	6.4E+06	4.7E+02					
2.2E+01	I	6.3E-03	C					V	1.0E+00		1.1E+05	1.4E+09	1.2											

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Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	ke y	IUR (ug/m ³ -y) ⁻¹	ke y	RfD _o (mg/kg-day)	ke y	RfC _i (mg/m ³)	ke y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
				5.0E-02	I				1.0E+00	6.0E-03		1.4E+09		Octahydro-1,3,5,7-tetrahydro-1,3,5,7-tetrazocine (HMX)	2691-41-0					5.8E+04	2.3E+06		5.7E+04	
				2.0E-03	H				1.0E+00	1.0E-01		1.4E+09		Octamethylpyrophosphoramide	152-16-9					2.3E+03	5.5E+03		1.6E+03	
				5.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Oryzalin	19044-88-3					5.8E+04	1.4E+05		4.1E+04	
				5.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Oxadiazon	19666-30-9					5.8E+03	1.4E+04		4.1E+03	
				2.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Oxamyl	23135-22-0					2.9E+04	6.9E+04		2.1E+04	
				1.3E-02	I				1.0E+00	1.0E-01		1.4E+09		Paclitaxel	76738-62-0					1.5E+04	3.6E+04		1.1E+04	
				4.5E-03	I				1.0E+00	1.0E-01		1.4E+09		Paraquat Dichloride	1910-42-5					5.3E+03	1.2E+04		3.7E+03	
				6.0E-03	H				1.0E+00	1.0E-01		1.4E+09		Parathion	56-38-2					7.0E+03	1.7E+04		4.9E+03	
				5.0E-02	H			V	1.0E+00			4.5E+04		Pebulate	1114-71-2					5.8E+04			5.8E+04	
				4.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Pendimethalin	40487-42-1					4.7E+04	1.1E+05		3.3E+04	
				2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Pentabromodiphenyl Ether	32534-81-9					2.3E+03	5.5E+03		1.6E+03	
				1.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Pentabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-99)	60348-60-9					1.2E+02	2.8E+02		8.2E+01	
				8.0E-04	I			V	1.0E+00			8.1E+04		Pentachlorobenzene	608-93-5					9.3E+02			9.3E+02	
9.0E-02	P							V	1.0E+00		4.5E+02	1.4E+09	9.7E+03	Pentachloroethane	76-01-7	3.6E+01			3.6E+01					
2.6E-01	H			3.0E-03	I			V	1.0E+00			4.3E+05	Pentachloronitrobenzene	82-68-8	1.3E+01				1.3E+01	3.5E+03			3.5E+03	
4.0E-01	I	5.1E-06	C	5.0E-03	I				1.0E+00	2.5E-01		1.4E+09	Pentachlorophenol	87-86-5	8.2E+00	7.7E+00	3.3E+06	4.0E+00	5.8E+03	5.5E+03			2.8E+03	
4.0E-03	X			2.0E-03	P				1.0E+00	1.0E-01		1.4E+09	Pentaerythritol tetranitrate (PETN)	78-11-5	8.2E+02	1.9E+03		5.7E+02	2.3E+03	5.5E+03			1.6E+03	
						1.0E+00	P	V	1.0E+00		3.9E+02	1.4E+09	7.8E+02	Pentane, n-	109-68-0							3.4E+03	3.4E+03	
														Perchlorates										
				7.0E-04	I				1.0E+00			1.4E+09		--Ammonium Perchlorate	7790-98-9					8.2E+02				8.2E+02
				7.0E-04	I				1.0E+00			1.4E+09		--Lithium Perchlorate	7791-03-9					8.2E+02				8.2E+02
				7.0E-04	I				1.0E+00			1.4E+09		--Perchlorate and Perchlorate Salts	14797-73-0					8.2E+02				8.2E+02
				7.0E-04	I				1.0E+00			1.4E+09		--Potassium Perchlorate	7778-74-7					8.2E+02				8.2E+02
				7.0E-04	I				1.0E+00			1.4E+09		--Sodium Perchlorate	1601-89-0					8.2E+02				8.2E+02
				2.0E-02	P			V	1.0E+00			1.4E+09	1.3E+05	Perfluorobutane Sulfonate	375-73-5					2.3E+04				2.3E+04
2.2E-03	C	6.3E-07	C	5.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Permethrin	52645-53-1	1.5E+03	3.5E+03	2.6E+07	1.0E+03	5.8E+04	1.4E+05			4.1E+04
				2.5E-01	I				1.0E+00	1.0E-01		1.4E+09		Phenmedipham	13884-63-4					2.9E+05	6.9E+05			2.1E+05
				3.0E-01	I	2.0E-01	C		1.0E+00	1.0E-01		1.4E+09		Phenol	108-95-2					3.5E+05	8.3E+05	1.2E+09		2.5E+05
				5.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Phenothiazine	92-84-2					5.8E+02	1.4E+03			4.1E+02
				6.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Phenylenediamine, m-	108-45-2					7.0E+03	1.7E+04			4.9E+03
				1.9E-01	H				1.0E+00	1.0E-01		1.4E+09		Phenylenediamine, o-	95-54-5	7.0E+01	1.6E+02		4.9E+01					
				1.9E-01	H				1.0E+00	1.0E-01		1.4E+09		Phenylenediamine, p-	106-50-3					2.2E+05	5.2E+05			1.6E+05
1.9E-03	H			2.0E-04	H				1.0E+00	1.0E-01		1.4E+09		Phenylphenol-2	90-43-7	1.7E+03	4.0E+03		1.2E+03					
				3.0E-04	I	V			1.0E+00	1.0E-01	1.6E+03	1.4E+09	9.8E+02	Phorate	298-02-2					2.3E+02	5.5E+02	1.3E+00		1.6E+02
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Phosgene	75-44-5									1.3E+00
				4.9E+01	P				1.0E+00			1.4E+09		Phosmet	732-11-6					2.3E+04	5.5E+04			1.6E+04
				4.9E+01	P				1.0E+00			1.4E+09		Phosphates, Inorganic										
				4.9E+01	P				1.0E+00			1.4E+09		--Aluminum metaphosphate	13176-88-0					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Ammonium polyphosphate	68333-79-9					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Calcium pyrophosphate	7790-76-3					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Diammonium phosphate	7783-28-0					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Dicalcium phosphate	7757-93-9					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Dimagnesium phosphate	7782-75-4					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Dipotassium phosphate	7758-11-4					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Disodium phosphate	7558-79-4					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monoaluminum phosphate	13530-50-2					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monoammonium phosphate	7722-76-1					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monocalcium phosphate	7758-23-8					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monomagnesium phosphate	7757-86-0					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monopotassium phosphate	7778-77-0					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monosodium phosphate	7558-80-7					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Polyphosphoric acid	8017-16-1					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Potassium triphosphate	13845-36-8					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium acid pyrophosphate	7758-16-9					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium aluminum phosphate (acidic)	7785-88-8					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium aluminum phosphate (anhydrous)	10279-59-1					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium aluminum phosphate (tetrahydrate)	10305-76-7					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium hexametaphosphate	10124-56-8					5.7E+07				5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium polyphosphate	68915-31-1					5.7E+07				

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; J = New Jersey; O = EPA Office of Water; F = See FAQ; E = Environmental Criteria and Assessment Office; S = see user guide Section 5; L = see user guide on lead; M = mutagen; V = volatile; R = RBA applied (See User Guide for Arsenic notice); c = cancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; n = noncancer; m = Concentration may exceed ceiling limit (See User Guide); s = Concentration may exceed Csat (See User Guide); SSL values are based on DAF=1

Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1										
SFO (mg/kg-day) ⁻¹	ke y	IUR (ug/m ³ -y) ⁻¹	ke y	RfD _o (mg/kg-day)	ke y	RfC _i (mg/m ³)	ke y	ke y	ke y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
4.9E+01	P			4.9E+01	P						1.0E+00			1.4E+09		--Trialuminum sodium tetra decahydrogenoctaorthophosphate (dihydrate)	15136-87-5					5.7E+07				5.7E+07
4.9E+01	P			4.9E+01	P						1.0E+00			1.4E+09		--Tricalcium phosphate	7758-87-4					5.7E+07				5.7E+07
4.9E+01	P			4.9E+01	P						1.0E+00			1.4E+09		--Trimagnesium phosphate	7757-87-1					5.7E+07				5.7E+07
4.9E+01	P			4.9E+01	P						1.0E+00			1.4E+09		--Tripotassium phosphate	7778-53-2					5.7E+07				5.7E+07
4.9E+01	P			4.9E+01	P						1.0E+00			1.4E+09		--Trisodium phosphate	7801-54-9					5.7E+07				5.7E+07
3.0E-04	I	3.0E-04	I	3.0E-04	I	V					1.0E+00			1.4E+09		Phosphine	7803-51-2					3.5E+02		1.8E+06		3.5E+02
4.9E+01	P	1.0E-02	I	4.9E+01	P						1.0E+00			1.4E+09		Phosphoric Acid	7664-38-2					5.7E+07		6.0E+07		2.9E+07
2.0E-05	I		V	2.0E-05	I						1.0E+00			6.9E+03		Phosphorus, White	7723-14-0					2.3E+01				2.3E+01
1.4E-02	I	2.4E-06	C	2.0E-02	I						1.0E+00	1.0E-01		1.4E+09		Phthalates						2.3E+04	5.5E+04			1.6E+04
				1.0E+00	I						1.0E+00	1.0E-01		1.4E+09		--Bis(2-ethylhexyl)phthalate	117-81-7	2.3E+02	5.5E+02	6.9E+06	1.6E+02	1.2E+06	2.8E+06			8.2E+05
				1.0E-01	I						1.0E+00	1.0E-01		1.4E+09		--Butylphthalyl Butylglycolate	85-70-1					1.2E+05	2.8E+05			8.2E+04
				8.0E-01	I						1.0E+00	1.0E-01		1.4E+09		--Diethyl Phthalate	84-74-2					9.3E+05	2.2E+05			6.6E+05
				1.0E-01	I		V				1.0E+00			2.1E+04		--Dimethylterephthalate	120-61-6					1.2E+05				1.2E+05
				1.0E-02	P						1.0E+00	1.0E-01		1.4E+09		--Octyl Phthalate, di-N-	117-84-0					1.2E+04	2.8E+04			8.2E+03
				1.0E+00	H						1.0E+00	1.0E-01		1.4E+09		--Phthalic Acid, P-	100-21-0					1.2E+06	2.8E+06			8.2E+05
				2.0E+00	I	2.0E-02	C				1.0E+00	1.0E-01		1.4E+09		--Phthalic Anhydride	85-44-9					2.3E+06	5.5E+06	1.2E+08		1.6E+06
				7.0E-02	I						1.0E+00	1.0E-01		1.4E+09		Picloram	1918-02-1					8.2E+04	1.9E+05			5.7E+04
				1.0E-04	X						1.0E+00	1.0E-01		1.4E+09		Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3					1.2E+02	2.8E+02			8.2E+01
				1.0E-02	I						1.0E+00	1.0E-01		1.4E+09		Pirimphos, Methyl	29232-93-7					1.2E+04	2.8E+04			8.2E+03
3.0E+01	C	8.6E-03	C	7.0E-06	H						1.0E+00	1.0E-01		1.4E+09		Polybrominated Biphenyls	59536-65-1	1.1E-01	2.6E-01	1.9E+03	7.7E-02	8.2E+00	1.9E+01			5.7E+00
7.0E-02	S	2.0E-05	S	7.0E-05	I		V				1.0E+00	1.4E-01		1.4E+09	5.9E+05	Polychlorinated Biphenyls (PCBs)						8.2E+01	1.4E+02			5.1E+01
				2.0E+00	S	5.7E-04	S				1.0E+00	1.4E-01		1.4E+09	1.1E+05	--Aroclor 1016	12674-11-2	4.7E+01	7.9E+01	3.6E+02	2.7E+01					5.1E+01
				2.0E+00	S	5.7E-04	S				1.0E+00	1.4E-01		1.4E+09	1.1E+05	--Aroclor 1221	11104-28-2	1.6E+00	2.8E+00	2.4E+00	7.2E-01					
				2.0E+00	S	5.7E-04	S				1.0E+00	1.4E-01		1.4E+09	1.1E+05	--Aroclor 1232	11141-16-5	1.6E+00	2.8E+00	2.4E+00	7.2E-01					
				2.0E+00	S	5.7E-04	S				1.0E+00	1.4E-01		1.4E+09	7.9E+05	--Aroclor 1242	53489-21-9	1.6E+00	2.8E+00	1.7E+01	9.7E-01					
				2.0E+00	S	5.7E-04	S				1.0E+00	1.4E-01		1.4E+09	5.1E+05	--Aroclor 1248	12672-29-6	1.6E+00	2.8E+00	1.7E+01	9.4E-01					
				2.0E+00	S	5.7E-04	S	2.0E-05	I		1.0E+00	1.4E-01		1.4E+09	8.4E+05	--Aroclor 1254	11097-69-1	1.6E+00	2.8E+00	1.8E+01	9.7E-01	2.3E+01	3.9E+01			1.5E+01
				2.0E+00	S	5.7E-04	S				1.0E+00	1.4E-01		1.4E+09	1.3E+06	--Aroclor 1280	11096-82-5	1.6E+00	2.8E+00	2.8E+01	9.9E-01					
				6.0E-04	X		V				1.0E+00	1.4E-01		1.4E+09	7.2E+05	--Aroclor 5460	11126-42-4	8.4E-01	1.4E+00	2.2E+01	5.1E-01	7.0E+02	1.2E+03			4.4E+02
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V			1.0E+00	1.4E-01		1.4E+09	2.0E+06	--Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	8.4E-01	1.4E+00	1.5E+01	5.1E-01	2.7E+01	4.6E+01	1.2E+04		1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V			1.0E+00	1.4E-01		1.4E+09	1.4E+06	--Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167)	52663-72-6	8.4E-01	1.4E+00	1.5E+01	5.1E-01	2.7E+01	4.6E+01	8.4E+03		1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V			1.0E+00	1.4E-01		1.4E+09	1.5E+06	--Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	8.4E-01	1.4E+00	1.6E+01	5.1E-01	2.7E+01	4.6E+01	8.5E+03		1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V			1.0E+00	1.4E-01		1.4E+09	1.5E+06	--Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 156)	38380-08-4	8.4E-01	1.4E+00	1.7E+01	5.1E-01	2.7E+01	4.6E+01	9.0E+03		1.7E+01
3.9E+03	E	1.1E+00	E	2.3E-08	E	1.3E-06	E	V			1.0E+00	1.4E-01		1.4E+09	1.4E+06	--Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 189)	32174-16-6	8.4E-04	1.4E-03	1.5E-02	5.1E-04	2.7E-02	4.6E-02	8.4E+00		1.7E-02
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V			1.0E+00	1.4E-01		1.4E+09	1.0E+06	--Pentachlorobiphenyl, 2,3,4,4',5'- (PCB 123)	65510-44-3	8.4E-01	1.4E+00	1.1E+01	5.0E-01	2.7E+01	4.6E+01	6.0E+03		1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V			1.0E+00	1.4E-01		1.4E+09	8.3E+05	--Pentachlorobiphenyl, 2,3',4',4',5'- (PCB 118)	31508-00-6	8.4E-01	1.4E+00	8.9E+00	5.0E-01	2.7E+01	4.6E+01	4.8E+03		1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V			1.0E+00	1.4E-01		1.4E+09	8.5E+05	--Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	8.4E-01	1.4E+00	9.1E+00	5.0E-01	2.7E+01	4.6E+01	4.9E+03		1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V			1.0E+00	1.4E-01		1.4E+09	1.0E+06	--Pentachlorobiphenyl, 2,3,4,4',5'- (PCB 114)	74472-37-0	8.4E-01	1.4E+00	1.1E+01	5.0E-01	2.7E+01	4.6E+01	6.0E+03		1.7E+01
1.3E+04	E	3.8E+00	E	7.0E-09	E	4.0E-07	E	V			1.0E+00	1.4E-01		1.4E+09	1.0E+06	--Pentachlorobiphenyl, 3,3',4,4',5'- (PCB 126)	57485-28-8	2.5E+04	4.2E+04	3.3E-03	1.5E-04	8.2E-03	1.4E-02	1.8E+00		5.1E-03
2.0E+00	I	5.7E-04	I				V				1.0E+00	1.4E-01		1.4E+09	7.9E+05	--Polychlorinated biphenyls (high risk)	1336-36-3	1.6E+00	2.8E+00	1.7E+01	9.7E-01					
4.0E-01	I	1.0E-04	I				V				1.0E+00	1.4E-01		1.4E+09		--Polychlorinated biphenyls (low risk)	1336-36-3									
7.0E-02	I	2.0E-05	I				V				1.0E+00	1.4E-01		1.4E+09		--Polychlorinated biphenyls (lowest risk)	1336-36-3									
1.3E+01	E	3.8E-03	E	7.0E-06	E	4.0E-04	E				1.0E+00	1.4E-01		1.4E+09		--Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	2.5E-01	4.2E-01	4.4E+03	1.6E-01	8.2E+00	1.4E+01	2.4E+06		5.1E+00
3.9E+01	E	1.1E-02	E	2.3E-06	E	1.3E-04	E	V			1.0E+00	1.4E-01		1.4E+09	7.3E+05	--Tetrachlorobiphenyl, 3,4,4',5'- (PCB 81)	70382-50-4	8.4E-02	1.4E-01	7.8E-01	4.9E-02	2.7E+00	4.6E+00	4.2E+02		1.7E+00
				6.0E-04	I						1.0E+00	1.0E-01		1.4E+09		--Polymerc Methylen Diphenyl Diisocyanate (PMDI)	9016-87-9									3.6E+06
				6.0E-02	I		V				1.0E+00	1.3E-01		1.4E+09	1.4E+05	Polynuclear Aromatic Hydrocarbons (PAHs)										
				3.0E-01	I		V				1.0E+00	1.3E-01														

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Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1											
SFO (mg/kg-day) ⁻¹	ke (y)	IUR (ug/m ³) ⁻¹	ke (y)	RfD _o (mg/kg-day)	ke (y)	RfC _i (mg/m ³)	ke (y)	muta-gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)					
3.4E-05	C	2.0E-02	I	3.0E-03	I	V	1.0E+00	1.3E-01	1.0E+00	1.3E-01	1.4E+09	4.6E+04	-Naphthalene	91-20-3	1.7E+01	1.7E+01	2.3E+04	4.2E+04	6.1E+02	5.9E+02								
1.2E+00	C	1.1E-04	C	3.0E-02	I	V	1.0E+00	1.3E-01	1.0E+00	1.3E-01	1.4E+09	2.4E+06	-Nitropyrene, 4- -Pyrene Potassium Perfluorobutane Sulfonate	57835-92-4 129-00-0 29420-49-3	2.7E+00	5.0E+00	1.5E+05	1.8E+00	3.5E+04	6.4E+04	2.3E+04	5.5E+04	2.3E+04	1.6E+04				
1.5E-01	I	9.0E-03	I	6.0E-03	H	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	4.2E+05	Prochloraz Profuralin Prometon	67747-09-5 26399-36-0 1610-18-0	2.2E+01	5.2E+01	1.5E+01	1.1E+04	2.5E+04	7.4E+03	7.0E+03	1.8E+04	4.1E+04	7.0E+03	1.2E+04			
4.0E-03	I	1.3E-02	I	5.0E-03	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Prometryn Propachlor Propanil	7287-19-6 1918-16-7 709-98-8					4.7E+03	1.1E+04	3.6E+04	1.5E+03	1.4E+04	3.3E+03	1.1E+04	4.1E+03		
2.0E-02	I	2.0E-03	I	2.0E-02	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	6.3E+04	Propargite Propargyl Alcohol Propazine	2312-35-8 107-19-7 139-40-2					2.3E+04	5.5E+04	2.3E+03	2.3E+04	5.5E+04	1.6E+04	2.3E+03	1.6E+04		
2.0E-02	I	1.3E-02	I	8.0E-03	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	8.9E+03	Propam Propiconazole Propionaldehyde	122-42-9 60207-90-1 123-38-6					2.3E+04	5.5E+04	1.5E+04	3.6E+04	3.1E+02	1.6E+04	1.1E+04	3.1E+02		
1.0E-01	X	3.0E+00	C	1.0E+00	X	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	2.6E+02	7.0E+03	Propyl benzene Propylene Propylene Glycol	103-65-1 115-07-1 57-55-6					1.2E+05	3.1E+04	9.3E+03	1.6E+07	3.1E+04	2.4E+04	9.3E+03	1.6E+07		
2.4E-01	I	3.7E-06	I	3.0E-02	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	7.8E+04	1.4E+09	1.0E+04	Propylene Glycol Dinitrate Propylene Glycol Monoethyl Ether Propylene Glycol Monomethyl Ether	6423-43-4 1569-02-4 107-98-2					8.2E+05	8.2E+05	6.9E+05	1.6E+06	6.9E+05	3.7E+05	1.6E+06	3.7E+05	
2.4E-01	I	3.7E-06	I	3.0E-02	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	7.8E+04	1.4E+09	1.0E+04	Propylene Oxide Pursut Pydrin	75-56-9 81335-77-5 51630-58-1	1.4E+01	3.4E+01	9.7E+00	2.9E+05	2.9E+04	6.9E+04	1.4E+03	2.1E+05	2.1E+04	2.1E+05	2.1E+04		
3.0E+00	I	5.0E-04	I	1.0E+00	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	5.3E+05	1.4E+09	5.5E+04	Pyridine Quinalphos Quinoline	110-86-1 3593-03-8 91-22-5	1.1E+00	2.6E+00	7.7E-01	1.2E+03	5.8E+02	1.4E+03	4.1E+02	1.2E+03	4.1E+02	1.2E+03	4.1E+02		
2.2E-01	C	6.3E-05	C	3.0E-02	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	4.7E+05	Refractory Ceramic Fibers Resmethrin Ronnel	NA 10463-86-8 299-84-3					3.5E+04	8.3E+04	1.8E+08	1.8E+08	1.8E+08	1.8E+08	1.8E+08	1.8E+08		
2.2E-01	C	6.3E-05	C	4.0E-03	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Rotenone Safrole Saxby	83-79-4 94-59-7 78587-05-0	1.5E+01	3.5E+01	2.6E+05	1.0E+01	4.7E+03	1.1E+04	2.9E+04	6.9E+04	2.1E+04	3.3E+03	3.3E+03			
5.0E-03	I	5.0E-03	I	5.0E-03	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Selenious Acid Selenium Selenium Sulfide	7783-00-8 7782-49-2 446-34-6					5.8E+03	5.8E+03	5.8E+03	1.2E+08	1.2E+08	5.8E+03	5.8E+03			
1.2E-01	H	5.0E-03	I	1.3E-02	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Sethoxydim Silica (crystalline, respirable) Silver	74051-80-2 7631-86-9 7440-22-4	2.7E+01	6.4E+01	1.9E+01	5.8E+03	1.5E+04	3.6E+04	4.7E+03	1.1E+05	2.5E+05	1.8E+07	5.8E+03	7.4E+04	1.8E+07	5.8E+03
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M	2.5E-02	1.0E+00	1.4E+09	1.4E+09	Sodium Dichromate Sodium Diethylthiocarbamate Sodium Fluoride	10699-81-0 148-18-5 7681-49-4	8.5E+00	1.1E+02	6.2E+00	2.3E+04	3.5E+04	8.3E+04	7.7E+07	1.2E+06	2.3E+04	2.3E+04	2.3E+04	2.3E+04		
2.7E-01	H	5.0E-02	A	3.0E-02	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Sodium Fluoroacetate Sodium Metavanadate Stirofos (Tetrachlorovinphos)	62-74-8 13718-26-8 961-11-5	1.4E+02	3.2E+02	9.6E+01	2.3E+01	5.5E+01	8.3E+04	7.7E+07	1.2E+06	2.3E+04	2.3E+04	2.3E+04			
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M	2.5E-02	1.0E+00	1.4E+09	1.4E+09	Strontium Chromate Strontium, Stable Strychnine	7789-06-2 7440-24-6 57-24-9	6.5E+00	1.1E+02	6.2E+00	2.3E+04	7.0E+05	3.5E+02	8.3E+02	1.2E+06	2.3E+04	2.3E+04	2.3E+04	2.3E+04		
2.0E-01	I	1.0E+00	I	1.3E-02	P	2.0E-03	X	1.0E+00	1.0E-01	1.0E+00	1.4E+09	9.4E+03	Styrene Styrene-Acrylonitrile (SAN) Trimer Sulfolane	100-42-5 NA 126-33-0				2.3E+05	3.5E+03	8.3E+03	1.2E+07	4.1E+04	3.5E+04	2.5E+03	8.2E+02			
8.0E-04	P	1.0E-03	C	1.0E+00	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Sulfonylbis(4-chlorobenzene), 1,1'- Sulfur Trioxide Sulfuric Acid	80-07-9 7446-11-9 7664-93-9				9.3E+02	2.2E+03	6.0E+06	6.0E+06	6.0E+06	6.0E+06	6.0E+06	6.0E+06			
2.5E-02	I	3.0E-02	H	7.0E-02	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Systhane TCMTB Tebuthiuron	88671-89-0 21564-17-0 34014-18-1				2.9E+04	3.5E+04	8.3E+04	1.9E+05	6.9E+04	8.3E+04	5.7E+04	2.1E+04	2.5E+04	5.7E+04	
2.0E-02	H	1.3E-02	I	2.5E-05	H	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	3.1E+01	1.4E+09	2.6E+05	Temephos Terbacil Terbufos	3383-96-8 5902-51-2 13071-79-9				2.3E+04	1.5E+04	3.6E+04	2.9E+01	5.5E+04	3.6E+04	1.8E+04	1.1E+04	2.9E+01	

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Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	ke y	IUR (ug/m ³ -y) ⁻¹	ke y	RfD _o (mg/kg-day)	ke y	RfC _i (mg/m ³ -y)	ke y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)
				1.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Terbutryn	886-50-0					1.2E+03	2.8E+03		8.2E+02
				1.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1					1.2E+02	2.8E+02		8.2E+01
				3.0E-04	I			V	1.0E+00			1.4E+09	5.1E+04	Tetrachlorobenzene, 1,2,4,5-	95-94-3					3.5E+02			3.5E+02
2.6E-02	I	7.4E-06	I	3.0E-02	I			V	1.0E+00		6.8E+02	1.4E+09	5.7E+03	Tetrachloroethane, 1,1,1,2-	630-20-6	1.3E+02		9.4E+00	8.8E+00	3.5E+04			3.5E+04
2.0E-01	I	5.8E-05	C	2.0E-02	I			V	1.0E+00		1.9E+03	1.4E+09	1.5E+04	Tetrachloroethane, 1,1,2,2-	79-34-5	1.6E+01		2.7E+00		2.3E+04			2.3E+04
2.1E-03	I	2.6E-07	I	6.0E-03	I	4.0E-02	I	V	1.0E+00		1.7E+02	1.4E+09	2.4E+03	Tetrachloroethylene	127-18-4	1.6E+03		1.1E+02	1.0E+02	7.0E+03		4.1E+02	3.9E+02
2.0E+01	H			3.0E-02	I			V	1.0E+00	1.0E-01		1.4E+09		Tetrachlorophenol, 2,3,4,6-	58-90-2				1.6E-01	3.5E+04	8.3E+04		2.5E+04
				5.0E-04	I			V	1.0E+00	1.0E-01		1.4E+09	1.1E+05	Tetrachlorotoluene, p- alpha, alpha- Tetraethyl Dithiopyrophosphate	5216-25-1 3689-24-5				1.6E-01	5.8E+02	1.4E+03		4.1E+02
				8.0E+01	I			V	1.0E+00		1.1E+03	1.4E+09	1.2E+03	Tetrafluoroethane, 1,1,1,2-	811-97-2							4.3E+05	
				2.0E-03	P				1.0E+00	6.5E-04		1.4E+09		Tetryl (Trinitrophenylmethyl)nitramine)	479-45-8					2.3E+03	8.5E+05		2.3E+03
				7.0E-06	X				1.0E+00			1.4E+09		Thallium (I) Nitrate	10102-45-1					8.2E+00			8.2E+00
				1.0E-05	X				1.0E+00			1.4E+09		Thallium (Soluble Salts)	7440-28-0					1.2E+01			1.2E+01
				6.0E-06	X			V	1.0E+00	1.0E-01		1.4E+09		Thallium Acetate	563-68-8					7.0E+00	1.7E+01		4.9E+00
				2.0E-05	X				1.0E+00	1.0E-01		1.4E+09		Thallium Carbonate	6533-73-9					2.3E+01	5.5E+01		1.6E+01
				6.0E-06	X				1.0E+00			1.4E+09		Thallium Chloride	7791-12-0					7.0E+00			7.0E+00
				2.0E-05	X				1.0E+00			1.4E+09		Thallium Sulfate	7446-18-6					2.3E+01			2.3E+01
				1.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Thiobencarb	28249-77-6					1.2E+04	2.8E+04		8.2E+03
				7.0E-02	X				1.0E+00	7.5E-03		1.4E+09		Thiodiglycol	111-48-8					8.2E+04	2.6E+06		7.9E+04
				3.0E-04	H				1.0E+00	1.0E-01		1.4E+09		Thiofanox	39196-18-4					3.5E+02	8.3E+02		2.5E+02
				8.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Thiophanate, Methyl	23564-05-8					9.3E+04	2.2E+05		6.6E+04
				5.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Thiram	137-26-8					5.8E+03	1.4E+04		4.1E+03
				6.0E-01	H				1.0E+00			1.4E+09		Tin	7440-31-5					7.0E+05			7.0E+05
				1.0E-04	A	V			1.0E+00			1.4E+09		Titanium tetrachloride	7550-45-0							6.0E+05	6.0E+05
1.8E-01	X			8.0E-02	I	5.0E+00	I	V	1.0E+00		8.2E+02	1.4E+09	4.3E+03	Toluene	108-88-3					9.3E+04		9.4E+04	4.7E+04
3.0E-02	P			2.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Toluene-2,5-diamine	95-70-5	1.8E+01	4.3E+01		1.3E+01	2.3E+02	5.5E+02		1.6E+02
				4.0E-03	X				1.0E+00	1.0E-01		1.4E+09		Toluidine, p-	106-49-0	1.1E+02	2.6E+02		7.7E+01	4.7E+03	1.1E+04		3.3E+03
				3.0E+00	P			V	1.0E+00		3.4E-01	1.4E+09	1.1E+03	Total Petroleum Hydrocarbons (Aliphatic High)	NA					3.5E+06			3.5E+06
				6.0E-01	P	V			1.0E+00		1.4E+02	1.4E+09	8.3E+02	Total Petroleum Hydrocarbons (Aliphatic Low)	NA						2.2E+03		2.2E+03
				1.0E-02	X	1.0E-01	P	V	1.0E+00		6.9E+00	1.4E+09	1.0E+03	Total Petroleum Hydrocarbons (Aliphatic Medium)	NA					1.2E+04		4.6E+02	4.4E+02
				4.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Total Petroleum Hydrocarbons (Aromatic High)	NA					4.7E+04	1.1E+05		3.3E+04
				4.0E-03	P	3.0E-02	P	V	1.0E+00		1.8E+03	1.4E+09	3.5E+03	Total Petroleum Hydrocarbons (Aromatic Low)	NA					4.7E+03		4.6E+02	4.2E+02
				4.0E-03	P	3.0E-03	P	V	1.0E+00		1.4E+09	5.2E+04		Total Petroleum Hydrocarbons (Aromatic Medium)	NA					4.7E+03		6.9E+02	6.0E+02
1.1E+00	I	3.2E-04	I						1.0E+00	1.0E-01		1.4E+09		1,2,4-trioxaphene	8001-35-2	3.0E+00	7.0E+00	5.2E+04	2.1E+00				
				7.5E-03	I				1.0E+00	1.0E-01		1.4E+09		Trialomethrin	89841-25-6					8.8E+03	2.1E+04		6.2E+03
				3.0E-04	A			V	1.0E+00			1.4E+09	3.4E+03	Tri-n-butyltin	898-73-3					3.5E+02			3.5E+02
				8.0E+01	X				1.0E+00	1.0E-01		1.4E+09		Triacetin	102-76-1					9.3E+07	2.2E+08		6.6E+07
				1.3E-02	I			V	1.0E+00			1.4E+09	3.6E+05	Triallate	2303-17-5					1.5E+04			1.5E+04
				1.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Inasuturon	82091-50-5					1.2E+04	2.8E+04		8.2E+03
9.0E-03	P			5.0E-03	I			V	1.0E+00			1.4E+09	4.5E+04	Tribromobenzene, 1,2,4-	615-54-3	3.6E+02	8.6E+02		2.6E+02	5.8E+03			5.8E+03
				1.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Tributyl Phosphate	126-73-8					1.2E+04	2.8E+04		8.2E+03
				3.0E-04	P				1.0E+00	1.0E-01		1.4E+09		Tributyltin Compounds	NA					3.5E+02	8.3E+02		2.5E+02
				3.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Tributyltin Oxide	56-35-9					3.5E+02	8.3E+02		2.5E+02
7.0E-02	I			3.0E+01	I	3.0E+01	H	V	1.0E+00		9.1E+02	1.4E+09	1.3E+03	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1	4.7E+01	1.1E+02			3.5E+07		1.7E+05	1.7E+05
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Trichloroacetic Acid	76-03-9					2.3E+04	5.5E+04		1.6E+04
2.9E-02	H								1.0E+00	1.0E-01		1.4E+09		Trichloroaniline HCl, 2,4,6-	33663-50-2	1.1E+02	2.7E+02		7.9E+01				
7.0E-03	X			3.0E-05	X				1.0E+00	1.0E-01		1.4E+09		Trichloroaniline, 2,4,6-	634-93-5	4.7E+02	1.1E+03		3.3E+02	3.5E+01	8.3E+01		2.5E+01
				8.0E-04	X				1.0E+00			1.4E+09	3.2E+04	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	1.0E+00		4.0E+02	1.4E+09	3.0E+04	Trichlorobenzene, 1,2,4-	120-82-1	1.1E+02			1.1E+02	1.2E+04		2.6E+02	2.6E+02
				2.0E+00	I	5.0E+00	I	V	1.0E+00		6.4E+02	1.4E+09	1.7E+03	Trichloroethane, 1,1,1-	71-55-6					2.3E+06			3.6E+04
5.7E-02	I	1.6E-05	I	4.0E-03	I	2.0E-04	X	V	1.0E+00		2.2E+03	1.4E+09	7.2E+03	Trichloroethane, 1,1,2-	79-00-5	5.7E+01		5.5E+00	5.0E+00	4.7E+03			6.3E+00
4.6E-02	I	4.1E-06	I	5.0E-04	I	2.0E-03	I	V	1.0E+00		6.9E+02	1.4E+09	2.2E+03	Trichloroethylene	79-01-6	7.1E+01		6.6E+00	6.0E+00	5.8E+02		1.9E+01	1.9E+01
				3.0E-01	I	7.0E-01	H	V	1.0E+00		1.2E+03	1.4E+09	1.0E+03	Trichlorofluoromethane	75-69-4					3.5E+05		3.2E+03	3.1E+03
				1.0E-01	I				1.0E+00	1.0E-01		1.4E+09		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				1.0E+00	1.0E-01		1.4E+09		Trichlorophenol, 2,4,6-	88-06-2	3.0E+02	7.0E+02	5.4E+06	2.1E+02	1.2E+03	2.8E+03		8.2E+02
				1.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5								

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Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1									
SFO (mg/kg-day) ⁻¹	Key ¹	IUR (ug/m ³ -y) ⁻¹	Key ²	RfD _o (mg/kg-day)	Key ³	RfC _i (mg/m ³)	Key ⁴	Vol ⁵	muta-gen	GIABS	ABS	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
7.7E-03 2.0E-02	I P			7.5E-03 1.0E-02	I P			V		1.0E+00 1.0E+00	1.0E-01		1.4E+09 1.4E+09	5.1E+05	Trifluralin Trimethyl Phosphate	1582-09-8 512-56-1	4.2E+02 1.6E+02			4.2E+02 1.1E+02	8.8E+03 1.2E+04				8.8E+03 8.2E+03
				5.0E-03 7.0E-03 1.0E-02	P V X			V		1.0E+00 1.0E+00 1.0E+00		2.9E+02 2.2E+02 1.8E+02	1.4E+09 1.4E+09 1.4E+09	9.4E+03 7.9E+03 6.6E+03	Trimethylbenzene, 1,2,3- Trimethylbenzene, 1,2,4- Trimethylbenzene, 1,3,5-	526-73-8 95-63-6 108-67-8					1.2E+04			2.1E+02 2.4E+02 1.2E+04	
3.0E-02	I			3.0E-02 5.0E-04 2.0E-02	I I P					1.0E+00 1.0E+00 1.0E+00	1.9E-02 3.2E-02 1.0E-01		1.4E+09 1.4E+09 1.4E+09		Trinitrobenzene, 1,3,5- Trinitrotoluene, 2,4,6- Triphenylphosphine Oxide	99-35-4 118-96-7 791-28-6	1.1E+02	8.0E+02		9.6E+01	3.5E+04 5.8E+02 2.3E+04	4.4E+05 4.3E+03 5.5E+04		3.2E+04 5.1E+02 1.6E+04	
2.3E+00	C	6.6E-04	C	2.0E-02 1.0E-02	A X			V		1.0E+00 1.0E+00	1.0E-01	4.7E+02	1.4E+09 1.4E+09	9.0E+05	Tris(1,3-Dichloro-2-propyl) Phosphate Tris(1-chloro-2-propyl)phosphate Tris(2,3-dibromopropyl)phosphate	13674-87-8 13674-84-5 126-72-7	1.4E+00		1.7E+01	1.3E+00	2.3E+04 1.2E+04	5.5E+04 2.8E+04		1.6E+04 8.2E+03	
2.0E-02 3.2E-03	P P			7.0E-03 1.0E-01 3.0E-03	P P I					1.0E+00 1.0E+00 1.0E+00	1.0E-01		1.4E+09 1.4E+09 1.4E+09		Tris(2-chloroethyl)phosphate Tris(2-ethylhexyl)phosphate Uranium (Soluble Salts)	144-98-8 78-42-2 NA	1.6E+02 1.0E+03	3.9E+02 2.4E+03		1.1E+02 7.2E+02	8.2E+03 1.2E+05 3.5E+03	1.9E+04 2.8E+05		5.7E+03 8.2E+04 3.5E+03	
1.0E+00	C	2.9E-04 8.3E-03	C P	9.0E-03 1.0E+00	I H	7.0E-06 2.0E-01	P I		M	2.6E-02 2.6E-02	1.0E-01		1.4E+09 1.4E+09		Urethane Vanadium Pentoxide Vanadium and Compounds	51-79-6 1314-62-1 7440-62-2	3.3E+00	7.7E+00	5.7E+04 2.0E+03	2.3E+00 2.0E+03	1.1E+04 5.9E+03		4.2E+04 6.0E+05	8.4E+03 5.8E+03	
				1.0E-03 2.5E-02 1.0E+00	I I H			V		1.0E+00 1.0E+00 1.0E+00	1.0E-01	2.8E+03	1.4E+09 1.4E+09 4.4E+03	1.2E+05	Vermolate Vinclozolin Vinyl Acetate	1929-77-7 50471-44-8 108-05-4					1.2E+03 2.9E+04 1.2E+06	6.9E+04		1.2E+03 2.1E+04 3.8E+03	
7.2E-01	I	4.4E-06	I	3.0E-03 3.0E-04	I I	1.0E-01 1.0E-01	I V	M		1.0E+00 1.0E+00		3.4E+03 3.9E+03	1.4E+09 1.4E+09	1.4E+03 9.6E+02	Vinyl Bromide Vinyl Chloride Warfarin	593-60-2 75-01-4 81-81-2	4.5E+00		5.2E-01 2.7E+00	5.2E-01 1.7E+00	3.5E+03 3.5E+02		1.8E+01 4.2E+02 8.3E+02	1.8E+01 3.7E+02 2.5E+02	
				2.0E-01 2.0E-01 2.0E-01	S S S	1.0E-01 1.0E-01 1.0E-01	S S S	V		1.0E+00 1.0E+00 1.0E+00		3.9E+02 3.9E+02 4.3E+02	1.4E+09 1.4E+09 1.4E+09	5.6E+03 5.5E+03 6.5E+03	Xylene, p- Xylene, m- Xylene, o-	106-42-3 108-38-3 95-47-6				2.3E+05 2.3E+05 2.3E+05		2.4E+03 2.4E+03 2.8E+03	2.4E+03 2.4E+03 2.8E+03		
				2.0E-01 3.0E-04 3.0E-01	I I I	1.0E-01 1.0E+00 1.0E+00	I I I	V		1.0E+00 1.0E+00 1.0E+00		2.6E+02	1.4E+09 1.4E+09	6.5E+03	Xylenes Zinc Phosphide Zinc and Compounds	1330-20-7 1314-84-7 7440-66-6					2.3E+05 3.5E+02 3.5E+05		2.8E+03	2.8E+03 3.5E+02 3.5E+05	
				5.0E-02 8.0E-05	I X					1.0E+00 1.0E+00	1.0E-01		1.4E+09 1.4E+09		Zinc Zirconium	12122-67-7 7440-67-7					5.8E+04 9.3E+01	1.4E+05		4.1E+04 9.3E+01	

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Toxicity and Chemical-specific					Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1	
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y	v o l a t i l e	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
5.1E-06	C					ALAR	1596-84-5	2.4E+00	
2.2E-06	I	9.0E-03	I	V		Acephate	30560-19-1		
						Acetaldehyde	75-07-0	5.6E+00	3.9E+01
		3.1E+01	A	V		Acetochlor	34256-82-1		
		2.0E-03	X	V		Acetone	67-64-1		1.4E+05
						Acetone Cyanohydrin	75-86-5		8.8E+00
		6.0E-02	I	V		Acetonitrile	75-05-8		2.6E+02
1.3E-03	C			V		Acetophenone	98-86-2		
						Acetylaminofluorene, 2-	53-96-3	9.4E-03	
		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
		1.0E-03	I	V		Acrylic Acid	79-10-7		4.4E+00
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-03	
		1.0E-04	X	V		Allyl	74223-64-6		
						Allyl Alcohol	107-18-6		4.4E-01
6.0E-06	C	1.0E-03	I	V		Allyl Chloride	107-05-1	2.0E+00	4.4E+00
		5.0E-03	P			Aluminum	7429-90-5		2.2E+01
						Aluminum Phosphide	20859-73-8		
						Amdro	67485-29-4		
6.0E-03	C					Ametryn	834-12-8	2.0E-03	
						Aminobiphenyl, 4-	92-67-1		
						Aminophenol, m-	591-27-5		
						Aminophenol, p-	123-30-8		
						Amitraz	38089-61-1		
1.0E-01	I	V				Ammonia	7664-41-7		4.4E+02
		3.0E-03	X	V		Ammonium Sulfamate	7773-06-0		1.3E+01
						Amyl Alcohol, tert-	75-95-4		
1.6E-06	C	1.0E-03	I			Aniline	62-53-3	7.7E+00	4.4E+00
						Anthraquinone, 9,10-	84-65-1		
						Antimony (metallic)	7440-36-0		
						Antimony Pentoxide	1314-60-9		
						Antimony Potassium Tartrate	11071-15-1		
						Antimony Trioxide	1332-81-6		
		2.0E-04	I			Antimony Trioxide	1309-64-4		8.8E-01
7.1E-06	I					Apollo	74115-24-5	1.7E+00	
						Aramite	140-57-8		
4.3E-03	I	1.5E-05	C			Arsenic, Inorganic	7440-38-2	2.9E-03	6.6E-02
		5.0E-05	I			Arsine	7784-42-1		2.2E-01
						Assure	76578-14-8		
						Asulam	3337-71-1		
2.5E-04	C					Atrazine	1912-24-9	4.9E-02	
						Auramine	492-80-8		
						Avermectin B1	65195-55-3		
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
1.5E-01	C	2.0E-04	C		M	Barium Chromate	10294-40-3	8.2E-05	8.8E-01
						Baygon	114-26-1		
						Bayleton	43121-43-3		
						Baythroid	68359-37-5		
						Benefin	1861-40-1		
						Benomyl	17804-35-2		
						Bentazon	25057-89-0		
						Benzaldehyde	100-52-7		
7.8E-06	I	3.0E-02	I	V		Benzene	71-43-2	1.6E+00	1.3E+02
						Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1		
						Benzenethiol	108-98-5		
6.7E-02	I				M	Benzydine	92-87-5	1.8E-04	
						Benzoic Acid	65-85-0		
						Benzotrithloride	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-03	8.8E-02
						Bidrin	141-66-2		
						Bifenox	42576-02-3		
						Biphenthrin	82657-04-3		
		4.0E-04	X	V		Biphenyl, 1,1'-	92-52-4		1.8E+00
1.0E-05	H			V		Bis(2-chloro-1-methylethyl) ether	108-60-1	1.2E+00	
						Bis(2-chloroethoxy)methane	111-91-1		

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Toxicity and Chemical-specific					Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1	
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y c	v o l a t i l e	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
3.3E-04	I			V		Bis(2-chloroethyl)ether	111-44-4	3.7E-02	
6.2E-02	I			V		Bis(chloromethyl)ether	542-88-1	2.0E-04	
						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-04	X			V		Bromo-2-chloroethane, 1-	107-04-0	2.0E-02	
		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+01	
		5.0E-03	I	V		Bromomethane	74-83-9		2.2E+01
				V		Bromophos	2104-96-3		
				V		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		
		3.0E+01	P	V		Butyl Benzyl Phthlate	85-68-7		
				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+02	
				V		Butylated hydroxytoluene	128-37-0		
				V		Butylbenzene, n-	104-51-8		
				V		Butylbenzene, sec-	135-98-8		
				V		Butylbenzene, tert-	98-06-6		
				V		Cacodylic Acid	75-60-5		
1.8E-03	I	1.0E-05	A			Cadmium (Diet)	7440-43-9		
1.8E-03	I	1.0E-05	A			Cadmium (Water)	7440-43-9	6.8E-03	4.4E-02
1.5E-01	C	2.0E-04	C		M	Calcium Chromate	13765-19-0	8.2E-05	8.8E-01
		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					Captaol	133-06-2	1.9E+01	
						Carbaryl	62-25-2		
						Carbaryl	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+00	4.4E+02
						Carbosulfan	55285-14-8		
						Carboxin	5234-68-4		
		9.0E-04	I			Chloral Hydrate	1306-38-3		3.9E+00
				V		Chloral Hydrate	302-17-0		
						Chloramben	133-90-4		
						Chloranil	118-75-2		
1.0E-04	I	7.0E-04	I	V		Chlordane	12789-03-6	1.2E-01	3.1E+00
4.6E-03	C					Chlordecone (Kepone)	143-50-0	2.7E-03	
						Chlorfenvinphos	470-90-6		
		1.5E-04	A	V		Chlorimuron, Ethyl-	90982-32-4		6.4E-01
						Chlorine	7782-50-5		
		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-02	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		
		3.0E-05	I			Chloroacetic Acid	79-11-8		1.3E-01
						Chloroacetophenone, 2-	532-27-4		
		5.0E-02	P	V		Chloroaniline, p-	106-47-8		
3.1E-05	C					Chlorobenzene	108-90-7	4.0E-01	2.2E+02
						Chlorobenzilate	510-15-6		
		3.0E-01	P	V		Chlorobenzoic Acid, p-	74-11-3		
				V		Chlorobenzotrifluoride, 4-	98-56-6		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-02	
		1.0E-05	X			Chloronitrobenzene, o-	88-73-3		4.4E-02
		6.0E-04	P			Chloronitrobenzene, p-	100-00-5		2.6E+00
				V		Chlorophenol, 2-	95-57-8		
		4.0E-04	C	V		Chloropicrin	76-06-2		1.8E+00
8.9E-07	C					Chlorothalonil	1897-45-6	1.4E+01	
				V		Chlorotoluene, o-	95-49-8		
				V		Chlorotoluene, p-	106-43-4		

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Toxicity and Chemical-specific					Contaminant	Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1	
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y	v o l a t i l e	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
6.9E-02	C				Chlorozotocin Chlorpropham Chlorpyrifos	54749-90-5 101-21-3 2921-88-2	1.8E-04	
					Chlorpyrifos Methyl Chlorsulfuron Chlorthiophos	5598-13-0 64902-72-3 60238-56-4		
8.4E-02	S	1.0E-04	I		Chromium(III), Insoluble Salts Chromium(VI) Chromium, Total	16065-83-1 18540-29-9 7440-47-3	1.5E-04	4.4E-01
9.0E-03	P	6.0E-06	P		Cobalt	7440-48-4	1.4E-03	2.6E-02
6.2E-04	I		V	M	Coke Oven Emissions Copper	8007-45-2 7440-50-8	2.0E-02	
		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
		6.0E-01	C		Cresol, p-chloro-m-	59-50-7		2.6E+03
			V		Cresols Crotonaldehyde, trans-	1319-77-3 123-73-9		
6.3E-05	C	4.0E-01	I	V	Cumene Cupferron Cyanazine	98-82-8 135-20-6 21725-46-2	1.9E-01	1.8E+03
					Cyanides ~Calcium Cyanide ~Copper Cyanide	592-01-8 544-92-3		
8.0E-04	S		V		~Cyanide (CN-) ~Cyanogen ~Cyanogen Bromide	57-12-5 460-19-5 506-68-3		3.5E+00
			V		~Cyanogen Chloride ~Hydrogen Cyanide ~Potassium Cyanide	506-77-4 74-90-8 151-50-8		3.5E+00
8.0E-04	I		V		~Potassium Silver Cyanide ~Silver Cyanide ~Sodium Cyanide	506-61-6 506-64-9 143-33-9		
			V		~Thiocyanates ~Thiocyanic Acid ~Zinc Cyanide	NA 483-56-9 557-21-1		
6.0E+00	I		V		Cyclohexane Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro- Cyclohexanone	110-82-7 87-84-3 108-94-1		2.6E+04 3.1E+03
7.0E-01	P		V		Cyclohexene Cyclohexylamine Cyhalothrin/karate	110-83-8 108-91-8 68085-85-8		4.4E+03
1.0E+00	X		V		Cypermethrin Cyromazine DDD	52315-07-8 66215-27-8 72-54-8		
6.9E-05	C				DDE, p,p'- DDT Dacthal	72-55-9 50-29-3 1861-32-1	1.8E-01	
9.7E-05	C		V		Dalapon Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209) Demeton	75-99-0 1163-19-5 8065-48-3	1.3E-01 1.3E-01	
9.7E-05	I				Di(2-ethylhexyl)adipate Diallate Diazinon	103-23-1 2303-16-4 333-41-5		
6.0E-03	P	2.0E-04	I	V	Dibenzothiophene Dibromo-3-chloropropane, 1,2- Dibromobenzene, 1,3-	132-65-0 96-12-8 108-36-1	2.0E-03	8.8E-01
2.7E-05	C		V		Dibromobenzene, 1,4- Dibromochloromethane Dibromoethane, 1,2-	106-37-6 124-48-1 106-93-4	4.5E-01 2.0E-02	3.9E+01
6.0E-04	I	9.0E-03	I	V	Dibromomethane (Methylene Bromide) Dibutyltin Compounds Dicamba	74-95-3 NA 1918-00-9		1.8E+01
4.2E-03	P		V		Dichloro-2-butene, 1,4- Dichloro-2-butene, cis-1,4- Dichloro-2-butene, trans-1,4-	764-41-0 1476-11-5 110-57-6	2.9E-03 2.9E-03 2.9E-03	
2.0E-01	H		V		Dichloroacetic Acid Dichlorobenzene, 1,2- Dichlorobenzene, 1,4-	79-43-6 95-50-1 106-46-7		8.8E+02 3.5E+03
1.1E-05	C	8.0E-01	I	V	Dichlorobenzidine, 3,3'- Dichlorobenzophenone, 4,4'- Dichlorodifluoromethane	91-94-1 90-98-2 75-71-8	3.6E-02	4.4E+02
3.4E-04	C		X	V	Dichloroethane, 1,1- Dichloroethane, 1,2- Dichloroethylene, 1,1-	75-34-3 107-06-2 75-35-4	7.7E+00 4.7E-01	3.1E+01 8.8E+02

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Toxicity and Chemical-specific					Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1	
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y	v o l a t i l e	mutagen	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
						Dichloroethylene, 1,2-cis-	156-59-2		
						Dichloroethylene, 1,2-trans-	156-60-5		
						Dichlorophenol, 2,4-	120-83-2		
1.0E-05	C	4.0E-03	I	V		Dichlorophenoxy Acetic Acid, 2,4-	94-75-7	1.2E+00	1.8E+01
						Dichlorophenoxybutyric Acid, 4-(2,4-	94-82-6		
						Dichloropropane, 1,2-	78-87-5		
4.0E-06	I	2.0E-02	I	V		Dichloropropane, 1,3-	142-28-9	3.1E+00	8.8E+01
						Dichloropropanol, 2,3-	616-23-9		
						Dichloropropene, 1,3-	542-75-6		
8.3E-05	C	5.0E-04	I			Dichlorvos	62-73-7	1.5E-01	2.2E+00
4.6E-03	I	3.0E-04	X	V		Dicyclopentadiene	77-73-6		
						Dieldrin	60-57-1	2.7E-03	1.3E+00
3.0E-04	C	5.0E-03	I			Diesel Engine Exhaust	NA	4.1E-02	2.2E+01
						Diethanolamine	111-42-2		
						Diethylene Glycol Monobutyl Ether	112-34-5		
						Diethylene Glycol Monoethyl Ether	111-90-0		
1.0E-01	C					Diethylformamide	617-84-5	1.2E-04	1.3E+00
						Diethylstilbestrol	56-53-1		
						Difenzoquat	43222-48-6	9.4E-01	1.8E+05
						Diffubenzuron	35367-38-5		
						Diffluoroethane, 1,1-	75-37-6		
1.3E-05	C					Dihydrosafrole	94-58-6	9.4E-01	3.1E+03
						Diisopropyl Ether	108-20-3		
						Diisopropyl Methylphosphonate	1445-75-6		
						Dimethipin	55290-64-7	9.4E-03	
						Dimethoate	60-51-5		
						Dimethoxybenzidine, 3,3'-	119-90-4		
1.3E-03	C					Dimethyl methylphosphonate	756-79-6	9.4E-03	
						Dimethylamino azobenzene [p-]	60-11-7		
						Dimethylaniline HCl, 2,4-	21436-96-4		
						Dimethylaniline, 2,4-	95-68-1		
						Dimethylaniline, N,N-	121-69-7	7.7E-05	1.3E+02
						Dimethylbenzidine, 3,3'-	119-93-7		
						Dimethylformamide	68-12-2		
						Dimethylhydrazine, 1,1-	57-14-7		
1.6E-01	C					Dimethylhydrazine, 1,2-	540-73-6	7.7E-05	8.8E-03
						Dimethylphenol, 2,4-	105-67-9		
						Dimethylphenol, 2,6-	576-26-1		
						Dimethylphenol, 3,4-	95-65-8	9.4E-01	
1.3E-05	C					Dimethylvinylchloride	513-37-1		
						Dinitro-o-cresol, 4,6-	534-52-1		
						Dinitro-d-cyclohexyl Phenol, 4,6-	131-89-5	1.4E-01	
						Dinitrobenzene, 1,2-	528-29-0		
						Dinitrobenzene, 1,3-	99-65-0		
						Dinitrobenzene, 1,4-	100-25-4	1.4E-01	
						Dinitrophenol, 2,4-	51-28-5		
						Dinitrotoluene Mixture, 2,4/2,6-	NA		
8.9E-05	C					Dinitrotoluene, 2,4-	121-14-2	2.5E+00	1.3E+02
						Dinitrotoluene, 2,6-	606-20-2		
						Dinitrotoluene, 2-Amino-4,6-	35572-78-2		
						Dinitrotoluene, 4-Amino-2,6-	19406-51-0		
5.0E-06	I	3.0E-02	I	V		Dinitrotoluene, Technical grade	25321-14-6	3.2E-07	1.8E-04
						Dinoseb	88-85-7		
						Dioxane, 1,4-	123-91-1	2.5E+00	1.3E+02
1.3E+00	I					Dioxins		9.4E-06	1.8E-04
3.8E+01	C	4.0E-08	C	V		~Hexachlorodibenzo-p-dioxin, Mixture	NA		
						~TCDD, 2,3,7,8-	1746-01-6	3.2E-07	1.8E-04
						Diphenamid	957-51-7	5.6E-02	
						Diphenyl Sulfone	127-63-9		
						Diphenylamine	122-39-4		
2.2E-04	I					Diphenylhydrazine, 1,2-	122-66-7	8.8E-05	
						Diquat	85-00-7		
						Direct Black 38	1937-37-7		
1.4E-01	C					Direct Blue 6	2602-46-2	8.8E-05	
1.4E-01	C					Direct Brown 95	16071-86-6		
						Disulfoton	298-04-4	1.0E+01	4.4E+00
						Dithiane, 1,4-	505-29-3		
						Diuron	330-54-1		
						Dodine	2439-10-3		
						EPTC	759-94-4	1.0E+01	8.8E+01
						Endosulfan	115-29-7		
						Endothall	145-73-3		
1.2E-06	I	1.0E-03	I	V		Endrin	72-20-8	1.0E+01	4.4E+00
						Epichlorohydrin	106-89-8		
						Epoxybutane, 1,2-	106-88-7		

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Toxicity and Chemical-specific					Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y	v o l u t e	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
6.0E-02	P	V			Ethephon	16672-87-0		
					Ethion	563-12-2		
					Ethoxyethanol Acetate, 2-	111-15-9		2.6E+02
2.0E-01	I	V			Ethoxyethanol, 2-	110-80-5		8.8E+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
1.0E+01	I	V			Ethyl Chloride (Chloroethane)	75-00-3		4.4E+04
					Ethyl Ether	60-29-7		
3.0E-01	P	V			Ethyl Methacrylate	97-63-2		1.3E+03
2.5E-06	C	1.0E+00	I	V	Ethyl-p-nitrophenyl Phosphonate	2104-64-5		
					Ethylbenzene	100-41-4	4.9E+00	4.4E+03
					Ethylene Cyanohydrin	109-78-4		
					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
8.8E-05	C	3.0E-02	C	V	Ethylene Oxide	75-21-8	1.4E-01	1.3E+02
1.3E-05	C				Ethylene Thiourea	96-45-7	9.4E-01	
1.9E-02	C				Ethyleneimine	151-56-4	6.5E-04	
					Ethylphthalyl Ethyl Glycolate	84-72-0		
					Express	101200-48-0		
					Fenamiphos	22224-92-6		
					Fenpropathrin	39515-41-8		
1.3E-02	C				Fluometuron	2164-17-2		
					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		
					Flutolanil	66332-96-5		
					Fluvalinate	69409-94-5		
					Folpet	133-07-3		
1.3E-05	I	9.8E-03	A	V	Fomesafen	72178-02-0		
					Fopofos	944-22-9		
					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			
					~Dibenzofuran	132-64-9		
					~Furan	110-00-9		
2.0E+00	I	V			~Tetrahydrofuran	109-99-9		8.8E+03
					Furazolidone	67-45-8		
4.3E-04	C	5.0E-02	H	V	Furfural	98-01-1		2.2E+02
					Furium	531-82-8	2.9E-02	
8.6E-06	C				Furmecyclo	60568-05-0	1.4E+00	
					Glufosinate, Ammonium	77182-82-2		
8.0E-05	C				Glutaraldehyde	111-30-8		3.5E-01
1.0E-03	H	V			Glycidyl	765-34-4		4.4E+00
					Glyphosate	1071-83-6		
					Goal	42874-03-3		
					Guanidine	113-00-8		
1.0E-02	A				Guanidine Chloride	50-01-1		
					Guthion	86-50-0		4.4E+01
					Haloxypof, Methyl	69806-40-2		
1.3E-03	I				Harmony	79277-27-3	9.4E-03	
					Heptachlor	76-44-8		
2.6E-03	I				Heptachlor Epoxide	1024-57-3	4.7E-03	
					Hexabromobenzene	87-82-1		
					Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BDE-153)	68631-49-2		
4.6E-04	I				Hexachlorobenzene	118-74-1	2.7E-02	
2.2E-05	I				Hexachlorobutadiene	87-68-3	5.6E-01	
1.8E-03	I				Hexachlorocyclohexane, Alpha-	319-84-6	6.8E-03	
5.3E-04	I				Hexachlorocyclohexane, Beta-	319-85-7	2.3E-02	
3.1E-04	C				Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	4.0E-02	
5.1E-04	I				Hexachlorocyclohexane, Technical	608-73-1	2.4E-02	
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+00	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
					Hexamethylphosphoramide	680-31-9		
7.0E-01	I	V			Hexane, N-	110-54-3		3.1E+03
					Hexanedioic Acid	124-04-9		
3.0E-02	I	V			Hexanone, 2-	591-78-6		1.3E+02
					Hexazinone	51235-04-2		
4.9E-03	I	3.0E-05	P	V	Hydrazine	302-01-2	2.5E-03	1.3E-01
4.9E-03	I				Hydrazine Sulfate	10034-93-2	2.5E-03	

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Toxicity and Chemical-specific					Contaminant	Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y c	v o l a t i l e m u t a g e n	Analyte CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
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		
					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 NA		1.3E+03
					Kerb 23950-58-5		
					Lactofen 77501-63-4		
					Lead Compounds		
1.5E-01	C	2.0E-04	C	M	~Lead Chromate 7758-97-6	8.2E-05	8.8E-01
1.2E-05	C				~Lead Phosphate 7446-27-7	1.0E+00	
8.0E-05	C				~Lead acetate 301-04-2	1.5E-01	
1.2E-05	C				~Lead and Compounds 7439-92-1		
					~Lead subacetate 1335-32-6	1.0E+00	
	V				~Tetraethyl Lead 78-00-2		
					Linuron 330-55-2		
					Lithium 7439-93-2		
					Londax 83055-99-6		
					MCPA 94-74-6		
					MCPB 94-81-5		
					MCPP 93-65-2		
7.0E-04	C				Malathion 121-75-5		3.1E+00
					Maleic Anhydride 108-31-6		
					Maleic Hydrazide 123-33-1		
					Malonitrile 109-77-3		
					Mancozeb 8018-01-7		
5.0E-05	I				Maneb 12427-38-2		
5.0E-05	I				Manganese (Diet) 7439-96-5		2.2E-01
					Manganese (Non-diet) 7439-96-5		
					Mephistofar 950-10-7		
					Mepiquat Chloride 24307-26-4		
					Mercury Compounds		
3.0E-04	S				~Mercuric Chloride (and other Mercuric 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		
	V				~Phenylmercuric Acetate 62-38-4		
					Merphos 150-50-5		
					Merphos Oxide 78-48-8		
3.0E-02	P	V			Metaxyl 57837-19-1		1.3E+02
					Methacrylonitrile 120-90-7		
					Methamidophos 10265-92-6		
2.0E+01	I	V			Methanol 67-56-1		8.8E+04
					Methidathion 950-37-8		
					Methomyl 16752-77-5		
1.4E-05	C				Methoxy-5-nitroaniline, 2- 99-59-2	8.8E-01	
					Methoxychlor 72-43-5		
1.0E-03	P	V			Methoxyethanol Acetate, 2- 110-49-6		4.4E+00
2.0E-02	I	V			Methoxyethanol, 2- 109-86-4		8.8E+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-02	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		
4.0E-02	H	V			Methyl Phosphonic Acid 993-13-5		1.8E+02
2.8E-05	C				Methyl Styrene (Mixed Isomers) 25013-15-4		
					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+01	1.3E+04
					Methyl-1,4-benzenediamine dihydrochloride, 2- 615-45-2		
					Methyl-5-Nitroaniline, 2- 99-55-8		
2.4E-03	C				Methyl-N-nitro-N-nitrosoguanidine, N- 70-25-7	5.1E-03	
3.7E-05	C				Methylaniline Hydrochloride, 2- 636-21-5	3.3E-01	
					Methylarsonic acid 124-58-3		

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Toxicity and Chemical-specific					Contaminant	Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1		
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y	v o l a t i l e	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)	
6.3E-03	C				Methylbenzene, 1,4-diamine monohydrochloride, 2- Methylbenzene-1,4-diamine sulfate, 2- Methylcholanthrene, 3-	74612-12-7 615-50-9 56-49-5	1.9E-03		
1.0E-08	I	6.0E-01	I	V	Methylene Chloride	75-09-2	1.2E+03	2.6E+03	
4.3E-04	C				Methylene-bis(2-chloroaniline), 4,4'- Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-14-4 101-61-1	2.9E-02 9.4E-01		
4.6E-04	C	2.0E-02	C		Methylenbisbenzenamine, 4,4'- Methylenediphenyl Diisocyanate Methylstyrene, Alpha-	101-77-9 101-68-8 98-83-9	2.7E-02	8.8E+01 2.6E+00	
				V	Metolachlor Metribuzin Mineral oils	51218-45-2 21087-64-9 8012-95-1			
5.1E-03	C			V	Mirex Molinate Molybdenum	2385-85-5 2212-67-1 7439-98-7	2.4E-03		
				V	Monochloramine Monomethylaniline N,N'-Diphenyl-1,4-benzenediamine	10599-90-3 100-61-8 74-31-7			
0.0E+00	C	1.0E-01	P	V	Naled Naphtha, High Flash Aromatic (HFAN) Naphthylamine, 2-	300-76-5 64742-95-6 91-59-8		4.4E+02	
2.6E-04	C	1.4E-05	C		Napropamide	15299-99-7			
2.6E-04	C	1.4E-05	C		Nickel Acetate Nickel Carbonate	373-02-4 3333-67-3	4.7E-02 4.7E-02	6.1E-02 6.1E-02	
2.6E-04	C	1.4E-05	C	V	Nickel Carbonyl	13463-39-3	4.7E-02	6.1E-02	
2.6E-04	C	1.4E-05	C		Nickel Hydroxide	12054-48-7	4.7E-02	6.1E-02	
2.6E-04	C	2.0E-05	C		Nickel Oxide	1313-99-1	4.7E-02	8.8E-02	
2.4E-04	I	1.4E-05	C		Nickel Refinery Dust	NA	5.1E-02	6.1E-02	
2.6E-04	C	9.0E-05	A		Nickel Soluble Salts	7440-02-0	4.7E-02	3.9E-01	
4.8E-04	I	1.4E-05	C		Nickel Sulfide	12035-72-2	2.6E-02	6.1E-02	
2.6E-04	C	1.4E-05	C		Nickelocene Nitrate Nitrate + Nitrite (as N)	1271-28-9 14797-55-8 NA	4.7E-02	6.1E-02	
		5.0E-05	X		Nitrite Nitroaniline, 2- Nitroaniline, 4-	14797-85-0 89-74-4 100-01-6		2.2E-01 2.6E+01	
4.0E-05	I	9.0E-03	I	V	Nitrobenzene Nitrocellulose Nitrofurantoin	98-95-3 9004-70-0 67-20-9	3.1E-01	3.9E+01	
3.7E-04	C				Nitrofurazone Nitroglycerin Nitroquajidine	59-87-0 55-63-0 556-88-7	3.3E-02		
8.8E-06	P	5.0E-03	P	V	Nitromethane	75-52-5	1.4E+00	2.2E+01	
2.7E-03	H	2.0E-02	I	V	Nitropropane, 2-	79-46-9	4.5E-03	8.8E+01	
7.7E-03	C			M	Nitroso-N-ethylurea, N-	759-73-9	1.6E-03		
3.4E-02	C			M	Nitroso-N-methylurea, N-	684-93-5	3.6E-04		
1.6E-03	I			V	Nitroso-di-N-butylamine, N-	924-16-3	7.7E-03		
2.0E-03	C				Nitroso-di-N-propylamine, N-	621-64-7	6.1E-03		
8.0E-04	C				Nitrosodihethanolamine, N-	1116-54-7	1.5E-02		
4.3E-02	I			M	Nitrosodithylamine, N-	55-18-5	2.9E-04		
1.4E-02	I	4.0E-05	X	V	M	Nitrosodimethylamine, N-	62-75-9	8.8E-04	1.8E-01
2.6E-06	C				Nitrosodiphenylamine, N-	86-30-6	4.7E+00		
6.3E-03	C			V	Nitrosomethylethylamine, N-	10595-95-6	1.9E-03		
1.9E-03	C				Nitrosomorpholine [N-]	59-89-2	6.5E-03		
2.7E-03	C				Nitrosopiperidine [N-]	100-75-4	4.5E-03		
6.1E-04	I				Nitrosopyrrolidine, N- Nitrotoluene, m-	930-55-2 99-08-1	2.0E-02		
				V	Nitrotoluene, o- Nitrotoluene, p- Nonane, n-	88-72-2 99-99-0 111-84-2		8.8E+01	
					Norflurazon Nustar Octabromodiphenyl Ether	27314-13-2 85509-19-9 32536-52-0			
					Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX) Octamethylpyrophosphoramide Oryzalin	2691-41-0 152-16-9 19044-88-3			
					Oxadiazon Oxamyl Paclobutrazol	19666-30-9 23135-22-0 76738-62-0			
				V	Paraquat Dichloride Parathion Pebulate	1910-42-5 56-38-2 1114-71-2			
					Pendimethalin Pentabromodiphenyl Ether Pentabromodiphenyl ether, 2,2',4,4',5'- (BDE-99)	40487-42-1 32534-81-9 60348-60-9			

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Toxicity and Chemical-specific					Contaminant	Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1	
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y c	v o l a t i l e m u t a g e n	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
				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+00	
		1.0E+00	P	V	Pentaerythritol tetranitrate (PETN)	78-11-5		
					Pentane, n-	109-66-0		4.4E+03
					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		
				V	Perfluorobutane Sulfonate	375-73-5		
6.3E-07	C				Permethrin	52645-53-1	1.9E+01	
					Phenacetin	62-44-2		
		2.0E-01	C		Phenmedipham	13684-63-4		
					Phenol	108-95-2		8.8E+02
					Phenothiazine	92-84-2		
					Phenylenediamine, m-	108-45-2		
					Phenylenediamine, o-	95-54-5		
					Phenylenediamine, p-	106-50-3		
		3.0E-04	I	V	Phenylphenol, 2-	90-43-7		
					Phorate	298-02-2		
					Phosgene	75-44-5		1.3E+00
					Phosmet	732-11-6		
					Phosphates, Inorganic			
					~Aluminum metaphosphate	13776-88-0		
					~Ammonium polyphosphate	68333-79-9		
					~Calcium pyrophosphate	7790-76-3		
					~Diammonium phosphate	7783-28-0		
					~Dicalcium phosphate	7757-93-9		
					~Dimagnesium phosphate	7782-75-4		
					~Dipotassium phosphate	7758-11-4		
					~Disodium phosphate	7558-79-4		
					~Monocalcium phosphate	13530-50-2		
					~Monodiammonium phosphate	7722-76-1		
					~Monocalcium phosphate	7758-23-8		
					~Monomagnesium phosphate	7757-86-0		
					~Monopotassium phosphate	7778-77-0		
					~Monosodium phosphate	7558-80-7		
					~Polyphosphoric acid	8017-16-1		
					~Potassium tripolyphosphate	13845-36-8		
					~Sodium acid pyrophosphate	7758-16-9		
					~Sodium aluminum phosphate (acidic)	7785-88-8		
					~Sodium aluminum phosphate (anhydrous)	10279-59-1		
					~Sodium aluminum phosphate (tetrahydrate)	10305-76-7		
					~Sodium hexametaphosphate	10124-56-8		
					~Sodium polyphosphate	68915-31-1		
					~Sodium trimetaphosphate	7785-84-4		
					~Sodium tripolyphosphate	7758-29-4		
					~Tetrapotassium phosphate	7320-34-5		
					~Tetrasodium pyrophosphate	7722-88-5		
					~Trialuminum sodium tetra decahydrogenoctaorthophosphate (dihydrate)	15136-87-5		
					~Tricalcium phosphate	7758-87-4		
					~Trimagnesium phosphate	7757-87-1		
					~Tripotassium phosphate	7778-53-2		
					~Trisodium phosphate	7601-54-9		
		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		
		2.4E-06	C		Phthalates		5.1E+00	
					~Bis(2-ethylhexyl)phthalate	117-81-7		
					~Butylphthalyl Butylglycolate	85-70-1		
					~Dibutyl Phthalate	84-74-2		
				V	~Diethyl Phthalate	84-66-2		
					~Dimethylterephthalate	120-61-6		
					~Octyl Phthalate, di-N-	117-84-0		
		2.0E-02	C		~Phthalic Acid, P-	100-21-0		
					~Phthalic Anhydride	85-44-9		8.8E+01
					Picloram	1918-02-1		
					Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3		
					Pirimiphos, Methyl	29232-93-7		
8.6E-03	C				Polybrominated Biphenyls	59536-65-1	1.4E-03	
2.0E-05	S			V	Polychlorinated Biphenyls (PCBs)		6.1E-01	
					~Aroclor 1016	12674-11-2		

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Toxicity and Chemical-specific					Contaminant	Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1	
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k o y c	v o l a t i l e	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
5.7E-04	S			V	~Aroclor 1221	11104-28-2	2.1E-02	
5.7E-04	S			V	~Aroclor 1232	11141-16-5	2.1E-02	
5.7E-04	S			V	~Aroclor 1242	53469-21-9	2.1E-02	
5.7E-04	S			V	~Aroclor 1248	12672-29-6	2.1E-02	
5.7E-04	S			V	~Aroclor 1254	11097-69-1	2.1E-02	
5.7E-04	S			V	~Aroclor 1260	11096-82-5	2.1E-02	
				V	~Aroclor 5460	11126-42-4		
1.1E-03	E	1.3E-03	E	V	~Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	1.1E-02	5.8E+00
1.1E-03	E	1.3E-03	E	V	~Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167)	52663-72-6	1.1E-02	5.8E+00
1.1E-03	E	1.3E-03	E	V	~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	1.1E-02	5.8E+00
1.1E-03	E	1.3E-03	E	V	~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 156)	38390-08-4	1.1E-02	5.8E+00
1.1E+00	E	1.3E-06	E	V	~Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	1.1E-05	5.8E-03
1.1E-03	E	1.3E-03	E	V	~Pentachlorobiphenyl, 2',3,4,4',5'- (PCB 123)	65510-44-3	1.1E-02	5.8E+00
1.1E-03	E	1.3E-03	E	V	~Pentachlorobiphenyl, 2,3',4,4',5'- (PCB 118)	31508-00-6	1.1E-02	5.8E+00
1.1E-03	E	1.3E-03	E	V	~Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	1.1E-02	5.8E+00
1.1E-03	E	1.3E-03	E	V	~Pentachlorobiphenyl, 2,3,4,4',5'- (PCB 114)	74472-37-0	1.1E-02	5.8E+00
3.8E+00	E	4.0E-07	E	V	~Pentachlorobiphenyl, 3,3',4,4',5'- (PCB 126)	57465-28-8	3.2E-06	1.8E-03
5.7E-04	I			V	~Polychlorinated Biphenyls (high risk)	1336-36-3	2.1E-02	
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	E	4.0E-04	E	V	~Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	3.2E-03	1.8E+00
1.1E-02	E	1.3E-04	E	V	~Tetrachlorobiphenyl, 3,4,4',5'- (PCB 81)	70362-50-4	1.1E-03	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		
1.1E-04	C			V	~Benz[a]anthracene	56-55-3	1.1E-01	
1.1E-04	C				~Benzo[j]fluoranthene	205-82-3	1.1E-01	
1.1E-03	C			M	~Benzo[a]pyrene	50-32-8	1.1E-02	
1.1E-04	C			M	~Benzo[b]fluoranthene	205-99-2	1.1E-01	
1.1E-04	C			M	~Benzo[k]fluoranthene	207-08-9	1.1E-01	
				V	~Chloronaphthalene, Beta-	91-58-7		
1.1E-05	C			M	~Chrysene	218-01-9	1.1E+00	
1.2E-03	C			M	~Dibenz[a,h]anthracene	53-70-3	1.0E-02	
1.1E-03	C			M	~Dibenz[a,e]pyrene	192-65-4	1.1E-02	
7.1E-02	C			M	~Dimethylbenz[a]anthracene, 7,12-	57-97-6	1.7E-04	
				V	~Fluoranthene	208-44-0		
1.1E-04	C			M	~Fluorene	86-73-7	1.1E-01	
				V	~Indeno[1,2,3-cd]pyrene	193-39-5		
				V	~Methylnaphthalene, 1-	90-12-0		
3.4E-05	C	3.0E-03	I	V	~Methylnaphthalene, 2-	91-57-6	3.6E-01	1.3E+01
				V	~Naphthalene	91-20-3		
1.1E-04	C			V	~Nitrofluorene, 4-	57835-92-4	1.1E-01	
				V	~Pyrene	129-00-0		
				V	Potassium Perfluorobutane Sulfonate	29420-49-3		
				V	Prochloraz	67747-09-5		
				V	Profluralin	26399-36-0		
				V	Prometon	1610-18-0		
				V	Prometryn	7287-19-6		
				V	Propachlor	1918-16-7		
				V	Propanil	709-98-8		
				V	Propargite	2312-35-8		
				V	Propargyl Alcohol	107-19-7		
				V	Propazine	139-40-2		
				V	Propam	122-42-9		
8.0E-03	I	V		V	Propiconazole	60207-90-1		
				V	Propionaldehyde	123-38-6		3.5E+01
1.0E+00	X	V		V	Propyl benzene	103-65-1		4.4E+03
3.0E+00	C	V		V	Propylene	115-07-1		1.3E+04
				V	Propylene Glycol	57-55-6		
2.7E-04	A			V	Propylene Glycol Dinitrate	6423-43-4		1.2E+00
				V	Propylene Glycol Monoethyl Ether	1569-02-4		
2.0E+00	I	V		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+00	1.3E+02
				V	Pursuit	81335-77-5		
				V	Pydrin	51630-58-1		
				V	Pyridine	110-86-1		
				V	Quinalphos	13593-03-8		
				V	Quinoline	91-22-5		
3.0E-02	A			V	Refractory Ceramic Fibers	NA		1.3E+02
				V	Resmethrin	10453-86-8		
				V	Ronnell	299-84-3		
6.3E-05	C			M	Rotenone	83-79-4	1.9E-01	
				V	Safrole	94-59-7		
				V	Savey	78587-05-0		

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Toxicity and Chemical-specific					Contaminant	Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1	
IUR (ug/m ³) ⁻¹	k _e y	RfC _i (mg/m ³)	k _e o c	v o l u t e m u t a g e n	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
2.0E-02	C				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
3.0E-03	C				Sethoxydim	74051-80-2		
					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		
1.5E-01	C	2.0E-04	C	M	Sodium Dichromate	10588-01-9	8.2E-05	8.8E-01
					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		
					Stirofos (Tetrachlorovinphos)	961-11-5		
1.5E-01	C	2.0E-04	C	M	Strontium Chromate	7789-06-2	8.2E-05	8.8E-01
					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	NA		
2.0E-03	X				Sulfolane	126-33-0		8.8E+00
1.0E-03	C			V	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9		
1.0E-03	C				Sulfur Trioxide	7446-11-9		4.4E+00
					Sulfuric Acid	7664-93-9		4.4E+00
					Systhane	88671-89-0		
					TCMTB	21564-17-0		
					Tebuthiuron	34014-18-1		
				V	Temephos	3383-96-8		
					Terbacil	5902-51-2		
					Terbufos	13071-79-9		
				V	Terbutryn	886-50-0		
					Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1		
					Tetrachlorobenzene, 1,2,4,5-	95-94-3		
7.4E-06	I			V	Tetrachloroethane, 1,1,1,2-	630-20-6	1.7E+00	
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+01	1.8E+02
				V	Tetrachlorophenol, 2,3,4,6-	58-90-2		
					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
					Tetryl (Trinitrophenylmethyl)nitramine	479-45-8		
					Thallium (I) Nitrate	10102-45-1		
				V	Thallium (Soluble Salts)	7440-28-0		
					Thallium Acetate	563-68-8		
					Thallium Carbonate	6533-73-9		
					Thallium Chloride	7791-12-0		
					Thallium Sulfate	7446-18-6		
					Thiobencarb	28249-77-6		
					Thiodiglycol	111-48-8		
					Thioflanzox	39196-18-4		
					Thiophanate, Methyl	23564-05-8		
					Thiram	137-26-8		
1.0E-04	A			V	Tin	7440-31-5		4.4E-01
5.0E+00	I			V	Titanium Tetrachloride	7550-45-0		
					Toluene	108-88-3		2.2E+04
					Toluene-2,5-diamine	95-70-5		
					Toluidine, p-	106-49-0		
				V	Total Petroleum Hydrocarbons (Aliphatic High)	NA		
6.0E-01	P			V	Total Petroleum Hydrocarbons (Aliphatic Low)	NA		2.6E+03
1.0E-01	P			V	Total Petroleum Hydrocarbons (Aliphatic Medium)	NA		4.4E+02
					Total Petroleum Hydrocarbons (Aromatic High)	NA		
3.0E-02	P			V	Total Petroleum Hydrocarbons (Aromatic Low)	NA		1.3E+02
3.0E-03	P			V	Total Petroleum Hydrocarbons (Aromatic Medium)	NA		1.3E+01
3.2E-04	I				Toxaphene	8001-35-2	3.8E-02	
				V	Tralometrin	66841-25-6		
					Tri-n-butyltin	688-73-3		
				V	Triacetin	102-76-1		
					Triallate	2303-17-5		
					Triasulfuron	82097-50-5		
				V	Tribromobenzene, 1,2,4-	615-54-3		
					Tributyl Phosphate	126-73-8		
					Tributyltin Compounds	NA		
3.0E+01	H			V	Tributyltin Oxide	56-35-9		
					Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1		1.3E+05
					Trichloroacetic Acid	76-03-9		

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Toxicity and Chemical-specific					Contaminant		Carcinogenic Target Risk (TR) = 1E-06	Noncancer Hazard Index (HI) = 1
IUR (ug/m ³) ⁻¹	k e y	RfC _i (mg/m ³)	k e y	v o l u t i l e	Analyte	CAS No.	Carcinogenic SL TR=1.0E-6 (ug/m ³)	Noncarcinogenic SL HI=1 (ug/m ³)
					Trichloroaniline HCl, 2,4,6- Trichloroaniline, 2,4,6- Trichlorobenzene, 1,2,3-	33663-50-2 634-93-5 87-61-6		
	V							
2.0E-03 5.0E+00 1.6E-05	P I I				Trichlorobenzene, 1,2,4- Trichloroethane, 1,1,1- Trichloroethane, 1,1,2-	120-82-1 71-55-6 79-00-5		8.8E+00 2.2E+04 8.8E-01
4.1E-06	I	2.0E-03 7.0E-01			Trichloroethylene Trichlorofluoromethane Trichlorophenol, 2,4,5-	79-01-6 75-69-4 95-95-4	3.0E+00	8.8E+00 3.1E+03
3.1E-06	I				Trichlorophenol, 2,4,6- Trichlorophenoxyacetic Acid, 2,4,5- Trichlorophenoxypropionic acid, -2,4,5	88-06-2 93-76-5 93-72-1	4.0E+00	
	V							
3.0E-04 3.0E-04	I P				Trichloropropane, 1,1,2- Trichloropropane, 1,2,3- Trichloropropene, 1,2,3-	598-77-6 96-18-4 96-19-5		1.3E+00 1.3E+00
7.0E-03	I				Tricresyl Phosphate (TCP) Tridiphane Triethylamine	1330-78-5 58138-08-2 121-44-8		3.1E+01
	V							
5.0E-03 7.0E-03	P P				Triethylene Glycol Trifluralin Trimethyl Phosphate Trimethylbenzene, 1,2,3- Trimethylbenzene, 1,2,4- Trimethylbenzene, 1,3,5-	142-27-6 1582-09-8 512-56-1 526-73-8 95-83-6 108-67-8		2.2E+01 3.1E+01
	V							
6.6E-04	C				Trinitrobenzene, 1,3,5- Trinitrotoluene, 2,4,6- Triphenylphosphine Oxide Tris(1,3-Dichloro-2-propyl) Phosphate Tris(1-chloro-2-propyl)phosphate Tris(2,3-dibromopropyl)phosphate	99-35-4 118-96-7 791-28-6 13674-87-8 13674-84-5 126-72-7	1.9E-02	
4.0E-05	A				Tris(2-chloroethyl)phosphate Tris(2-ethylhexyl)phosphate Uranium (Soluble Salts)	115-96-8 78-42-2 NA		1.8E-01
2.9E-04 8.3E-03 1.0E-04	C P A				Urethane Vanadium Pentoxide Vanadium and Compounds	51-79-6 1314-62-1 7440-62-2	4.2E-02 1.5E-03	3.1E-02 4.4E-01
	V							
2.0E-01	I				Vinclozolin Vinyl Acetate	1929-77-7 50471-44-8 108-05-4		8.8E+02
3.2E-05 4.4E-06	H I	3.0E-03 1.0E-01			Vinyl Bromide Vinyl Chloride Warfarin	593-60-2 75-01-4 81-81-2	3.8E-01 2.8E+00	1.3E+01 4.4E+02
1.0E-01 1.0E-01 1.0E-01	S S S				Xylene, p- Xylene, m- Xylene, o-	106-42-3 108-38-3 95-47-6		4.4E+02 4.4E+02 4.4E+02
1.0E-01	I				Xylenes Zinc Phosphide Zinc and Compounds	1330-20-7 1314-84-7 7440-66-6		4.4E+02
	V							
					Zineb Zirconium	12122-67-7 7440-67-7		

Contaminant	Molecular Weight	Volatility Parameters						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	Vapor Pressure (g/m ³)	VP Ref	Density (g/cm ³)	Density Ref	D ₁₀ (cm ² /s)	D _w (cm ² /s)	Dia Ref	K _{ow}	K _{ow} Ref	K _{oc} (L/kg)	K _{oc} Ref	log K _{ow}	log K _{oc} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/yr)	t (hr)	K _p (cm ² /hr)	K _p Ref
ALAR	1596-84-5	1.6E+02	EPI	1.7E-08	4.2E-10	EPI	2.0E-04	EPI	1.4E+00	CRC89	6.4E-02	7.5E-06	WATER9 (U.S. EPA, 2001)	1.0E+01	EPI	-1.5E+00	EPI	1.0E+05	EPI	1.0E+05	EPI	1.0E+05	EPI	9.7E-05	8.3E-01	2.0E+00	2.0E-05	EPI
Acephate	30560-19-1	1.8E+02	EPI	2.0E-11	5.0E-13	EPI	1.7E-06	EPI	1.4E+00	CRC89	3.7E-02	8.0E-06	WATER9 (U.S. EPA, 2001)	1.0E+01	EPI	-8.5E-01	EPI	8.2E+05	EPI	8.2E+05	EPI	8.2E+05	EPI	2.1E-04	1.1E+00	2.7E+00	4.0E-05	EPI
Acetaldelyde	75-07-0	4.4E+01	EPI	2.7E-03	6.7E-05	EPI	9.0E-02	EPI	7.8E-01	CRC89	1.3E-01	1.4E-05	WATER9 (U.S. EPA, 2001)	3.0E+00	EPI	-3.4E-01	EPI	1.0E+06	EPI	1.0E+06	EPI	1.0E+06	EPI	1.3E-03	1.9E-01	4.5E-01	5.3E-04	EPI
Acetochlor	34268-82-1	2.7E+02	EPI	9.1E-07	2.2E-08	EPI	2.8E-05	EPI	1.1E+00	PubChem	2.2E-02	5.6E-06	WATER9 (U.S. EPA, 2001)	3.0E+00	EPI	3.0E+00	EPI	2.2E+02	EPI	2.2E+02	EPI	2.2E+02	EPI	3.1E-02	3.4E+00	8.2E+00	5.0E-03	EPI
Acetol	67-64-1	5.8E+01	EPI	1.4E-03	3.5E-05	EPI	2.3E+02	EPI	7.8E-01	CRC89	1.4E-03	5.0E-06	WATER9 (U.S. EPA, 2001)	2.4E+00	EPI	2.4E+00	EPI	1.0E+06	EPI	1.0E+06	EPI	1.0E+06	EPI	1.5E-03	2.2E-01	5.3E-01	1.5E-04	EPI
Acetone Cyanohydril	75-86-5	8.5E+01	EPI	5.3E-04	1.3E-05	EPI	3.4E-01	EPI	9.3E-01	CRC89	8.6E-02	1.0E-05	WATER9 (U.S. EPA, 2001)	1.0E+00	EPI	-3.0E-02	EPI	1.0E+06	EPI	1.0E+06	EPI	1.0E+06	EPI	1.8E-03	3.2E-01	7.6E-01	5.0E-04	EPI
Acetonitrile	125-06-8	4.1E+01	EPI	1.4E-03	3.5E-05	EPI	8.9E+01	EPI	7.9E-01	CRC89	1.3E-01	1.4E-05	WATER9 (U.S. EPA, 2001)	4.7E+00	EPI	-3.4E-01	EPI	1.0E+06	EPI	1.0E+06	EPI	1.0E+06	EPI	1.4E-03	1.8E-01	4.3E-01	5.5E-04	EPI
Acetophenone	98-86-2	1.2E+02	EPI	4.3E-04	1.0E-05	EPI	4.0E-01	EPI	1.0E+00	CRC89	6.5E-02	8.7E-06	WATER9 (U.S. EPA, 2001)	5.2E+01	EPI	1.6E+00	EPI	6.1E+03	EPI	6.1E+03	EPI	6.1E+03	EPI	1.6E-02	5.0E-01	1.2E+00	3.7E-03	EPI
Aceylaminofluorene, 2	53-96-3	2.2E+02	EPI	7.8E-09	1.9E-10	EPI	2.9E-05	EPI	2.9E-05	EPI	2.2E-02	6.0E-06	WATER9 (U.S. EPA, 2001)	2.2E+03	EPI	3.1E+00	EPI	2.2E+02	EPI	2.2E+02	EPI	2.2E+02	EPI	7.2E-02	1.9E+00	4.5E+00	1.2E-02	RAGSE
Acrolein	107-02-8	5.6E+01	EPI	5.0E-03	1.2E-04	EPI	2.7E-02	EPI	8.4E-01	CRC89	1.1E-01	1.2E-05	WATER9 (U.S. EPA, 2001)	1.0E+00	EPI	-1.0E-02	EPI	2.1E+05	EPI	2.1E+05	EPI	2.1E+05	EPI	2.2E-03	2.2E-01	5.2E-01	7.5E-04	EPI
Acrylamide	79-06-1	7.1E+01	EPI	7.0E-08	1.7E-09	EPI	7.0E-03	EPI	1.2E+00	LANGE	1.1E-01	1.3E-05	WATER9 (U.S. EPA, 2001)	5.7E+00	EPI	-6.7E-01	EPI	3.9E+05	EPI	3.9E+05	EPI	3.9E+05	EPI	7.3E-04	2.6E-01	6.3E-01	2.2E-04	EPI
Acrylic Acid	79-10-7	7.2E+01	EPI	1.5E-05	3.7E-07	EPI	4.0E+00	EPI	1.1E+00	CRC89	1.0E-01	1.2E-05	WATER9 (U.S. EPA, 2001)	1.4E+00	EPI	3.5E-01	EPI	1.0E+06	EPI	1.0E+06	EPI	1.0E+06	EPI	3.4E-03	2.7E-01	6.4E-01	1.1E-03	EPI
Acrylonitrile	107-13-1	5.6E+01	EPI	5.6E-03	1.4E-04	EPI	1.1E+02	EPI	8.0E-01	CRC89	1.1E-01	1.2E-05	WATER9 (U.S. EPA, 2001)	8.5E+00	EPI	2.5E-01	EPI	7.5E+04	EPI	7.5E+04	EPI	7.5E+04	EPI	3.2E-03	2.1E-01	5.0E-01	1.2E-03	EPI
Adiponitrile	111-69-3	1.1E+02	EPI	4.9E-08	1.2E-09	EPI	6.8E-04	EPI	9.7E-01	CRC89	7.1E-02	9.0E-06	WATER9 (U.S. EPA, 2001)	2.0E+01	EPI	-3.2E-01	EPI	8.0E+04	EPI	8.0E+04	EPI	8.0E+04	EPI	9.5E-04	4.2E-01	1.0E+00	2.4E-04	EPI
Alachlor	15972-60-8	2.7E+02	EPI	3.4E-07	8.3E-09	EPI	2.1E-05	EPI	1.1E+00	CRC89	2.3E-02	5.7E-06	WATER9 (U.S. EPA, 2001)	3.1E+02	EPI	3.5E+00	EPI	2.4E+02	EPI	2.4E+02	EPI	2.4E+02	EPI	6.6E-02	3.4E+00	8.2E+00	1.1E-02	EPI
Aldicarb	116-06-3	1.9E+02	EPI	5.9E-08	1.4E-09	EPI	3.5E-05	EPI	1.2E+00	CRC89	3.2E-02	7.2E-06	WATER9 (U.S. EPA, 2001)	2.5E+01	EPI	1.1E+00	EPI	6.0E+03	EPI	6.0E+03	EPI	6.0E+03	EPI	4.0E-03	1.2E+00	2.9E+00	7.6E-04	EPI
Aldicarb Sulfone	1646-88-4	2.2E+02	EPI	1.4E-07	3.4E-09	EPI	9.0E-05	EPI	1.2E+00	CRC89	5.2E-02	6.1E-06	WATER9 (U.S. EPA, 2001)	1.0E+01	EPI	-5.7E-01	EPI	1.0E+04	EPI	1.0E+04	EPI	1.0E+04	EPI	2.1E-04	1.8E+00	4.4E+00	3.7E-05	EPI
Aldicarb sulfoxide	1646-87-3	2.1E+02	EPI	4.0E-08	9.7E-10	EPI	1.0E-04	EPI	1.2E+00	CRC89	5.4E-02	6.4E-06	WATER9 (U.S. EPA, 2001)	1.0E+01	EPI	-7.8E-01	EPI	2.8E+04	EPI	2.8E+04	EPI	2.8E+04	EPI	1.8E-04	1.5E+00	3.9E+00	3.3E-05	EPI
Aldrin	309-90-2	3.3E+02	EPI	9.8E-03	4.4E-05	EPI	1.2E-04	EPI	1.6E+00	PubChem	3.8E-02	5.8E-06	WATER9 (U.S. EPA, 2001)	8.2E+04	EPI	6.5E+00	EPI	1.7E+02	EPI	1.7E+02	EPI	1.7E+02	EPI	2.2E-05	1.3E+01	4.8E+01	2.3E-01	EPI
Allyl	74223-64-6	5.8E+01	EPI	5.4E-15	1.3E-16	EPI	2.9E-11	EPI	2.9E-11	EPI	3.9E-01	4.2E-06	WATER9 (U.S. EPA, 2001)	9.3E+01	EPI	9.3E+01	EPI	9.5E+03	EPI	9.5E+03	EPI	9.5E+03	EPI	2.5E-03	1.4E+00	3.9E-01	3.3E-04	EPI
Allyl Alcohol	107-18-6	5.8E+01	EPI	2.0E-04	5.0E-06	EPI	2.6E+01	EPI	8.5E-01	CRC89	1.1E-01	1.2E-05	WATER9 (U.S. EPA, 2001)	1.9E+00	EPI	1.7E-01	EPI	1.0E+06	EPI	1.0E+06	EPI	1.0E+06	EPI	2.8E-03	2.2E-01	5.3E-01	9.6E-04	EPI
Allyl Chloride	107-05-1	7.7E+01	EPI	4.5E-01	1.1E-02	EPI	3.7E+02	EPI	9.4E-01	CRC89	9.4E-02	1.1E-05	WATER9 (U.S. EPA, 2001)	1.5E+03	BAES	4.0E+01	EPI	1.9E+00	YAWS	3.4E+03	EPI	3.4E+03	EPI	3.8E-02	2.8E-01	6.8E-01	1.1E-02	EPI
Aluminum	7429-90-5	2.7E+01	CRC89	0.0E+00	0.0E+00	NIOSH	2.7E+00	CRC89	2.7E+00	CRC89	2.4E+00	2.7E+00	WATER9 (U.S. EPA, 2001)	1.5E+03	BAES	4.0E+01	EPI	1.9E+00	YAWS	3.4E+03	EPI	3.4E+03	EPI	2.0E-03	1.5E-01	3.6E-01	1.0E-03	RAGSE
Aluminum Phosphide	20859-73-8	5.8E+01	EPI	0.0E+00	0.0E+00	NIOSH	2.4E+00	CRC89	2.4E+00	CRC89	2.4E+00	2.4E+00	WATER9 (U.S. EPA, 2001)	1.5E+03	BAES	4.0E+01	EPI	1.9E+00	YAWS	3.4E+03	EPI	3.4E+03	EPI	2.9E-03	2.2E-01	5.3E-01	1.0E-03	RAGSE
Amdro	67485-29-4	4.9E+02	EPI	9.0E-05	2.2E-06	EPI	2.0E-08	EPI	2.0E-08	EPI	3.0E-02	3.6E-06	WATER9 (U.S. EPA, 2001)	1.8E+08	EPI	2.3E+00	EPI	6.0E+03	EPI	6.0E+03	EPI	6.0E+03	EPI	7.7E-04	2.0E+01	1.5E+02	9.0E-05	EPI
Ametryn	834-12-8	2.3E+02	EPI	9.9E-08	2.4E-09	EPI	2.7E-06	EPI	4.3E+02	EPI	5.1E-02	6.0E-06	WATER9 (U.S. EPA, 2001)	3.0E+00	EPI	3.0E+00	EPI	2.1E+02	EPI	2.1E+02	EPI	2.1E+02	EPI	4.6E-02	6.2E+00	4.7E+00	7.9E-03	EPI
Aminobiphenyl, 4	92-67-1	1.7E+02	EPI	7.1E-06	1.7E-07	EPI	9.6E-04	EPI	9.6E-04	EPI	6.2E-02	7.3E-06	WATER9 (U.S. EPA, 2001)	2.5E+03	EPI	2.9E+00	EPI	1.3E+02	EPI	1.3E+02	EPI	1.3E+02	EPI	7.0E-02	9.3E-01	2.2E+00	1.4E-02	EPI
Aminophenol, m-	591-27-5	1.1E+02	EPI	1.1E-08	2.7E-10	EPI	1.9E-03	EPI	8.0E-01	CRC89	8.3E-02	9.7E-06	WATER9 (U.S. EPA, 2001)	2.1E-01	EPI	2.1E-01	EPI	2.7E+04	EPI	2.7E+04	EPI	2.7E+04	EPI	2.1E-03	4.3E-01	1.0E+00	5.3E-04	EPI
Aminophenol, p-	123-30-8	1.1E+02	EPI	1.5E-08	3.6E-10	EPI	4.0E-05	EPI	8.0E-01	CRC89	8.3E-02	9.7E-06	WATER9 (U.S. EPA, 2001)	9.0E+01	EPI	4.0E-02	EPI	1.6E+04	EPI	1.6E+04	EPI	1.6E+04	EPI	1.6E-03	4.3E-01	1.0E+00	4.1E-04	EPI
Amirbaz	33089-61-1	2.9E+02	EPI	4.0E-04	9.9E-06	EPI	2.0E-06	EPI	1.1E+00	CRC89	2.2E-02	5.4E-06	WATER9 (U.S. EPA, 2001)	9.0E+05	EPI	5.5E+00	EPI	1.0E+00	EPI	1.0E+00	EPI	1.0E+00	EPI	1.1E+00	4.6E+00	1.8E+01	1.6E-01	EPI
Ammonia	7664-41-7	1.7E+01	CRC89	6.6E-04	1.6E-05	PHYSPROP	7.5E+03	PHYSPROP	7.0E-01	CRC89	2.3E-01	2.2E-05	WATER9 (U.S. EPA, 2001)	2.3E-01	OTHER	9.0E+05	PERRY	9.0E+05	PERRY	9.0E+05	PERRY	9.0E+05	PERRY	1.6E-03	1.3E-01	3.1E-01	1.0E-03	RAGSE
Ammonium Sulfamate	7773-06-0	1.1E+02	CRC89	0.0E+00	0.0E+00	NIOSH	1.8E+00	PubChem	1.8E+00	PubChem	7.9E-02	9.1E-06	WATER9 (U.S. EPA, 2001)	4.1E+00	EPI	8.9E-01	EPI	1.1E+05	EPI	1.1E+05	EPI	1.1E+05	EPI	4.1E-03	4.6E-01	1.1E+00	1.0E-03	RAGSE
Amly Alcohol, tert	75-85-4	8.9E+01	EPI	5.6E-04	1.4E-05	EPI	1.7E+01	EPI	8.1E-01	CRC89	3.8E-02	9.1E-06	WATER9 (U.S. EPA, 2001)	7.0E+01	EPI	9.0E-01	EPI	3.6E+04	EPI	3.6E+04	EPI	3.6E+04	EPI	7.9E-03	3.3E-01	7.9E-01	2.0E-03	EPI
Aniline	62-53-3	9.3E+01	EPI	8.3E-05	2.0E-06	EPI	6.7E-01	EPI	1.0E+00	CRC89	8.3E-02	1.0E-05	WATER9 (U.S. EPA, 2001)	7.0E+01	EPI	9.0E-01	EPI	3.6E+04	EPI	3.6E+04	EPI	3.6E+04	EPI	6.9E-03	3.3E-01	8.4E-01	1.3E-03	EPI
Anthraquinone, 9,10-	84-65-1	2.1E+02	EPI	9.6E-07	2.4E-08	EPI	1.2E-07	EPI	1.2E-07	EPI	5.4E-02	6.3E-06	WATER9 (U.S. EPA, 2001)	5.0E+03	EPI	3.4E+00	EPI	1.4E+00	EPI	1.4E+00	EPI	1.4E+00	EPI	1.1E-01	1.5E+00	3.7E+00	1.9E-02	EPI
Antimony (metallic)	7440-36-0	1.2E+02	CRC89	0.0E+00	0.0E+00	NIOSH	6.7E+00	CRC89	6.7E+00	CRC89	6.7E+00	6.7E+00	WATER9 (U.S. EPA, 2001)															

Regional Screening Level (RSL) Chemical-specific Parameters Supporting Table June 2015

Contaminant		Molecular Weight		Volatility Parameters				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	Vapor Pressure (cmHg)	VP Ref	Density (g/cm ³)	Density Ref	D ₁₀ (cm ² /s)	D _w (cm ² /s)	Dia Ref	K _{ow} (L/kg)	K _{oc} Ref	(K _{oc} /K _{ow})	K _{oc} Ref	log K _{ow} (L/kg)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/vent)	t (hr)	K _p (cm ² /hr)	KPREF				
Bromofom	75-25-2	2.5E+02	EPI	2.2E-02	5.4E-04	EPI	5.4E+00	EPI	2.9E+00	CRC89	3.6E-02	1.0E-05	WATERS (U.S. EPA, 2001)							3.2E+01	EPI	2.4E+00	EPI	3.1E+03	EPI	1.4E-02	2.7E+00	6.6E+00	2.4E+03	EPI
Bromomethane	74-83-9	9.5E+01	EPI	3.0E-01	7.3E-03	EPI	1.0E-01	1.4E-05	1.7E+00	CRC89	1.0E-01	1.4E-05	WATERS (U.S. EPA, 2001)							1.3E+01	EPI	1.2E+00	EPI	1.5E+04	EPI	1.1E-02	3.6E-01	8.6E-01	2.8E-03	EPI
Bromophos	2104-96-3	3.7E+02	EPI	8.4E-03	2.1E-04	EPI	1.3E-04	EPI	1.7E+00	LookChem	2.3E-02	6.1E-06	WATERS (U.S. EPA, 2001)							2.0E+03	EPI	5.2E+00	EPI	3.0E-01	EPI	3.0E-01	1.2E+01	2.9E+01	4.0E-02	EPI
Bromoxynil	1689-94-5	2.9E+02	EPI	5.4E-09	1.3E-10	EPI	4.7E-08	EPI	1.7E+00	LookChem	4.5E-02	5.2E-06	WATERS (U.S. EPA, 2001)							3.3E+02	EPI	3.4E+00	EPI	1.3E+02	EPI	5.0E-02	3.7E+00	9.0E+00	7.8E-02	EPI
Bromoxynil Octanoate	1689-99-2	4.0E+02	EPI	1.3E-03	3.2E-05	EPI	4.8E-06	EPI	1.5E+00	LookChem	2.1E-02	5.4E-06	WATERS (U.S. EPA, 2001)							4.3E+03	EPI	5.4E+00	EPI	8.0E-02	EPI	2.6E-01	1.3E+01	4.6E+01	3.8E-02	EPI
Butadiene, 1,3-	109-69-1	5.4E+01	EPI	3.0E-00	7.4E-02	EPI	2.1E+03	EPI	6.1E-01	CRC89	1.0E-01	1.0E-05	WATERS (U.S. EPA, 2001)							4.0E+01	EPI	2.0E+00	EPI	7.4E+02	EPI	4.8E-02	2.1E+01	5.1E-01	1.8E-02	EPI
Butanol, n	71-36-3	7.4E+01	EPI	3.6E-04	8.8E-06	EPI	6.7E+00	EPI	8.1E-01	CRC89	9.0E-02	1.0E-05	WATERS (U.S. EPA, 2001)							3.5E+00	EPI	8.8E-01	EPI	6.3E+04	EPI	7.6E-03	2.7E-01	6.8E-01	2.3E-03	EPI
Butyl Benzyl Phthalate	85-68-7	3.1E+02	EPI	5.2E-05	1.3E-06	EPI	8.3E-06	EPI	1.1E+00	CRC89	9.1E-02	5.2E-06	WATERS (U.S. EPA, 2001)							7.2E+03	EPI	4.7E+00	EPI	2.7E+00	EPI	2.6E-01	5.9E+00	1.4E+01	3.9E-02	EPI
Butyl alcohol, sec-	78-92-2	7.4E+01	EPI	3.7E-04	9.1E-06	EPI	1.8E+01	EPI	8.1E-01	CRC89	2.0E-02	1.0E-05	WATERS (U.S. EPA, 2001)							2.9E+00	EPI	6.1E-01	EPI	1.8E+05	EPI	5.1E-03	2.7E-01	6.6E-01	1.5E-03	EPI
Butylate	2008-41-5	2.2E+02	EPI	3.5E-03	8.5E-05	EPI	1.3E-02	EPI	9.4E-01	CRC89	2.3E-02	5.8E-06	WATERS (U.S. EPA, 2001)							3.9E+02	EPI	4.2E+00	EPI	4.5E+01	EPI	3.1E-01	1.7E+00	4.2E+00	5.4E-02	EPI
Butylated hydroxyanisole	25013-16-5	1.8E+02	EPI	4.8E-05	1.2E-06	EPI	2.3E-03	EPI	6.0E-02	WATERS (U.S. EPA, 2001)	6.0E-02	7.0E-06	WATERS (U.S. EPA, 2001)							8.4E+02	EPI	3.5E+00	EPI	7.4E+02	EPI	1.7E-01	1.1E+00	2.6E+00	3.3E-02	EPI
Butylated hydroxytoluene	128-37-0	2.2E+02	EPI	1.4E-04	3.4E-06	EPI	5.2E-03	EPI	8.9E-01	CRC89	2.3E-02	5.6E-06	WATERS (U.S. EPA, 2001)							1.5E+04	EPI	5.1E+00	EPI	6.0E-01	EPI	1.3E+00	1.8E+00	7.1E+00	2.2E-01	EPI
Butylbenzene, tert-	104-51-8	1.3E+02	EPI	6.5E-01	1.6E-02	EPI	1.1E+00	EPI	8.6E-01	CRC89	5.3E-02	7.3E-06	WATERS (U.S. EPA, 2001)							1.5E+03	EPI	4.4E+00	EPI	1.2E+01	EPI	1.0E+00	5.9E-01	2.3E+00	2.3E-01	EPI
Butylbenzene, sec-	135-98-8	1.3E+02	EPI	7.2E-01	1.8E-02	EPI	1.8E+00	EPI	8.6E-01	LANGE	5.3E-02	7.3E-06	WATERS (U.S. EPA, 2001)							1.3E+03	EPI	4.6E+00	EPI	1.8E+01	EPI	1.3E+00	5.9E-01	2.3E+00	3.0E-01	EPI
Butylbenzene, tert-	98-06-6	1.3E+02	EPI	5.4E-01	1.3E-02	EPI	2.2E+00	EPI	8.7E-01	CRC89	5.3E-02	7.4E-06	WATERS (U.S. EPA, 2001)							1.0E+03	EPI	4.1E+00	EPI	3.0E+01	EPI	6.6E-01	5.9E-01	2.3E+00	1.5E-01	EPI
Cacodylic Acid	75-60-5	1.4E+02	EPI				4.6E-03				7.1E-02	8.3E-06	WATERS (U.S. EPA, 2001)							4.4E+01	EPI	3.6E-01	EPI	2.0E+06	EPI	2.1E-03	6.2E-01	1.5E+00	4.6E-04	EPI
Cadmium (Diet)	7440-43-9	1.1E+02	EPI				0.0E+00	NIOSH	8.7E+00	CRC89				7.5E+01	SSL									4.1E-03	4.5E-01	1.1E+00	1.0E-03	RAGSE		
Cadmium (Water)	7440-43-9	1.1E+02	EPI				0.0E+00	NIOSH	8.7E+00	CRC89				7.5E+01	SSL									4.4E-03	4.5E-01	1.1E+00	1.0E-03	RAGSE		
Calcium Chromate	13765-19-0	1.6E+02	CRC89																					4.8E-03	7.9E-01	1.9E+00	1.0E-03	RAGSE		
Caprocladim	105-60-2	1.1E+02	EPI	1.0E-08	2.5E-08	EPI	1.6E-03	EPI	1.0E+00	LANGE	6.9E-02	8.0E-06	WATERS (U.S. EPA, 2001)							2.5E+01	EPI	1.9E-01	YAWS	7.7E+05	EPI	4.1E-03	4.5E-01	1.1E+00	1.0E-03	EPI
Carbafol	2426-06-1	3.5E+02	EPI	2.0E-07	4.9E-09	EPI	1.5E-08	EPI	1.0E+00		3.8E-02	4.5E-06	WATERS (U.S. EPA, 2001)							7.8E+02	EPI	3.8E+00	EPI	1.4E+00	EPI	4.1E-02	9.5E+00	2.3E+01	5.8E-03	EPI
Capitan	133-06-2	3.0E+02	EPI	2.9E-07	7.0E-09	EPI	9.0E-08	EPI	1.7E+00	CRC89	2.6E-02	6.9E-06	WATERS (U.S. EPA, 2001)							2.5E+02	EPI	2.8E+00	EPI	5.1E+00	EPI	1.6E-02	5.1E+00	1.2E+01	2.3E-03	EPI
Carbaryl	63-25-2	2.0E+02	EPI	1.3E-07	3.3E-09	EPI	1.4E-06	EPI	1.2E+00	CRC89	2.7E-02	7.1E-06	WATERS (U.S. EPA, 2001)							3.5E+02	EPI	2.4E+00	EPI	1.1E+02	EPI	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI
Carbaryl	1563-66-2	2.2E+02	EPI	1.3E-07	3.1E-09	EPI	4.9E-06	EPI	1.2E+00	CRC89	2.6E-02	6.6E-06	WATERS (U.S. EPA, 2001)							9.5E+01	EPI	2.3E+00	EPI	3.2E+02	EPI	1.8E-02	1.8E+00	4.4E+00	3.1E-03	EPI
Carbon Disulfide	75-15-0	7.6E+01	EPI	5.9E-01	1.4E-02	EPI	3.6E+02	EPI	1.3E+00	CRC89	1.1E-01	1.3E-05	WATERS (U.S. EPA, 2001)							2.2E+01	EPI	1.9E+00	EPI	2.2E+03	EPI	3.8E-02	2.8E-01	6.7E-01	1.1E-02	EPI
Carbon Tetrachloride	56-23-5	1.5E+02	EPI	1.1E+00	2.8E-02	EPI	1.2E+02	EPI	1.6E+00	CRC89	5.7E-02	9.8E-06	WATERS (U.S. EPA, 2001)							4.4E+01	EPI	2.8E+00	EPI	7.9E+02	EPI	7.8E-02	7.6E-01	1.8E+00	1.6E-02	EPI
Carbosulfan	55285-14-8	3.8E+02	EPI	2.1E-05	5.1E-07	EPI	3.1E-07	EPI	1.1E+00	CRC89	1.8E-02	4.4E-06	WATERS (U.S. EPA, 2001)							1.2E+04	EPI	3.8E+00	OTHER	3.0E-01	EPI	4.3E-01	1.4E+01	3.4E+01	5.8E-02	EPI
Carboxin	5234-68-4	2.4E+02	EPI	1.3E-08	3.2E-10	EPI	1.5E-07	EPI	1.0E+00		5.0E-02	5.8E-06	WATERS (U.S. EPA, 2001)							1.7E+02	EPI	2.1E+00	EPI	1.5E+02	EPI	1.2E-02	2.2E+00	5.2E+00	2.0E-03	EPI
Ceric oxide	1306-38-3	1.7E+02	CRC89						7.2E+00	CRC89														5.0E-03	9.7E-01	2.3E+00	1.0E-03	RAGSE		
Chloral Hydrate	302-17-0	1.7E+02	EPI	4.5E-09	1.1E-10	EPI	1.5E+01	EPI	1.9E+00	CRC89	5.4E-02	1.0E-05	WATERS (U.S. EPA, 2001)							1.0E+00	EPI	9.9E-01	EPI	7.9E+05	EPI	4.2E-03	8.9E-01	2.1E+00	8.4E-04	EPI
Chloramben	133-90-4	2.1E+02	EPI	1.6E-09	3.9E-11	EPI	1.5E-05	EPI	1.0E+00		5.4E-02	6.4E-06	WATERS (U.S. EPA, 2001)							2.1E+01	EPI	1.9E+00	EPI	7.0E+02	EPI	1.1E-02	1.5E+00	3.6E+00	2.0E-03	EPI
Chloraniol	118-75-2	2.5E+02	EPI	1.3E-08	3.3E-10	EPI	2.3E-06	EPI	1.0E+00	PubChem	4.8E-02	5.7E-06	WATERS (U.S. EPA, 2001)							3.1E+02	EPI	2.2E+00	EPI	2.5E+02	EPI	1.2E-02	2.5E+00	6.0E+00	1.9E-03	EPI
Chlorane	12789-03-6	4.1E+02	EPI	2.9E-03	7.0E-05	EPI	1.0E-05	EPI	1.6E+00	PubChem	2.1E-02	5.4E-06	WATERS (U.S. EPA, 2001)							3.4E+04	EPI	6.3E+00	OTHER	1.3E-02	PhysProp	8.3E-01	2.1E+01	8.0E+01	1.1E-01	EPI
Chloroacene (Kepone)	143-50-0	4.9E+02	EPI	2.2E-06	5.4E-08	EPI	2.3E-07	EPI	1.6E+00	CRC89	2.9E-02	4.9E-06	WATERS (U.S. EPA, 2001)							1.8E+04	EPI	5.4E+00	EPI	2.7E+00	EPI	9.3E-02	5.5E-01	1.4E+02	1.1E-02	EPI
Chlorfenvinphos	470-90-6	3.6E+02	EPI	1.2E-06	2.9E-08	EPI	7.5E-06	EPI	1.0E+00		3.9E-02	4.4E-06	WATERS (U.S. EPA, 2001)							1.3E+03	EPI	3.8E+00	EPI	1.2E+02	EPI	3.7E-02	1.1E+01	2.6E+01	5.1E-03	EPI
Chlorfomum, Ethyl-	90980-32-4	4.1E+02	EPI	7.4E-14	1.8E-16	EPI	4.0E-12	EPI	1.0E+00		3.4E-14	4.0E-06	WATERS (U.S. EPA, 2001)							7.2E+01	EPI	2.5E+00	EPI	1.2E+03	EPI	2.6E-03	2.2E+01	5.3E-01	3.4E-04	EPI
Chlorine	7782-50-5	7.1E+01	EPI	4.8E-01	1.2E-02	PHYSPROP	5.9E+03	EPI	2.9E+00	CRC89	1.1E-01	2.2E-05	WATERS (U.S. EPA, 2001)	2.5E-01	BAES					8.5E-01	OTHER			6.3E+03	EPI	3.2E-03	2.8E-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	2.8E+00	CRC89	1.6E-01	2.2E-05	WATERS (U.S. EPA, 2001)											3.2E-03	2.5E-01	6.0E-01	1.0E-03	RAGSE		

Contaminant		Molecular Weight		Volatility Parameters				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	Vapor Pressure	VP Ref	Density (g/cm ³)	Density Ref	D ₁₀ (cm ² /s)	D _w (cm ² /s)	Dia Ref	K _{ow}	K _{oc} Ref	(K _{oc} /K _{ow})	K _{oc} Ref	log K _{ow} (L/kg)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr)	t (hr)	K _p (cm ² /hr)	KPREF		
Glutaraldehyd	111-30-8	1.0E+02	EPI	9.8E-07	2.4E-08	EPI	6.0E-01	EPI			8.8E-02	1.0E-05	WATERS (U.S. EPA, 2001)					1.0E+00	EPI	-1.8E-01	EPI	7.1E+05	EPI	1.3E-03	3.8E-01	9.2E-01	3.3E-04	EPI
Glycidyl	765-34-4	7.2E+01	EPI	3.2E-05	7.8E-07	EPI	2.2E+01	EPI	1.1E+00	CRC89	1.1E-01	1.3E-05	WATERS (U.S. EPA, 2001)					1.0E+00	EPI	-1.2E-01	EPI	1.0E+06	EPI	1.7E-03	2.7E-01	6.4E-01	5.2E-04	EPI
Glyphosate	1071-83-6	1.7E+02	EPI	8.6E-11	2.1E-12	EPI	9.8E-08	EPI			6.2E-02	7.3E-06	WATERS (U.S. EPA, 2001)					2.1E+03	USDA ARS	-3.4E+00	EPI	1.1E+04	EPI	2.3E-07	9.3E-01	2.2E+00	4.5E-08	EPI
Coal	42874-03-3	3.9E+02	EPI	3.4E-05	8.2E-07	EPI	2.0E-07	EPI	1.4E+00	CRC89	2.4E-02	5.3E-06	WATERS (U.S. EPA, 2001)					4.0E+04	EPI	4.7E+00	EPI	1.2E-01	EPI	1.5E-01	1.1E+01	2.7E+01	2.0E-02	EPI
Quarantine	115-00-8	5.9E+01	EPI	9.6E-10	2.3E-11	EPI	2.2E+00	EPI	1.0E+00	GuideChem	1.4E-01	1.7E-05	WATERS (U.S. EPA, 2001)					1.2E+01	EPI	1.8E+00	EPI	1.8E+03	EPI	3.8E-04	2.3E-01	5.4E-01	6.3E-05	EPI
Quarantine Chloride	50-01-1	9.6E+01	EPI	7.3E-17	1.8E-18	EPI	2.4E-06	EPI	1.4E+00	CRC89	9.3E-02	1.2E-05	WATERS (U.S. EPA, 2001)					1.0E+00	OTHER	1.0E+06	EPI	1.5E-07	EPI	3.6E-01	8.7E-01	3.9E-08	EPI	
Guthion	86-50-0	3.2E+02	EPI	9.8E-07	2.4E-08	EPI	1.6E-06	EPI	1.4E+00	CRC89	2.3E-02	6.0E-06	WATERS (U.S. EPA, 2001)					5.2E+01	EPI	2.8E+00	EPI	2.1E+01	EPI	1.2E-02	6.3E+00	1.5E+01	1.8E-03	EPI
Haloxyflor, Methyl	69896-40-2	3.8E+02	EPI	1.3E-05	3.2E-07	EPI	6.0E-06	EPI			3.6E-02	4.3E-06	WATERS (U.S. EPA, 2001)					5.5E+03	EPI	4.1E+00	EPI	9.3E+00	EPI	4.5E-02	1.3E+01	3.2E+01	6.0E-03	EPI
Harmony	79277-27-3	3.9E+02	EPI	1.7E-12	4.1E-14	EPI	1.3E-10	EPI			3.6E-02	4.2E-06	WATERS (U.S. EPA, 2001)					5.1E+01	EPI	1.6E+00	EPI	2.2E+03	EPI	8.6E-04	1.6E+01	3.7E+01	1.1E-04	EPI
Heptachlor	76-44-8	3.7E+02	EPI	1.2E-02	2.9E-04	EPI	4.0E-04	EPI	1.6E+00	CRC89	2.2E-02	5.7E-06	WATERS (U.S. EPA, 2001)					4.1E+04	EPI	6.1E+00	EPI	1.8E-01	EPI	1.1E+00	1.3E+01	5.0E+01	1.4E-01	EPI
Heptachlor Epoxide	1024-57-3	3.9E+02	EPI	8.6E-04	2.1E-05	EPI	2.0E-05	EPI	1.9E+00	LookChem	2.4E-02	6.2E-06	WATERS (U.S. EPA, 2001)					1.0E+04	EPI	5.0E+00	EPI	2.0E-01	EPI	1.6E-01	1.6E+01	3.8E+01	2.1E-02	EPI
Hexabromobenzene	87-82-1	5.5E+02	EPI	1.1E-03	2.8E-05	EPI	1.7E-08	EPI	3.0E+00	LookChem	2.5E-02	6.6E-06	WATERS (U.S. EPA, 2001)					2.8E+03	EPI	6.1E+00	EPI	1.6E-04	EPI	1.2E-01	3.1E+02	3.1E+02	1.4E-02	EPI
Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BDE-153)	68631-49-2	6.4E+02	OTHER				5.8E-06	IRIS			2.5E-02	3.0E-06	WATERS (U.S. EPA, 2001)					3.8E+00	IRIS	9.0E-04	IRIS	1.0E-04	IRIS	4.2E+02	1.0E+03			
Hexachlorobenzene	118-74-1	2.8E+02	EPI	7.0E-02	1.7E-03	EPI	1.8E-05	EPI	2.0E+00	CRC89	2.9E-02	7.8E-06	WATERS (U.S. EPA, 2001)					6.2E+03	EPI	5.7E+00	EPI	6.2E-03	EPI	1.6E+00	4.1E+00	1.7E+01	2.6E-01	EPI
Hexachlorobutadiene	87-88-3	2.6E+02	EPI	4.2E-01	1.0E-02	EPI	2.2E-01	EPI	1.6E+00	CRC89	2.7E-02	7.0E-06	WATERS (U.S. EPA, 2001)					8.5E+02	EPI	4.8E+00	EPI	3.2E+00	EPI	5.0E-01	3.0E+00	7.3E+00	8.1E-02	EPI
Hexachlorocyclohexane, Alpha	319-84-6	2.9E+02	EPI	2.1E-04	5.1E-06	EPI	3.5E-05	EPI	1.9E+00	CRC89	4.3E-02	5.1E-06	WATERS (U.S. EPA, 2001)					2.8E+03	EPI	3.8E+00	EPI	2.0E+00	EPI	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclohexane, Beta	319-85-7	2.9E+02	EPI	2.1E-04	5.1E-06	EPI	3.5E-05	EPI	1.9E+00	CRC89	2.8E-03	7.4E-06	WATERS (U.S. EPA, 2001)					2.8E+03	EPI	3.8E+00	EPI	2.4E-01	EPI	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclohexane, Gamma-(Lindane)	58-89-9	2.9E+02	EPI	2.1E-04	5.1E-06	EPI	4.2E-05	EPI			4.3E-02	5.1E-06	WATERS (U.S. EPA, 2001)					2.8E+03	EPI	3.7E+00	EPI	7.3E+00	EPI	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclohexane, Technic	608-73-1	2.9E+02	EPI	2.1E-04	5.1E-06	EPI	3.5E-05	EPI			2.8E+03	5.1E-06	WATERS (U.S. EPA, 2001)					2.8E+03	EPI	4.1E+00	EPI	9.0E+00	EPI	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclopentadiene	77-47-4	2.7E+02	EPI	1.1E+00	2.7E-02	EPI	6.0E-02	EPI	1.7E+00	CRC89	3.7E-02	7.2E-06	WATERS (U.S. EPA, 2001)					1.4E+03	EPI	5.0E+00	EPI	1.8E+00	EPI	6.5E-01	2.3E+00	1.4E+01	1.0E-02	EPI
Hexachloroethane	67-72-1	2.4E+02	EPI	1.6E-01	3.9E-03	EPI	2.1E-01	EPI	2.1E+00	CRC89	2.0E-02	1.2E-05	WATERS (U.S. EPA, 2001)					2.0E+02	EPI	4.1E+00	EPI	5.0E-01	EPI	2.5E-01	2.2E+00	5.3E+00	4.2E-02	EPI
Hexachlorophene	70-30-4	4.1E+02	EPI	2.2E-11	5.5E-13	EPI	8.3E-11	EPI			3.5E-02	4.0E-06	WATERS (U.S. EPA, 2001)					6.7E+05	EPI	7.5E+00	EPI	1.4E+02	EPI	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	EPI	8.2E-10	2.0E-11	EPI	4.1E-09	EPI	1.8E+00	CRC89	3.1E-02	8.5E-06	WATERS (U.S. EPA, 2001)					8.9E+01	EPI	8.7E-01	EPI	6.0E+01	EPI	1.9E-03	1.8E+00	4.4E+00	3.4E-04	EPI
Hexamethylene Dicyclohexane, 1,6-	822-06-0	1.7E+02	EPI	2.0E-03	4.8E-05	EPI	2.1E-02	EPI	1.1E+00	CRC89	4.0E-02	7.2E-06	WATERS (U.S. EPA, 2001)					4.8E+03	EPI	3.2E+00	EPI	1.8E+02	EPI	1.2E-01	9.2E-01	2.2E+00	2.4E-02	EPI
Hexamethylphosphoramide	680-31-9	1.8E+02	EPI	2.9E-10	7.1E-12	EPI	4.6E-02	EPI	1.0E+00	CRC89	3.5E-02	6.9E-06	WATERS (U.S. EPA, 2001)					1.0E+01	EPI	2.8E-01	EPI	1.0E+06	EPI	1.2E-03	1.1E+00	2.5E+00	2.0E-04	EPI
Hexane, N-	110-54-3	8.6E+01	EPI	7.4E+01	1.8E+00	EPI	1.5E-02	EPI	6.6E-01	CRC89	7.3E-02	8.2E-06	WATERS (U.S. EPA, 2001)					1.3E+02	EPI	3.9E+00	EPI	9.5E+00	EPI	7.2E-01	3.2E-01	1.2E+00	2.0E-01	EPI
Hexanedioic Acid	124-04-9	1.5E+02	EPI	1.9E-10	4.7E-12	EPI	3.2E-07	EPI	1.4E+00	CRC89	5.8E-02	9.2E-06	WATERS (U.S. EPA, 2001)					2.4E+01	EPI	8.0E-02	EPI	3.1E+04	EPI	1.2E-03	6.9E-01	1.7E+00	2.7E-04	EPI
Hexanone, 2	591-78-6	1.0E+02	EPI	3.8E-03	9.3E-05	EPI	1.2E+01	EPI	8.1E-01	CRC89	7.0E-02	8.4E-06	WATERS (U.S. EPA, 2001)					1.5E+01	EPI	1.4E+00	EPI	1.7E+04	EPI	1.4E-02	3.8E-01	9.2E-01	3.6E-03	EPI
Hexazinone	51235-04-2	2.5E+02	EPI	9.2E-11	2.3E-12	PubChem	2.3E-07	EPI	1.3E+00	CRC89	2.5E-02	6.3E-06	WATERS (U.S. EPA, 2001)					1.3E+02	EPI	1.9E+00	EPI	3.3E+04	EPI	6.2E-03	2.7E+00	6.5E+00	1.0E-03	EPI
Hydrazine	302-01-2	3.2E+01	EPI	2.5E-05	6.1E-07	PubChem	1.4E+01	EPI	1.0E+00	CRC89	1.7E-01	1.9E-05	WATERS (U.S. EPA, 2001)					-2.1E+00	EPI	1.9E+00	EPI	3.3E+04	EPI	9.5E-05	1.6E-01	3.8E-01	4.4E-05	RAGSE
Hydrazine Sulfate	10034-93-2	1.3E+02	EPI						1.4E+00	CRC89											3.1E+04	PubChem	4.4E-03	5.5E-01	1.3E+00	1.0E-03	RAGSE	
Hydrogen Chloride	7647-01-0	3.5E+01	EPI	8.3E-07	2.0E+06	Toxnet HSDB	3.5E+04	PubChem	1.5E+00	CRC89	1.9E-01	2.3E-05	WATERS (U.S. EPA, 2001)					2.3E-01	OTHER	3.7E+05	Toxnet HSDB	2.3E-03	OTHER	1.7E-01	7.7E-01	4.0E-01	1.0E-03	RAGSE
Hydrogen Fluoride	7664-39-3	2.0E+01	CRC89	4.3E-03	1.0E-04	PhysProp	9.2E+02	PhysProp	8.2E-01	CRC89	2.2E-01	2.2E-05	WATERS (U.S. EPA, 2001)					2.3E-01	OTHER	1.0E+06	PhysProp	1.7E-03	OTHER	1.4E-01	1.6E-01	3.3E-01	1.0E-03	RAGSE
Hydrogen Sulfide	7783-06-4	3.4E+01	CRC89	3.5E-01	8.6E-03	PhysProp	1.6E+04	PhysProp	1.4E+00	CRC89	1.9E-01	2.2E-05	WATERS (U.S. EPA, 2001)					2.3E-01	OTHER	4.4E+06	PhysProp	2.2E-03	OTHER	1.6E-01	3.9E-01	1.0E-03	RAGSE	
Hydroquinone	123-31-9	1.1E+02	EPI	1.9E-09	4.7E-11	EPI	2.4E-05	EPI	1.3E+00	CRC89	8.0E-02	1.1E-05	WATERS (U.S. EPA, 2001)					2.4E+02	EPI	5.9E-01	EPI	7.2E+04	EPI	3.8E-03	4.3E-01	1.0E+00	3.9E-04	EPI
Imazali	35554-44-0	3.0E+02	EPI	1.1E-07	2.6E-09	EPI	1.2E-06	EPI	1.2E+00	CRC89	2.1E-02	3.7E-06	WATERS (U.S. EPA, 2001)					8.5E+03	EPI	1.9E+00	EPI	7.7E-02	EPI	7.7E-02	4.9E+00	1.2E-01	1.2E-02	EPI
Imazaliquin	81335-37-7	3.1E+02	EPI	2.8E-16	6.9E-18	EPI	3.4E-13	EPI			4.0E-02	4.8E-06	WATERS (U.S. EPA, 2001)					2.4E+03	EPI	1.9E+00	EPI	9.0E-01	EPI	3.3E-03	5.8E+00	1.4E+01	9.8E-04	EPI
Iodine	7553-56-2	2.5E+02	EPI				2.3E-01	EPI	4.9E+00	CRC89				6.0E+01	BAES						3.3E+02	EPI	6.1E-03	2.8E+00	6.7E+00	1.0E-03	RAGSE	
Iprodione	36734-19-7	3.3E+02	EPI	1.3E-07	3.1E-09	EPI	3.8E-09	EPI			4.0E-02	4.6E-06	WATERS (U.S. EPA, 2001)			</												

Regional Screening Level (RSL) Chemical-specific Parameters Supporting Table June 2015

Contaminant	Molecular Weight		Volatility Parameters				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	Vapor Pressure	VP Ref	Density (g/cm ³)	Density Ref	Dia (cm ² /s)	DW (cm ² /s)	Dia Ref	K _{ow} (L/kg)	K _{oc} Ref	K _{oc} (L/kg)	K _{oc} Ref	log K _{ow} (L/kg)	log K _{oc} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr)	t (hr)	K _p (cm ² /hr)	KPREF	
Methoxychlor	72-43-5	3.5E+02	EPI	8.3E-06	2.0E-07	EPI	4.2E-05	EPI	1.4E+00	CRC89	2.2E-02	5.6E-06	WATER9 (U.S. EPA, 2001)	2.7E+04	EPI	5.1E+00	EPI	1.0E-01	EPI	1.0E-01	EPI	3.1E-01	9.1E+00	2.2E+01	4.3E-02	EPI		
Methoxyethanol Acetate, 2	110-49-6	1.2E+02	EPI	1.2E-05	3.1E-07	EPI	2.0E+00	EPI	1.0E+00	CRC89	6.8E-02	8.7E-06	WATER9 (U.S. EPA, 2001)	2.5E+00	EPI	1.0E-01	EPI	1.0E-01	EPI	1.0E+06	EPI	1.7E-03	4.8E-01	1.2E+00	4.0E-04	EPI		
Methoxyethanol, 2-	109-86-4	7.6E+01	EPI	1.4E-05	3.3E-07	EPI	9.5E+00	EPI	9.6E-01	CRC89	9.5E-02	1.1E-05	WATER9 (U.S. EPA, 2001)	1.0E+00	EPI	-7.7E-01	EPI	1.0E+06	EPI	6.0E-04	EPI	2.8E-01	2.8E-01	6.7E-01	1.8E-04	EPI		
Methyl Acetate	79-20-9	7.4E+01	EPI	4.7E-03	1.2E-04	EPI	2.2E+02	EPI	9.3E-01	CRC89	9.6E-02	1.1E-05	WATER9 (U.S. EPA, 2001)	3.1E+00	EPI	1.8E-01	EPI	2.0E+05	EPI	2.6E-03	EPI	2.7E-01	6.6E-01	7.9E-04	1.9E-03	EPI		
Methyl Acrylate	96-33-3	9.6E+01	EPI	9.1E-03	2.0E-04	EPI	8.7E+01	EPI	9.5E-01	CRC89	9.1E-03	1.0E-05	WATER9 (U.S. EPA, 2001)	5.8E+00	EPI	8.9E-01	EPI	4.9E+04	EPI	6.2E-03	EPI	3.2E-01	7.7E-01	1.8E-03	1.9E-03	EPI		
Methyl Ethyl Ketone (2-Butanone)	78-93-3	7.2E+01	EPI	2.3E-03	5.7E-05	EPI	9.1E+01	EPI	8.0E-01	CRC89	9.1E-02	1.0E-05	WATER9 (U.S. EPA, 2001)	4.5E+00	EPI	2.9E-01	EPI	2.7E+05	EPI	3.1E-03	EPI	2.7E-01	6.4E-01	9.6E-04	1.9E-03	EPI		
Methyl Hydrazine	60-34-4	4.6E+01	EPI	1.3E-06	3.2E-08	EPI	5.0E+01	EPI	8.7E-01	LANGE	1.3E+01	1.4E-05	WATER9 (U.S. EPA, 2001)	1.3E+01	EPI	-1.1E+00	EPI	1.0E+06	EPI	4.5E-04	EPI	1.9E-01	4.6E-01	1.7E-04	1.7E-04	EPI		
Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1	1.0E+02	EPI	5.6E-03	1.4E-04	EPI	2.0E+01	EPI	8.0E-01	CRC89	7.0E-02	8.3E-06	WATER9 (U.S. EPA, 2001)	1.3E+01	EPI	1.3E+00	EPI	1.9E+04	EPI	1.2E-02	EPI	3.8E-01	9.2E-01	3.2E-03	2.0E-03	EPI		
Methyl Isocyanate	624-83-9	5.7E+01	EPI	3.8E-02	9.3E-04	EPI	3.5E+02	EPI	1.2E-01	CRC89	1.2E-01	1.3E-05	WATER9 (U.S. EPA, 2001)	4.0E+01	EPI	7.9E-01	YAW5	4.8E+04	EPI	7.3E-03	EPI	2.2E-01	2.5E-01	5.3E-01	2.5E-03	EPI		
Methyl Methacrylate	80-62-6	1.0E+02	EPI	1.3E-02	3.2E-04	EPI	3.9E+01	EPI	9.4E-01	CRC89	7.5E-02	9.2E-06	WATER9 (U.S. EPA, 2001)	9.1E+00	EPI	1.4E+00	EPI	1.5E+04	EPI	1.4E-02	EPI	3.8E-01	9.2E-01	3.8E-03	4.0E-03	EPI		
Methyl Parathion	298-00-0	2.6E+02	EPI	4.1E-06	1.0E-07	EPI	2.5E-06	EPI	1.4E+00	CRC89	2.5E-02	6.4E-06	WATER9 (U.S. EPA, 2001)	7.3E+02	EPI	2.9E+00	EPI	3.8E+01	EPI	2.6E-02	EPI	1.4E-02	3.1E+00	7.5E+00	4.2E-03	EPI		
Methyl Phosphonic Acid	993-13-5	9.6E+01	EPI	5.0E-10	1.2E-11	EPI	3.3E-04	EPI	3.3E-04	EPI	9.1E-02	1.1E-05	WATER9 (U.S. EPA, 2001)	1.4E+00	EPI	-1.0E+00	EPI	2.0E+04	EPI	3.7E-04	EPI	3.6E-01	8.7E-01	9.8E-05	1.9E-05	EPI		
Methyl Styrene (Mixed Isomers)	25013-15-4	1.2E+02	EPI	1.2E-01	3.0E-03	EPI	1.7E+00	EPI	8.9E-01	HSDB	6.2E-02	8.1E-06	WATER9 (U.S. EPA, 2001)	7.2E+02	EPI	3.4E+00	EPI	4.8E+01	EPI	2.8E-01	EPI	4.8E-01	1.2E+00	6.6E-02	6.0E-02	EPI		
Methyl methanesulfonate	66-27-3	1.1E+02	EPI	1.6E-04	4.0E-06	EPI	3.1E-01	EPI	1.3E+00	CRC89	7.9E-02	1.1E-05	WATER9 (U.S. EPA, 2001)	4.3E+00	EPI	-6.6E-01	EPI	2.0E+05	LANGE	5.6E-04	EPI	4.4E-01	1.0E+00	1.4E-04	1.4E-04	EPI		
Methyl tert-Butyl Ether (MTBE)	1634-04-4	8.8E+01	EPI	2.4E-02	5.9E-04	EPI	2.5E+02	EPI	7.4E-01	CRC89	7.5E-02	8.6E-06	WATER9 (U.S. EPA, 2001)	1.2E+01	EPI	9.4E-01	EPI	5.1E+04	EPI	7.6E-03	EPI	3.3E-01	7.9E-01	2.1E-03	2.1E-03	EPI		
Methyl-1,4-benzenediamine dichloride, 2-	615-45-2	2.0E+02	EPI	2.6E-16	6.4E-18	EPI	4.1E-12	EPI	1.2E+00	EPI	5.6E-02	6.6E-06	WATER9 (U.S. EPA, 2001)	2.0E+02	EPI	-2.1E+00	EPI	3.9E+02	EPI	2.9E-05	EPI	1.3E+00	3.1E+00	1.5E+00	5.4E-06	EPI		
Methyl-5-Nitroaniline, 2-	99-55-8	1.5E+02	EPI	7.9E-07	1.9E-08	EPI	5.9E-04	EPI	1.2E+00	EPI	6.7E-02	7.8E-06	WATER9 (U.S. EPA, 2001)	1.9E+00	EPI	1.9E+00	EPI	6.1E+02	EPI	1.8E-02	EPI	7.5E-01	1.8E+00	3.8E-03	3.8E-03	EPI		
Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	1.5E+02	EPI	5.0E-11	1.2E-12	EPI	1.2E-04	EPI	1.2E-04	EPI	6.8E-02	8.0E-06	WATER9 (U.S. EPA, 2001)	7.2E+01	EPI	-9.2E-01	EPI	1.0E+06	EPI	2.7E-04	EPI	7.0E-01	1.7E+00	5.7E-05	5.7E-05	EPI		
Methylaniline Hydrochloride, 2-	636-21-5	1.4E+02	EPI	8.0E-12	2.0E-13	EPI	1.6E-06	EPI	1.6E-06	EPI	6.9E-02	8.1E-06	WATER9 (U.S. EPA, 2001)	1.2E+02	EPI	-2.1E+00	EPI	4.8E+02	EPI	4.8E-05	EPI	6.7E-01	1.6E+00	1.1E-05	1.1E-05	EPI		
Methylanilinic acid	124-58-3	1.4E+02	EPI	1.4E+02		EPI	1.2E-04	EPI	1.2E-04	EPI	7.0E-02	8.2E-06	WATER9 (U.S. EPA, 2001)	6.4E+00	EPI	4.4E+01	EPI	1.2E+00	EPI	2.6E+05	EPI	1.9E-04	6.4E-01	1.5E+00	4.2E-05	EPI		
Methylbenzene, 1,4-diamine monohydrochloride, 2-	74172-12-7	1.6E+02	OTHER																									
Methylbenzene, 1,4-diamine sulfate, 2-	615-50-9	2.2E+02	OTHER																									
Methylchlorobenzene, 1	56-49-5	2.7E+02	EPI	2.1E-04	5.2E-06	EPI	4.3E-08	EPI	1.3E+00	CRC89	2.4E-02	6.1E-06	WATER9 (U.S. EPA, 2001)	9.6E+05	EPI	6.4E+00	EPI	2.9E-03	EPI	5.7E+00	EPI	3.3E+00	1.5E+01	9.0E-01	9.0E-01	EPI		
Methylene Chloride	75-09-2	8.5E+01	EPI	1.0E-01	3.3E-03	EPI	4.4E+02	EPI	1.3E+00	CRC89	1.0E-01	1.3E-05	WATER9 (U.S. EPA, 2001)	2.2E+01	EPI	1.3E+00	EPI	1.3E+04	EPI	1.3E-02	EPI	3.1E-01	7.5E-01	3.1E-01	3.5E-03	EPI		
Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	2.7E+02	EPI	4.7E-10	1.1E-11	EPI	3.9E-06	EPI	3.9E-06	EPI	4.6E-02	5.4E-06	WATER9 (U.S. EPA, 2001)	5.7E+03	EPI	3.9E+00	EPI	1.4E+01	EPI	1.2E-01	EPI	3.3E+00	7.9E+00	8.0E-02	8.0E-02	EPI		
Methylene-bis(N,N-dimethyl) Aniline, 4,4'	101-61-1	2.5E+02	EPI	4.9E-06	1.2E-07	EPI	3.3E-06	EPI	3.3E-06	EPI	4.7E-02	5.5E-06	WATER9 (U.S. EPA, 2001)	2.7E+03	EPI	4.4E+00	EPI	5.5E+00	EPI	5.2E-01	EPI	3.8E+00	6.7E+00	2.4E-02	RAGSE			
Methylenedibenzeneamine, 4,4'-	101-77-9	2.0E+02	EPI	6.5E-10	1.6E-11	EPI	2.1E-06	EPI	2.1E-06	EPI	5.6E-02	6.5E-06	WATER9 (U.S. EPA, 2001)	2.1E+03	EPI	1.6E+00	EPI	1.0E+03	EPI	7.5E-03	EPI	1.4E+00	3.3E+00	1.4E-03	1.4E-03	EPI		
Methylenediphenyl Disocyanate	101-68-8	2.5E+02	EPI	3.7E-05	9.0E-07	EPI	5.0E-06	EPI	1.2E+00	CRC89	2.4E-02	6.2E-06	WATER9 (U.S. EPA, 2001)	2.8E+05	EPI	5.2E+00	YAW5	1.8E+00	EPI	1.1E+00	EPI	2.7E+00	1.0E+01	1.8E-01	1.8E-01	EPI		
Methylstyrene, Alpha	98-83-9	1.2E+02	EPI	1.0E-01	2.6E-03	EPI	1.9E+00	EPI	9.1E-01	CRC89	6.3E-02	8.2E-06	WATER9 (U.S. EPA, 2001)	7.0E+02	EPI	3.5E+00	EPI	1.2E+02	EPI	2.9E-01	EPI	4.8E-01	1.2E+00	7.0E-02	7.0E-02	EPI		
Metalachlor	51218-45-2	2.8E+02	EPI	3.7E-07	9.0E-09	EPI	3.1E-05	EPI	1.1E-01	CRC89	3.2E-02	5.5E-06	WATER9 (U.S. EPA, 2001)	4.9E+02	EPI	3.1E+00	EPI	3.5E+02	EPI	2.2E-02	EPI	4.1E+00	9.8E+00	3.4E-03	3.4E-03	EPI		
Metrizolin	21087-64-9	2.1E+02	EPI	4.8E-09	1.2E-10	EPI	4.4E-07	EPI	1.3E+00	CRC89	2.7E-02	7.1E-06	WATER9 (U.S. EPA, 2001)	5.3E+01	EPI	1.7E+00	EPI	1.1E+03	EPI	7.4E-03	EPI	1.7E+00	4.0E+00	1.3E-03	1.3E-03	EPI		
Mineral oils	8012-95-1	1.7E+02	EPI	3.3E+02	8.2E+00	EPI	1.4E-01	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	EPI	9.5E-01	4.3E+00	2.0E+00	2.0E+00	EPI		
Mirex	2385-85-6	5.5E+02	EPI	3.3E-02	8.1E-04	EPI	8.0E-07	EPI	2.3E+00	ChemNet	3.2E-02	5.6E-06	WATER9 (U.S. EPA, 2001)	3.6E+05	EPI	6.9E+00	EPI	8.5E-02	EPI	4.6E-01	EPI	1.2E+02	2.9E+02	5.2E-02	5.2E-02	EPI		
Molinate	2212-67-1	1.9E+02	EPI	1.7E-04	4.1E-06	EPI	5.6E-03	EPI	1.1E+00	CRC89	2.2E-02	6.8E-06	WATER9 (U.S. EPA, 2001)	1.8E+02	EPI	3.2E+00	EPI	7.9E-02	EPI	9.9E-02	EPI	1.2E+00	2.9E+00	1.9E-02	1.9E-02	EPI		
Molybdenum	7439-98-7	9.6E+01	EPI				0.0E+00	EPI	1.0E+00	NIOSH				2.0E+01	BAES													
Monochloramine	10599-90-3	5.1E+01	EPI																									
Monomethylaniline	100-61-8	1.1E+02	EPI	3.6E-04	8.9E-06	EPI	4.5E-01	EPI	9.9E-01	CRC89	7.2E-02	9.1E-06	WATER9 (U.S. EPA, 2001)	8.2E+01	EPI	1.7E+00	EPI	5.6E+03	EPI	2.0E-02	EPI	4.2E-01	1.0E+00	5.0E-03	5.0E-03	EPI		
N,N-Diphenyl-1,4-benzenediamin	74-31-7	2.6E+02	EPI	8.4E-09	2.1E-10	EPI	6.4E-09	EPI	2.9E+00	EPI	4.7E-02	5.4E-06	WATER9 (U.S. EPA, 2001)	5.2E+04	EPI	4.0E+00	YAW5	1.6E+00	EPI	1.6E-01	EPI	3.0E+00	1.0E+00	7.2E+00	2.6E-02	EPI		
Naled	3007-75-5	3.8E+02	EPI	2.7E-03	6.5E-05	EPI	2.0E-04	EPI	9.0E-01	CRC89	2.5E-02	6.4E-06	WATER9 (U.S. EPA, 2001)	1.3E+02	EPI	1.4E+00	EPI	1.5E+00	EPI	1.7E-04	EPI	1.4E+00	3.4E+01	9.4E-05	9.4E-05	EPI		
Naphtha, High Flash Aromatic (HFAN)	64742-95-6	1.8E+02	EPI	1.8E-02	4.4E-04	EPI	8.5E-02	EPI	1.6E+00																			

Regional Screening Level (RSL) Chemical-specific Parameters Supporting Table June 2015

Contaminant		Molecular Weight		Volatility Parameters				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	Vapor Pressure	VP Ref	Density (g/cm ³)	Density Ref	D _{air} (cm ² /s)	D _w (cm ² /s)	Dia Ref	K _{ow} (L/kg)	K _{oc} Ref	K _{oc} (L/kg)	K _{oc} Ref	log K _{ow}	log K _{oc} Ref	S (mg/L)	S Ref	B (unitless)	t _{event} (hr)	t (hr)	K _p (cm/hr)	KPREF	
Oxadiazon	19666-30-9	3.5E+02	EPI	3.0E-06	7.3E-08	EPI	1.1E-07	EPI			3.9E-02	4.5E-06	WATERS (U.S. EPA, 2001)			5.0E+03	EPI	4.8E+00	EPI	7.0E-01	EPI	2.0E-01	9.0E+00	2.2E+01	2.8E-02	EPI	
Oxamyl	23135-22-0	2.2E+02	EPI	9.7E-09	2.4E-10	EPI	2.3E-04	EPI	9.7E-01	CRC89	2.3E-02	5.9E-06	WATERS (U.S. EPA, 2001)			1.0E+01	EPI	4.7E-01	EPI	2.8E+05	EPI	2.6E-04	1.8E+00	4.3E+00	4.5E-05	EPI	
Paclitaxel	76738-62-0	2.9E+02	EPI	3.4E-09	8.3E-11	EPI	7.5E-09	EPI	1.2E+00	CRC89	2.4E-02	5.7E-06	WATERS (U.S. EPA, 2001)			9.2E+02	EPI	3.2E+00	EPI	2.6E+01	EPI	3.1E-02	4.6E+00	1.1E+01	4.7E-03	EPI	
Parquat Dichloride	1910-42-5	2.6E+02	EPI	1.3E-11	3.2E-13	EPI	1.0E-07	EPI			4.7E-02	5.5E-06	WATERS (U.S. EPA, 2001)			6.8E+03	EPI	4.5E+00	EPI	7.7E+05	EPI	3.8E-07	2.9E+00	7.0E+00	5.8E-08	EPI	
Permethrin	5638-2	2.0E+02	EPI	1.2E-05	3.0E-07	EPI	6.7E-06	EPI	1.3E+00	CRC89	2.3E-02	5.8E-06	WATERS (U.S. EPA, 2001)			2.4E+03	EPI	3.8E+00	EPI	1.1E+01	EPI	9.4E-02	4.5E+00	1.1E+01	1.3E-02	EPI	
Perubate	1114-71-2	2.0E+02	EPI	9.7E-03	2.4E-04	EPI	8.9E-02	EPI	9.5E-01	CRC89	2.4E-02	6.1E-06	WATERS (U.S. EPA, 2001)			3.0E+02	EPI	3.8E+00	EPI	1.0E+02	EPI	2.2E-01	1.4E+00	3.5E+00	4.0E-02	EPI	
Pendimethalin	40487-42-1	2.8E+02	EPI	3.5E-05	8.6E-07	EPI	3.0E-05	EPI	1.2E+00	CRC89	2.3E-02	5.7E-06	WATERS (U.S. EPA, 2001)			5.6E+03	EPI	5.2E+00	EPI	3.0E-01	EPI	7.4E-01	4.0E+00	1.5E+01	1.2E-01	EPI	
Pentabromodiphenyl Ether	32534-81-9	5.6E+02	EPI	1.4E-04	3.5E-06	EPI	3.1E-08	EPI	2.2E+00	CRC89	2.8E-02	3.2E-06	WATERS (U.S. EPA, 2001)			2.2E+04	EPI	6.8E+00	EPI	9.0E-07	EPI	3.4E-01	1.5E+02	3.7E+02	3.7E-02	EPI	
Pentabromodiphenyl ether, 2,2',4,4',5,5'-BDE-99	60348-60-9	5.6E+02	EPI	1.4E-04	3.5E-06	EPI	3.1E-08	EPI	1.3E+00	IRIS	2.2E-02	5.6E-06	WATERS (U.S. EPA, 2001)			2.2E+04	EPI	7.7E+00	EPI	1.1E-02	EPI	3.4E-01	1.5E+02	3.7E+02	3.7E-02	EPI	
Pentachlorobenzene	609-93-5	2.5E+02	EPI	2.9E-02	7.0E-04	EPI	1.0E-03	EPI	1.8E+00	CRC89	2.9E-02	7.9E-06	WATERS (U.S. EPA, 2001)			3.7E+03	EPI	5.2E+00	EPI	6.3E-01	EPI	1.0E+00	2.7E+00	1.0E+01	1.7E-01	EPI	
Pentachloroethane	76-01-7	2.0E+02	EPI	7.9E-02	1.9E-03	EPI	3.5E+00	EPI	1.7E+00	CRC89	3.2E-02	8.6E-06	WATERS (U.S. EPA, 2001)			1.4E+02	EPI	4.8E+02	EPI	3.2E+00	EPI	8.6E-02	1.4E+00	3.4E+00	1.6E-02	EPI	
Pentachloronitrobenzene	82-68-8	3.0E+02	EPI	1.8E-03	4.4E-05	EPI	5.0E-05	EPI	1.7E+00	CRC89	2.6E-02	6.9E-06	WATERS (U.S. EPA, 2001)			6.0E+03	EPI	4.6E+00	EPI	4.4E-01	EPI	2.8E-01	4.7E+00	1.1E+01	4.2E-01	EPI	
Pentachlorophenol	87-86-5	2.7E+02	EPI	1.0E-06	2.5E-08	EPI	1.1E-04	EPI	2.0E+00	CRC89	3.0E-02	8.0E-06	WATERS (U.S. EPA, 2001)			5.0E+03	EPI	5.1E+00	EPI	1.4E+01	EPI	8.0E-01	3.3E+00	1.3E+01	1.3E-01	EPI	
Pentacythritol tetranitrate (PETN)	78-11-5	3.2E+02	EPI	4.9E-10	1.2E-11	EPI	5.5E-09	EPI	1.8E+00	CRC89	2.6E-02	6.8E-06	WATERS (U.S. EPA, 2001)			6.5E+02	EPI	2.4E+00	YAWS	4.3E+01	EPI	6.9E-03	6.2E+00	1.5E+01	1.0E-03	EPI	
Pentane, n-	109-66-0	7.2E+01	EPI	5.1E+01	1.3E+00	EPI	5.1E+02	EPI	6.3E-01	CRC89	8.2E-02	8.8E-06	WATERS (U.S. EPA, 2001)			7.2E+01	EPI	3.4E+00	EPI	3.8E+01	EPI	3.6E-01	2.7E-01	6.4E-01	1.1E-01	EPI	
Perchlorates																											
-Ammonium Perchlorate	7790-98-9	1.2E+02	EPI						2.0E+00	CRC89											2.0E+05	EPI	4.2E-03	4.8E-01	1.1E+00	1.0E-03	RAGSE
-Lithium Perchlorate	7791-03-9	1.1E+02	CRC89						2.4E+00	CRC89											5.9E+05	CRC89	4.0E-03	4.1E-01	1.0E+00	1.0E-03	RAGSE
-Perchlorate and Perchlorate Salts	14797-73-0	1.2E+02	CRC89																		2.5E+05	CRC89	3.8E-03	4.8E-01	1.1E+00	1.0E-03	RAGSE
-Potassium Perchlorate	7778-74-7	1.4E+02	EPI						2.5E+00	CRC89											1.5E+04	EPI	9.1E-03	6.3E-01	1.5E+00	2.0E-03	RAGSE
-Sodium Perchlorate	7601-89-0	1.2E+02	EPI						2.5E+00	CRC89											2.1E+06	EPI	4.3E-03	5.1E-01	1.2E+00	1.0E-03	RAGSE
Perfluorobutane Sulfonate	375-73-5	3.0E+02	EPI	5.9E-04	1.4E-05	EPI	5.2E-02	EPI	1.8E+00	LookChem	2.7E-02	7.2E-06	WATERS (U.S. EPA, 2001)			1.8E+02	EPI	1.8E+00	EPI	8.9E+03	EPI	8.7E-03	5.0E+00	1.2E+01	1.3E-03	EPI	
Permethrin	5245-53-1	3.9E+02	EPI	7.7E-05	1.9E-06	EPI	2.2E-08	EPI	1.2E+00	CRC89	1.9E-02	4.8E-06	WATERS (U.S. EPA, 2001)			1.2E+05	EPI	6.5E+00	EPI	6.7E-03	EPI	1.6E+00	1.6E+01	6.5E+01	2.1E-01	EPI	
Phenacetin	62-44-2	1.8E+02	EPI	8.7E-09	2.1E-10	EPI	6.9E-07	EPI	1.2E+00	CRC89	6.0E-02	8.6E-06	WATERS (U.S. EPA, 2001)			4.1E+01	EPI	1.6E+00	EPI	7.0E+02	EPI	8.9E-03	1.1E+00	2.5E+00	1.7E-03	EPI	
Phenmedipham	13684-63-4	3.0E+02	EPI	3.4E-11	8.4E-13	EPI	1.0E-11	EPI			4.2E-02	5.0E-06	WATERS (U.S. EPA, 2001)			2.6E+03	EPI	3.6E+00	EPI	4.7E+00	EPI	5.2E-02	5.1E+00	1.2E+01	7.9E-03	EPI	
Phenol	108-95-2	9.4E+01	EPI	1.4E-05	3.3E-07	EPI	3.5E-01	EPI	1.1E+00	CRC89	8.3E-02	1.0E-05	WATERS (U.S. EPA, 2001)			1.9E+02	EPI	1.5E+00	EPI	8.3E+04	EPI	1.6E-02	3.9E-01	8.5E-01	4.3E-03	EPI	
Phenothiazine	92-84-2	2.0E+02	EPI	1.1E-06	2.8E-08	EPI	8.3E-07	EPI	1.3E+00	PubChem	2.9E-02	7.5E-06	WATERS (U.S. EPA, 2001)			1.5E+03	EPI	4.2E+00	EPI	1.6E+00	EPI	3.7E-01	1.4E+00	3.3E+00	6.8E-02	EPI	
Phenylenediamine, m-	108-46-2	1.1E+02	EPI	5.1E-08	1.3E-09	EPI	2.1E-03	EPI	1.0E+00	CRC89	7.2E-02	9.2E-06	WATERS (U.S. EPA, 2001)			3.4E+01	EPI	-3.3E-01	EPI	2.4E+05	EPI	9.4E-04	4.2E-01	1.0E+00	2.3E-04	EPI	
Phenylenediamine, o-	95-54-5	1.1E+02	EPI	2.9E-07	7.2E-09	EPI	2.1E-03	EPI	1.3E+00	CRC89	8.4E-02	9.8E-06	WATERS (U.S. EPA, 2001)			3.5E+01	EPI	1.5E-01	EPI	4.0E+04	EPI	1.9E-03	4.2E-01	1.0E+00	4.9E-04	EPI	
Phenylenediamine, p-	106-50-3	1.1E+02	EPI	3.6E-08	8.9E-10	EPI	6.6E-04	EPI	1.4E+00	CRC89	8.4E-02	9.8E-06	WATERS (U.S. EPA, 2001)			3.4E+01	EPI	-3.0E-01	EPI	3.7E+04	EPI	9.8E-04	4.2E-01	1.0E+00	2.5E-04	EPI	
Phenylphenol, 2-	90-43-7	1.7E+02	EPI	4.3E-05	1.1E-06	EPI	2.0E-03	EPI	1.2E+00	CRC89	4.2E-02	7.8E-06	WATERS (U.S. EPA, 2001)			6.7E+03	EPI	3.1E+00	EPI	7.0E+02	EPI	9.8E-02	9.4E-01	2.3E+00	2.0E-02	EPI	
Phorate	298-02-2	2.6E+02	EPI	1.8E-04	4.4E-06	EPI	6.4E-04	EPI	1.2E+00	CRC89	2.3E-02	5.9E-06	WATERS (U.S. EPA, 2001)			4.6E+02	EPI	3.6E+00	EPI	5.0E+01	EPI	7.8E-02	3.0E+00	7.2E+00	1.3E-02	EPI	
Phosgene	75-44-5	9.9E+01	EPI	6.8E-01	1.7E-02	EPI	1.4E+03	EPI	1.4E+00	CRC89	8.9E-02	1.2E-05	WATERS (U.S. EPA, 2001)			1.0E+00	EPI	-7.1E-01	YAWS	6.8E+03	YAWS	5.6E-04	3.8E-01	9.0E-01	1.5E-04	EPI	
Phosmet	732-11-6	3.2E+02	EPI	3.4E-07	8.4E-09	EPI	4.9E-07	EPI			4.1E-02	4.8E-06	WATERS (U.S. EPA, 2001)			1.0E+01	EPI	2.8E+00	EPI	2.4E+01	EPI	1.3E-02	3.6E+00	1.5E+01	1.8E-03	EPI	
Phosphates, Inorganic																											
-Aluminum metaphosphate	13776-88-0	2.6E+02	CRC89						2.8E+00	CRC89													6.2E-03	3.2E+00	7.6E+00	1.0E-03	RAGSE
-Ammonium pyrophosphate	68333-79-9	3.1E+02	EPI																				4.4E-03	5.8E-01	1.4E+00	1.0E-03	RAGSE
-Calcium pyrophosphate	7790-76-3	2.5E+02	CRC89						3.1E+00	CRC89													6.1E-03	2.8E+00	6.7E+00	1.0E-03	RAGSE
-Diammonium phosphate	7783-28-0	1.3E+02	EPI																				4.4E-03	5.8E-01	1.4E+00	1.0E-03	RAGSE
-Dicalcium phosphate	7757-93-9	1.4E+02	EPI																				4.5E-03	6.1E-01	1.5E+00	1.0E-03	RAGSE
-Dimagnesium phosphate	7782-75-4	1.7E+02	CRC89						2.1E+00	CRC89													5.1E-03	1.0E+00	2.4E+00	1.0E-03	RAGSE
-Dipotassium phosphate	7758-11-4	1.7E+02	EPI																				5.1E-03	9.9E-01	2.4E+00	1.0E-03	RAGSE
-Disodium phosphate	7558-79-4	1.4E+02	EPI																				4.6E-03	6.6E-01	1.6E+00	1.0E-03	RAGSE
-Monoaluminum phosphate	13530-50-2	3.2E+02	CRC89																				6.9E-03	6.3E+00	1.5E+01	1.0E-03	RAGSE
-Monoammonium phosphate																											

Contaminant	Molecular Weight		Volatility Parameters				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	Vapor Pressure (g/cm ³)	VP Ref	Density (g/cm ³)	Density Ref	D ₁₀ (cm ² /s)	D _w (cm ² /s)	Dia Ref	K _{ow} (mg/L)	K _{oc} Ref	K _{oc} (L/kg)	K _{oc} Ref	log K _{ow} (L/kg)	log K _{oc} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/vent)	t (hr)	K _p (cm ² /hr)	K _p Ref	KPREF			
-Aroclor 1242	53469-21-9	2.9E+02	EPI	7.8E-03	1.9E-04	EPI	6.6E-05	EPI	1.4E+00	ATSDR Profile	2.4E-02	6.1E-06	WATERS (U.S. EPA, 2001)	7.8E+04	EPI	6.3E+00	EPI	2.8E-01	EPI	2.8E-01	EPI	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	EPI	4.9E-04	EPI	1.4E+00	HSDB	4.9E-04	6.2E-06	WATERS (U.S. EPA, 2001)	7.7E+04	EPI	6.2E+00	EPI	1.0E-01	EPI	1.0E-01	EPI	3.1E+00	4.5E+00	1.9E+01	4.8E-01	EPI					
-Aroclor 1254	11097-69-1	3.3E+02	EPI	1.2E-02	2.8E-04	EPI	6.5E-06	EPI	1.5E+00	ATSDR Profile	2.4E-02	6.1E-06	WATERS (U.S. EPA, 2001)	1.3E+05	EPI	6.5E+00	EPI	4.3E-02	EPI	4.3E-02	EPI	5.2E+00	7.1E+00	3.1E+01	7.5E-01	EPI					
-Aroclor 1260	11096-82-5	4.0E+02	EPI	1.4E-02	3.4E-04	EPI	1.3E-07	EPI	1.6E+00	ATSDR Profile	2.5E-02	5.0E-06	WATERS (U.S. EPA, 2001)	3.5E+05	EPI	7.6E+00	EPI	1.4E-02	EPI	1.4E-02	EPI	7.5E+00	1.7E+01	7.7E+01	9.3E-01	EPI					
-Aroclor 5480	11726-82-4	2.9E+02	EPI	9.1E-03	2.2E-04	EPI	8.5E-06	EPI	1.6E+00	LookChem	2.3E-02	6.8E-06	WATERS (U.S. EPA, 2001)	8.3E+04	EPI	6.3E+00	EPI	3.2E-02	EPI	3.2E-02	EPI	3.8E+00	4.5E+00	2.0E+01	5.3E-01	EPI					
+Hexachlorobiphenyl, 2,3,3',4',4',5'- (PCB 189)	39635-31-9	4.0E+02	EPI	6.6E-03	1.4E-04	EPI	5.5E-07	EPI	1.7E+00	LookChem	2.2E-02	6.7E-06	WATERS (U.S. EPA, 2001)	3.5E+05	EPI	8.3E+00	EPI	7.5E-04	EPI	7.5E-04	EPI	2.3E+01	1.7E+01	8.0E+01	3.0E+00	EPI					
+Hexachlorobiphenyl, 2,3,3',4',4',5'- (PCB 167)	52963-72-6	3.6E+02	EPI	6.6E-03	1.6E-04	EPI	5.8E-07	EPI	1.6E+00	LookChem	2.3E-02	5.9E-06	WATERS (U.S. EPA, 2001)	2.1E+05	EPI	7.5E+00	EPI	2.2E-03	EPI	2.2E-03	EPI	1.0E+01	1.1E+01	5.0E+01	1.4E+00	EPI					
+Hexachlorobiphenyl, 2,3,3',4',4',5'- (PCB 157)	69782-90-7	3.6E+02	EPI	6.6E-03	1.6E-04	EPI	5.8E-07	EPI	1.6E+00	I	2.3E-02	5.9E-06	WATERS (U.S. EPA, 2001)	2.1E+05	EPI	7.6E+00	EPI	1.6E-03	EPI	1.6E-03	EPI	1.2E+01	1.1E+01	5.0E+01	1.7E+00	EPI					
+Hexachlorobiphenyl, 2,3,3',4',4',5'- (PCB 156)	38380-08-4	3.6E+02	EPI	5.8E-03	1.4E-04	EPI	1.6E-06	EPI	1.6E+00	LookChem	2.3E-02	5.9E-06	WATERS (U.S. EPA, 2001)	2.1E+05	EPI	7.6E+00	EPI	5.3E-03	EPI	5.3E-03	EPI	1.2E+01	1.1E+01	5.0E+01	1.7E+00	EPI					
+Hexachlorobiphenyl, 3,3',4',4',5,5'- (PCB 166)	32774-16-6	3.6E+02	EPI	6.6E-03	1.6E-04	EPI	5.8E-07	EPI	1.6E+00	LookChem	2.3E-02	5.9E-06	WATERS (U.S. EPA, 2001)	2.1E+05	EPI	7.4E+00	EPI	5.1E-04	EPI	5.1E-04	EPI	9.1E+00	1.1E+01	5.0E+01	1.2E+00	EPI					
+Pentachlorobiphenyl, 2,3,4,4',5'- (PCB 123)	6510-44-3	3.3E+02	EPI	7.8E-03	1.9E-04	EPI	5.5E-06	EPI	1.5E+00	LookChem	2.4E-02	6.1E-06	WATERS (U.S. EPA, 2001)	1.3E+05	EPI	7.0E+00	EPI	1.6E-02	EPI	1.6E-02	EPI	6.9E+00	7.1E+00	3.2E+01	1.0E+00	EPI					
+Pentachlorobiphenyl, 2,3',4,4',5'- (PCB 118)	31508-00-6	3.3E+02	EPI	1.2E-02	2.9E-04	EPI	9.0E-06	EPI	1.5E+00	LookChem	2.4E-02	6.1E-06	WATERS (U.S. EPA, 2001)	1.3E+05	EPI	7.1E+00	EPI	1.3E-02	EPI	1.3E-02	EPI	8.6E+00	7.1E+00	3.2E+01	1.2E+00	EPI					
+Pentachlorobiphenyl, 2,3,3',4',4'- (PCB 105)	32598-14-4	3.3E+02	EPI	1.2E-02	2.8E-04	EPI	6.5E-06	EPI	1.5E+00	LookChem	2.4E-02	6.1E-06	WATERS (U.S. EPA, 2001)	1.3E+05	EPI	6.8E+00	EPI	3.4E-03	EPI	3.4E-03	EPI	5.2E+00	7.1E+00	3.1E+01	7.5E-01	EPI					
+Pentachlorobiphenyl, 2,3,4,4',5'- (PCB 114)	74472-37-0	3.3E+02	EPI	7.8E-03	1.9E-04	EPI	5.5E-06	EPI	1.5E+00	LookChem	2.4E-02	6.1E-06	WATERS (U.S. EPA, 2001)	1.3E+05	EPI	7.0E+00	EPI	1.6E-02	EPI	1.6E-02	EPI	6.9E+00	7.1E+00	3.2E+01	1.0E+00	EPI					
+Pentachlorobiphenyl, 3,3',4',4',5'- (PCB 126)	57465-28-8	3.3E+02	EPI	7.8E-03	1.9E-04	EPI	2.2E-06	EPI	1.5E+00	LookChem	2.4E-02	6.1E-06	WATERS (U.S. EPA, 2001)	1.3E+05	EPI	7.0E+00	EPI	7.3E-03	EPI	7.3E-03	EPI	6.9E+00	7.1E+00	3.2E+01	1.0E+00	EPI					
+Polychlorinated Biphenyls (high risk)	1336-36-3	2.9E+02	EPI	7.8E-03	1.9E-04	EPI	6.6E-05	EPI	1.4E+00	HSDB	2.4E-02	6.3E-06	WATERS (U.S. EPA, 2001)	7.8E+04	EPI	7.1E+00	EPI	7.0E-01	SSL	7.0E-01	SSL	3.6E+00	4.5E+00	1.9E+01	5.5E-01	EPI					
+Polychlorinated Biphenyls (low risk)	1336-36-3	2.9E+02	EPI	7.8E-03	1.9E-04	EPI	6.6E-05	EPI	1.4E+00	HSDB	2.4E-02	6.3E-06	WATERS (U.S. EPA, 2001)	7.8E+04	EPI	7.1E+00	EPI	7.0E-01	SSL	7.0E-01	SSL	3.6E+00	4.5E+00	1.9E+01	5.5E-01	EPI					
+Polychlorinated Biphenyls (lowest risk)	1336-36-3	2.9E+02	EPI	7.8E-03	1.9E-04	EPI	6.6E-05	EPI	1.4E+00	HSDB	2.4E-02	6.3E-06	WATERS (U.S. EPA, 2001)	7.8E+04	EPI	7.1E+00	EPI	7.0E-01	SSL	7.0E-01	SSL	3.6E+00	4.5E+00	1.9E+01	5.5E-01	EPI					
+Tetrachlorobiphenyl, 3,3',4',4'- (PCB 77)	32598-13-3	2.9E+02	EPI	3.8E-04	9.4E-06	EPI	1.6E-05	EPI	1.6E+00	LookChem	4.3E-02	5.0E-06	WATERS (U.S. EPA, 2001)	6.6E+00	EPI	6.6E+00	EPI	5.7E-04	EPI	5.7E-04	EPI	6.0E+00	4.5E+00	2.0E+01	9.2E-01	EPI					
+Tetrachlorobiphenyl, 3,4,4',5'- (PCB 81)	70362-50-4	2.9E+02	EPI	9.1E-03	2.2E-04	EPI	8.5E-06	EPI	1.4E+00	LookChem	2.4E-02	6.3E-06	WATERS (U.S. EPA, 2001)	7.8E+04	EPI	6.3E+00	EPI	3.2E-02	EPI	3.2E-02	EPI	3.8E+00	4.5E+00	2.0E+01	5.8E-01	EPI					
Polymeric Methylenes Diphenyl Diisocyanate (PMDI)	10760-87-9	5.1E+02	EPI	5.4E-10	1.3E-11	EPI	5.4E-13	EPI	1.0E+00	LookChem	3.0E-02	3.5E-06	WATERS (U.S. EPA, 2001)	1.0E+10	EPI	1.0E+01	EPI	1.8E-06	EPI	1.8E-06	EPI	1.6E+02	7.5E+01	3.7E+02	1.9E+01	EPI					
Polynuclear Aromatic Hydrocarbons (PAHs)																															
-Acenaphthene	83-32-9	1.5E+02	EPI	7.5E-03	1.8E-04	EPI	2.2E-03	EPI	1.2E+00	CRC89	5.1E-02	8.3E-06	WATERS (U.S. EPA, 2001)	5.0E+03	EPI	3.9E+00	EPI	3.9E+00	EPI	3.9E+00	EPI	4.1E-01	7.7E-01	1.8E+00	8.6E-02	EPI					
-Anthracene	120-12-7	1.8E+02	EPI	2.3E-03	5.6E-05	EPI	6.5E-06	EPI	1.3E+00	CRC89	6.5E-06	7.9E-06	WATERS (U.S. EPA, 2001)	1.6E+04	EPI	4.5E+00	EPI	4.3E-02	EPI	4.3E-02	EPI	7.3E-01	1.0E+00	4.1E+00	1.4E-01	EPI					
-Benzo[a]anthracene	56-55-3	2.3E+02	EPI	4.9E-04	1.2E-05	EPI	2.1E-07	EPI	1.3E+00	PubChem	2.6E-02	6.7E-06	WATERS (U.S. EPA, 2001)	1.8E+05	EPI	5.8E+00	EPI	9.4E-03	EPI	9.4E-03	EPI	3.2E+00	2.0E+00	8.5E+00	5.5E-01	EPI					
-Benzo[b]fluoranthene	206-82-3	2.5E+02	EPI	3.8E-06	2.0E-07	EPI	2.6E-08	EPI	1.4E+00	LookChem	4.8E-02	6.6E-06	WATERS (U.S. EPA, 2001)	6.0E+05	EPI	6.1E+00	EPI	2.5E-03	EPI	2.5E-03	EPI	4.2E+00	2.7E+00	1.2E+01	6.9E-01	EPI					
-Benzo[a]pyrene	50-32-8	2.5E+02	EPI	1.9E-05	4.6E-07	EPI	5.5E-09	EPI	1.5E+00	LookChem	4.8E-02	6.6E-06	WATERS (U.S. EPA, 2001)	5.9E+05	EPI	6.1E+00	EPI	1.6E-03	EPI	1.6E-03	EPI	4.4E+00	2.7E+00	1.2E+01	7.1E-01	EPI					
-Benzo[b]fluoranthene	206-99-2	2.5E+02	EPI	2.7E-05	6.6E-07	EPI	5.0E-07	EPI	1.4E+00	LookChem	4.8E-02	6.6E-06	WATERS (U.S. EPA, 2001)	6.0E+05	EPI	5.8E+00	EPI	1.5E-03	EPI	1.5E-03	EPI	2.5E+00	2.7E+00	1.1E+01	4.2E-01	EPI					
-Benzo[k]fluoranthene	207-08-9	2.5E+02	EPI	2.4E-05	5.8E-07	EPI	9.7E-10	EPI	1.4E+00	LookChem	4.8E-02	6.6E-06	WATERS (U.S. EPA, 2001)	5.9E+05	EPI	6.1E+00	EPI	8.0E-04	EPI	8.0E-04	EPI	4.2E+00	2.7E+00	1.2E+01	6.9E-01	EPI					
-Chloronaphthalene, Beta-	91-58-7	1.6E+02	EPI	1.3E-02	3.2E-04	EPI	1.2E-02	EPI	1.1E+00	CRC89	4.5E-02	7.7E-06	WATERS (U.S. EPA, 2001)	2.5E+03	EPI	3.9E+00	EPI	1.2E+01	EPI	1.2E+01	EPI	3.7E-01	8.8E-01	2.1E+00	7.5E-02	EPI					
-Chrysene	21801-9	2.3E+02	EPI	2.1E-04	5.2E-06	EPI	6.2E-09	EPI	1.3E+00	CRC89	2.6E-02	6.7E-06	WATERS (U.S. EPA, 2001)	1.8E+05	EPI	5.8E+00	EPI	2.0E-03	EPI	2.0E-03	EPI	3.5E+00	2.0E+00	8.5E+00	6.0E-01	EPI					
-Dibenz[a,h]anthracene	53-70-3	2.8E+02	EPI	5.8E-06	1.4E-07	EPI	9.6E-10	EPI	1.4E+00	LookChem	4.5E-02	5.2E-06	WATERS (U.S. EPA, 2001)	1.9E+06	EPI	6.8E+00	EPI	2.5E-03	EPI	2.5E-03	EPI	6.1E+00	3.0E+00	1.7E+01	9.5E-01	EPI					
-Dibenz[ghi]perylene	192-65-4	3.0E+02	EPI	5.8E-07	1.4E-08	EPI	7.0E-11	EPI	1.4E+00	LookChem	4.2E-02	4.9E-06	WATERS (U.S. EPA, 2001)	6.5E+06	EPI	7.7E+00	EPI	4.3E-05	EPI	4.3E-05	EPI	2.8E+01	5.2E+00	2.4E+01	4.2E+00	EPI					
+Dimethylbenz[a]anthracene, 7,12	57-97-6	2.6E+02	EPI	1.5E-04	3.8E-06	EPI	2.5E-07	EPI	1.4E+00	LookChem	4.7E-02	5.5E-06	WATERS (U.S. EPA, 2001)	4.9E+05	EPI	5.8E+00	EPI	6.1E-02	EPI	6.1E-02	EPI	2.5E+00	2.9E+00	1.2E+01	4.1E-01	EPI					
-Fluoranthene	206-44-0	2.5E+02	EPI	3.6E-04	8.9E-06	EPI	9.2E-06	EPI	1.3E+00	CRC89	3.6E-02	7.2E-06	WATERS (U.S. EPA, 2001)	5.5E+04	EPI	5.2E+00	EPI	2.6E-01	EPI	2.6E-01	EPI	1.7E+00	1.6E+00	5.7E+00	3.4E-01	EPI	</				

Contaminant		Molecular Weight		Volatility Parameters				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	Vapor Pressure	VP Ref	Density (g/cm ³)	Density Ref	Dia (cm ² /s)	Diw (cm ² /s)	Dia Ref	K _d (L/kg)	K _d Ref	K _{oc} (L/kg)	K _{oc} Ref	log K _{ow} (L/kg)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	t _{event} (hr/event)	t (hr)	K _p (cm/hr)	KPREF
Trinitrobenzene, 1,3,5-	99-35-4	2.1E+02	EPI	2.7E-07	6.5E-09	EPI	6.4E-06	EPI	1.5E+00	CRC89	2.9E-02	7.7E-06	WATER9 (U.S. EPA, 2001)			1.7E+03	EPI	1.2E+00	EPI	2.8E+02	EPI	3.4E-03	1.6E+00	3.9E+00	6.1E-04	EPI
Trinitrotoluene, 2,4,6-	118-96-7	2.3E+02	EPI	8.5E-07	2.1E-08	EPI	8.0E-06	EPI	1.7E+00	CRC89	3.0E-02	7.9E-06	WATER9 (U.S. EPA, 2001)			2.8E+03	EPI	1.6E+00	EPI	1.2E+02	EPI	5.6E-03	2.0E+00	4.7E+00	9.6E-04	EPI
Triphenylphosphine Oxide	791-28-6	2.8E+02	EPI	2.2E-08	5.3E-10	EPI	2.6E-09	EPI	1.2E+00	CRC89	2.3E-02	5.8E-06	WATER9 (U.S. EPA, 2001)			2.0E+03	EPI	2.8E+00	EPI	2.0E+02	EPI	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	EPI	1.1E-07	2.6E-09	EPI	2.9E-07	EPI			3.3E-02	3.9E-06	WATER9 (U.S. EPA, 2001)			1.1E+04	EPI	3.7E+00	EPI	7.0E+00	EPI	1.3E-02	2.7E+01	6.5E+01	1.5E-03	EPI
Tris(1-chloro-2-propyl)phosphate	13674-94-5	3.3E+02	EPI	2.4E-06	6.0E-08	EPI	5.6E-05	EPI			4.0E-02	4.7E-06	WATER9 (U.S. EPA, 2001)			1.6E+03	EPI	2.6E+00	EPI	1.2E+03	EPI	8.4E-03	7.2E+00	1.7E+01	1.2E-03	EPI
Tris(2,3-dibromopropyl)phosphate	126-72-7	7.0E+02	EPI	8.9E-04	2.2E-05	EPI	1.9E-04	EPI	2.3E+00	PubChem	1.9E-02	4.9E-06	WATER9 (U.S. EPA, 2001)			9.7E+03	EPI	4.3E+00	EPI	8.0E+00	EPI	1.4E-03	8.5E+02	2.0E+03	1.4E-04	EPI
Tris(2-chloroethyl)phosphate	115-96-8	2.9E+02	EPI	1.3E-04	3.3E-06	EPI	6.1E-02	EPI	1.4E+00	CRC89	2.4E-02	6.2E-06	WATER9 (U.S. EPA, 2001)			3.9E+02	EPI	1.4E+00	EPI	7.0E+03	EPI	2.3E-03	4.2E+00	1.0E+01	3.6E-04	EPI
Tris(2-ethylhexyl)phosphate	78-42-2	4.3E+02	EPI	3.2E-06	7.9E-08	EPI	8.3E-08	EPI	9.9E-01	CRC89	1.6E-02	3.9E-06	WATER9 (U.S. EPA, 2001)			2.5E+06	EPI	9.5E+00	EPI	6.0E-01	EPI	9.3E+01	2.9E+01	1.3E+02	1.2E+01	EPI
Uranium (Soluble Salts)	NA	2.4E+02	CRC89				0.0E+00	NIOSH	1.9E+01					4.5E+02	BAES							5.9E-03	2.3E+00	5.4E+00	1.0E-03	RAGSE
Urethane	51-79-6	8.9E+01	EPI	2.6E-06	6.4E-08	EPI	2.6E-01	EPI	9.9E-01	CRC89	8.5E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			1.2E+01	EPI	-1.5E-01	EPI	4.8E+06	EPI	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	3.4E+00	CRC89										7.0E+02	CRC89	5.2E-03	1.1E+00	2.6E+00	1.0E-03	RAGSE
Vanadium and Compounds	7440-62-2	5.1E+01	EPI						6.0E+00	CRC89				1.0E+03	SSL							2.7E-03	2.0E-01	4.9E-01	1.0E-03	RAGSE
Vernolate	1929-77-7	2.0E+02	EPI	1.3E-03	3.1E-05	EPI	1.0E-02	EPI	9.5E-01	CRC89	2.4E-02	6.1E-06	WATER9 (U.S. EPA, 2001)			3.0E+02	EPI	3.8E+00	EPI	9.0E+01	EPI	2.2E-01	1.4E+00	3.5E+00	4.0E-02	EPI
Vinclozolin	50471-44-8	2.9E+02	EPI	7.1E-07	1.7E-08	EPI	1.2E-07	EPI	1.5E+00	CRC89	2.5E-02	6.5E-06	WATER9 (U.S. EPA, 2001)			2.8E+02	EPI	3.1E+00	EPI	2.6E+00	EPI	2.9E-02	4.2E+00	1.0E+01	4.5E-03	EPI
Vinyl Acetate	108-05-4	8.6E+01	EPI	2.1E-02	5.1E-04	EPI	9.0E+01	EPI	9.3E-01	CRC89	8.5E-02	1.0E-05	WATER9 (U.S. EPA, 2001)			5.6E+00	EPI	7.3E-01	EPI	2.0E+04	EPI	5.6E-03	3.2E-01	7.7E-01	1.6E-03	EPI
Vinyl Bromide	593-60-2	1.1E+02	EPI	5.0E-01	1.2E-02	EPI	1.0E+03	EPI	1.5E+00	CRC89	8.6E-02	1.2E-05	WATER9 (U.S. EPA, 2001)			2.2E+01	EPI	1.6E+00	EPI	1.0E+04	EPI	1.7E-02	4.2E-01	1.0E+00	4.4E-03	EPI
Vinyl Chloride	75-01-4	6.3E+01	EPI	1.1E+00	2.8E-02	EPI	3.0E+03	EPI	9.1E-01	CRC89	1.1E-01	1.2E-05	WATER9 (U.S. EPA, 2001)			2.2E+01	EPI	1.6E+00	YAWS	8.8E+03	EPI	2.5E-02	2.4E-01	5.7E-01	8.4E-03	EPI
Warfarin	81-81-2	3.1E+02	EPI	1.1E-07	2.8E-09	EPI	1.2E-07	EPI			4.2E-02	4.9E-06	WATER9 (U.S. EPA, 2001)			4.3E+02	EPI	2.7E+00	EPI	1.7E+01	EPI	1.2E-02	5.6E+00	1.3E+01	1.8E-03	EPI
Xylene, P-	106-42-3	1.1E+02	EPI	2.8E-01	6.9E-03	EPI	8.8E+00	EPI	8.6E-01	CRC89	6.8E-02	8.4E-06	WATER9 (U.S. EPA, 2001)			3.8E+02	EPI	3.2E+00	EPI	1.6E+02	EPI	2.0E-01	4.1E-01	9.9E-01	4.9E-02	EPI
Xylene, m-	108-39-3	1.1E+02	EPI	2.9E-01	7.2E-03	EPI	8.3E+00	EPI	8.6E-01	CRC89	6.8E-02	8.4E-06	WATER9 (U.S. EPA, 2001)			3.8E+02	EPI	3.2E+00	EPI	1.6E+02	EPI	2.1E-01	4.1E-01	9.9E-01	5.3E-02	EPI
Xylene, o-	95-47-6	1.1E+02	EPI	2.1E-01	5.2E-03	EPI	6.6E+00	EPI	8.8E-01	CRC89	6.9E-02	8.5E-06	WATER9 (U.S. EPA, 2001)			3.8E+02	EPI	3.1E+00	EPI	1.8E+02	EPI	1.9E-01	4.1E-01	9.9E-01	4.7E-02	EPI
Xylenes	1330-20-7	1.1E+02	EPI	1.1E+02		EPI			8.6E-01	ATSDR Profile												2.0E-01	4.1E-01	9.9E-01	5.0E-02	EPI
Zinc Phosphide	1314-84-7	2.6E+02	CRC89	2.1E-01	5.2E-03	EPI	8.0E+00	EPI	4.6E+00	CRC89	6.9E-02	8.5E-06	WATER9 (U.S. EPA, 2001)			3.8E+02	EPI	3.2E+00	EPI	1.1E+02	EPI	3.7E-03	2.9E+00	7.0E+00	6.0E-04	RAGSE
Zinc and Compound	7440-66-6	6.5E+01	PERRY						7.1E+00	CRC89				6.2E+01	SSL							1.9E-03	2.4E-01	5.9E-01	6.0E-04	RAGSE
Znab	12122-67-7	2.8E+02	EPI	6.5E-09	1.6E-10	EPI	5.7E-09	EPI			4.5E-02	5.2E-06	WATER9 (U.S. EPA, 2001)			1.3E+03	EPI	1.3E+00	EPI	1.0E+01	EPI	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	6.5E+00	CRC89				3.0E+03	BAES							3.7E-03	3.4E-01	8.2E-01	1.0E-03	RAGSE