

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

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 _o (mg/m ³)	k e y ⁻¹	v o l u t e n	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS _d	Analyte	CAS No.	Ingestion SL TR=1E-06 (mg/kg)	Dermal SL TR=1E-06 (mg/kg)	Inhalation SL TR=1E-06 (mg/kg)	Carcinogenic SL TR=1E-06 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
		2.2E-06	I	3.0E-04	O	9.0E-03	I	V		1.07E+05	1.36E+09	8.72E+03	1	0.1	Acephate	30560-19-1				9.9E+01	2.3E+01	9.9E+01		1.9E+01
				2.0E-02	I						1.36E+09		1	0.1	Acetaldehyde	75-07-0			1.1E+01				8.2E+01	8.2E+01
				9.0E-01	I					1.14E+05	1.36E+09	1.37E+04	1		Acetochlor	34256-82-1			1.1E+01				6.6E+03	1.3E+03
						2.0E-03	X				1.36E+09		1	0.1	Acetone	67-64-1							7.0E+04	7.0E+04
						6.0E-02	I	V		1.28E+05	1.36E+09	1.30E+04	1		Acetone Cyanohydrin	75-86-5							2.8E+06	2.8E+06
										2.52E+03	1.36E+09	5.97E+04	1		Acetonitrile	75-05-8							8.1E+02	8.1E+02
3.8E+00	C	1.3E-03	C	1.0E-01	I						1.36E+09		1	0.1	Acetophenone	98-86-2	1.8E-01	6.5E-01	2.9E+03	1.4E-01	7.8E+03			7.8E+03
				5.0E-04	I	2.0E-05	I	V		2.27E+04	1.36E+09	6.91E+03	1		Acetylaminofluorene, 2-	53-96-3								
											1.36E+09		1	0.1	Acroline	107-02-8								
5.0E-01	I	1.0E-04	I	2.0E-03	I	6.0E-03	I		M		1.36E+09		1	0.1	Acrylamide	79-06-1	3.1E-01	1.2E+00	1.4E+04	2.4E-01	1.6E+02	6.6E+02	8.5E+06	1.3E+02
				5.0E-01	I	2.0E-04	P	V		1.09E+05	1.36E+09	9.53E+04	1		Acrylic Acid	79-10-7							2.0E+01	2.0E+01
5.4E-01	I	6.8E-05	I	1.0E-02	A	2.0E-03	I	V		1.13E+04	1.36E+09	7.69E+03	1		Acrylonitrile	107-13-1	1.3E+00		3.2E-01	2.5E-01	7.8E+02		1.6E+01	1.6E+01
						6.0E-03	P				1.36E+09		1	0.1	Adiponitrile	111-69-3							8.5E+06	8.5E+06
5.6E-02	C			1.0E-02	I						1.36E+09		1	0.1	Alachlor	15972-60-8	1.2E+01	4.4E+01		9.7E+00	7.8E+02	3.3E+03		6.3E+02
				1.0E-03	I						1.36E+09		1	0.1	Aldicarb	116-06-3							3.3E+02	6.3E+01
				1.0E-03	I						1.36E+09		1	0.1	Aldicarb Sulfone	1646-88-4							3.3E+02	6.3E+01
											1.36E+09		1	0.1	Aldicarb sulfoxide	1646-87-3								
1.7E+01	I	4.9E-03	I	3.0E-05	I			V			1.36E+09	1.72E+06	1		Aldrin	309-00-2	4.1E-02		9.8E-01	3.9E-02	2.3E+00			2.3E+00
				4.0E-03	P	1.0E-04	X	V		1.11E+05	1.36E+09	3.42E+04	1		Allyl Alcohol	107-18-6							3.6E+00	3.5E+00
2.1E-02	C	6.0E-06	C			1.0E-03	I	V		1.42E+03	1.36E+09	1.58E+03	1		Allyl Chloride	107-05-1	3.3E+01		7.4E-01	7.2E-01			1.7E+00	1.7E+00
				1.0E+00	P	5.0E-03	P				1.36E+09		1		Aluminum	7429-90-5							7.1E+06	7.7E+04
				4.0E-04	I						1.36E+09		1		Aluminum Phosphide	20859-73-8							3.1E+01	3.1E+01
2.1E+01	C	6.0E-03	C	9.0E-03	I						1.36E+09		1	0.1	Ametryn	834-12-8	3.3E-02	1.2E-01	6.4E+02	2.6E-02	7.0E+02	3.0E+03		5.7E+02
											1.36E+09		1	0.1	Aminobiphenyl, 4-	92-67-1								
				8.0E-02	P						1.36E+09		1	0.1	Aminophenol, m-	591-27-5							2.6E+04	5.1E+03
				4.0E-03	X						1.36E+09		1	0.1	Aminophenol, o-	95-55-6							1.3E+03	2.5E+02
				2.0E-02	P						1.36E+09		1	0.1	Aminophenol, p-	123-30-8							6.6E+03	1.3E+03
				2.5E-03	I						1.36E+09		1	0.1	Amtraz	33089-61-1							8.2E+02	1.6E+02
						5.0E-01	I	V			1.36E+09		1		Ammonia	7664-41-7							6.6E+02	1.3E+02
				2.0E-03	X						1.36E+09		1	0.1	Ammonium Picrate	131-74-8							6.6E+02	1.3E+02
				2.0E-01	I						1.36E+09		1		Ammonium Sulfamate	7773-06-0							1.6E+04	1.6E+04
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I			1.37E+04	1.36E+09	2.62E+04	1	0.1	Amyl Alcohol, tert-	75-85-4							8.2E+01	8.2E+01
4.0E-02	P			2.0E-03	X						1.36E+09		1	0.1	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-04	I	3.0E-04	A				1.36E+09	0.15	1	0.1	Anthraquinone, 9,10-	84-65-1	1.7E+01	6.2E+01		1.4E+01	1.6E+02	6.6E+02		1.3E+02
				5.0E-04	H						1.36E+09	0.15	1	0.1	Antimony (metallic)	7440-36-0							4.3E+05	3.1E+01
											1.36E+09		1	0.1	Antimony Pentoxide	1314-60-9							3.9E+01	3.9E+01
				4.0E-04	H						1.36E+09	0.15	1	0.1	Antimony Tetroxide	1332-81-6							2.8E+05	3.1E+01
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C				1.36E+09	0.15	1	0.03	Antimony Trioxide	1309-64-4							2.8E+05	2.8E+05
				3.5E-06	C	5.0E-05	I				1.36E+09		1		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
											1.36E+09		1		Arsine	7784-42-1							7.1E+04	2.7E-01
				3.6E-01	O						1.36E+09		1	0.1	Asbestos (units in fibers)	1332-21-4							1.2E+05	2.3E+04
2.3E-01	C			3.0E-03	A						1.36E+09		1	0.1	Asulam	3337-71-1							9.9E+02	1.9E+02
8.8E-01	C	2.5E-04	C								1.36E+09		1	0.1	Atrazine	1912-24-9	3.0E+00	1.1E+01		2.4E+00	2.3E+02	9.9E+02		
				4.0E-04	I						1.36E+09		1	0.1	Auramine	492-80-8	7.9E-01	2.8E+00	1.5E+04	6.2E-01				
				1.0E-01	A	1.0E-02	A				1.36E+09		1	0.1	Avermectin B1	65195-55-3							1.3E+02	2.5E+01
1.1E-01	I	3.1E-05	I	3.0E-03	A	1.0E-02	A				1.36E+09	5.23E+05	1	0.1	Azinphos-methyl	86-50-0	6.3E+00		4.7E+01	5.6E+00	2.3E+02	9.9E+02	1.4E+07	1.9E+02
				1.0E+00	P	7.0E-06	P				1.36E+09		1	0.1	Azobenzene	103-33-3								
											1.36E+09		1	0.1	Azodicarbonamide	123-77-3								
				2.0E-01	I	5.0E-04	H				1.36E+09	0.07	1		Barium	7440-39-3							7.1E+05	1.5E+04
				5.0E-03	O			V			1.36E+09	3.07E+05	1		Benfluralin	1861-40-1							3.9E+02	3.9E+02
				5.0E-02	I						1.36E+09		1	0.1	Benomyl	17804-35-2							1.6E+04	3.2E+03
4.0E-03	P			2.0E-01	I						1.36E+09		1	0.1	Bensulfuron-methyl	83055-99-6							1.6E+04	1.3E+04
				3.0E-02	I						1.36E+09		1	0.1	Benazone	25057-89-0							2.3E+03	9.9E+03
5.5E-02	I	7.8E-06	I	1.0E-01	I			V		1.16E+03	1.36E+09	2.25E+04	1		Benzaldehyde	100-52-7	1.7E+02			1.7E+02	7.8E+03			7.8E+03
1.0E-01	X			4.0E-03	I	3.0E-02	I	V		1.82E+03	1.36E+09	3.54E+03	1		Benzene	71-43-2	1.3E+01		1.3E+00	1.2E+00	3.1E+02		1.1E+02	8.2E+01
				1.0E-03	P			V		1.26E+03	1.36E+09	1.94E+04	1	0.1	Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1	7.0E+00	2.5E+01		5.4E+00	2.3E+01	9.9E+01		1.9E+01
2.3E+02	I	6.7E-02	I	3.0E-03	I				M		1.36E+09													

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; c = cancer; n = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; m = ceiling limit exceeded; s = Csat exceeded; V = volatile; M = mutagen.														Contaminant			Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1						
Toxicity and Chemical-specific Information														Analyte	CAS No.	Ingestion SL TR=1E-06 (mg/kg)	Dermal SL TR=1E-06 (mg/kg)	Inhalation SL TR=1E-06 (mg/kg)	Carcinogenic SL TR=1E-06 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)				
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	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS											ABS _d			
				3.0E-04	X						3.23E+02	1.36E+09	1.14E+04	1		Bromo-4-fluorobenzene, 1-	460-00-4					2.3E+01				2.3E+01	
				1.7E-03	C						1.36E+09			0.1	Bromoacetic acid	79-08-3							1.3E+02	5.6E+02			1.1E+02
				8.0E-03	I	6.0E-02	I	V			6.79E+02	1.36E+09	8.37E+03	1		Bromobenzene	108-86-1						6.3E+02		5.2E+02	2.9E+02	
						4.0E-02	X	V			4.04E+03	1.36E+09	3.58E+03	1		Bromochloromethane	74-97-5							1.5E+02		1.5E+02	
6.2E-02	I	3.7E-05	C	8.0E-03	P						9.32E+02	1.36E+09	3.97E+03	1		Bromodichloromethane	75-27-4	1.1E+01		3.0E-01	2.9E-01	6.3E+02				6.3E+02	
7.9E-03	I	1.1E-06	I	2.0E-02	I						9.15E+02	1.36E+09	9.70E+03	1		Bromoforn	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			3.59E+03	1.36E+09	1.40E+03	1		Bromomethane	74-83-9						1.1E+02		7.3E+00	6.8E+00	
				5.0E-03	H						1.36E+09	1.24E+05			Bromophos	2104-96-3						3.9E+02				3.9E+02	
1.0E-01	O	3.7E-06	C	1.5E-02	O	1.0E-01	A	V			9.66E+02	1.36E+09	2.14E+03	1		Bromopropane, 1-	106-94-5			1.6E+00	1.6E+00	2.3E+03		4.9E+03	2.2E+02	2.2E+02	
1.0E-01	O			1.5E-02	O						1.36E+09	4.74E+05		0.1	Bromoxynil	1689-94-5	6.7E+00	2.4E+01			5.3E+00	1.2E+03				1.2E+03	
6.0E-01	C	3.0E-05	I			2.0E-03	I	V			6.67E+02	1.36E+09	8.66E+02	1		Bromoxynil Octanoate	1689-99-2	6.7E+00			7.6E+00	1.2E+03				1.2E+03	
											7.64E+03	1.36E+09	3.00E+04	1		Butadiene, 1,3-	106-99-0	1.2E+00		8.1E-02	7.6E-02					1.8E+00	
				1.0E-01	I						1.36E+09				Butanol, n-	71-36-3						7.8E+03				7.8E+03	
5.0E-04	I			4.0E-01	I	5.0E+00	I	V			1.36E+09	2.87E+04		1	Butyl Alcohol, t-	75-65-0	1.4E+03			1.4E+03	3.1E+04			1.5E+05	2.6E+04		
				2.0E+00	P	3.0E+01	P	V			2.13E+04	1.36E+09	2.92E+04	1		Butyl alcohol, sec-	78-92-2				1.6E+05			9.1E+05	1.3E+05		
				5.0E-02	I						1.36E+09	8.63E+04		1	Butylate	2008-41-5					3.9E+03				3.9E+03		
2.0E-04	C	5.7E-08	C	3.0E-01	P						1.36E+09			0.1	Butylated hydroxyanisole	25013-16-5	3.5E+03	1.2E+04	6.7E+07	2.7E+03	2.3E+04	9.9E+04				1.9E+04	
3.6E-03	P			5.0E-02	P						1.36E+09			0.1	Butylated hydroxytoluene	128-37-0	1.9E+02	6.9E+02		1.5E+02	3.9E+03					3.9E+03	
				1.0E-01	X						1.45E+02	1.36E+09	7.35E+03	1		Butylbenzene, sec-	135-98-8				7.8E+03				7.8E+03		
				1.0E-01	X						1.83E+02	1.36E+09	7.36E+03	1		Butylbenzene, tert-	98-06-6				7.8E+03				7.8E+03		
				2.0E-02	A						1.36E+09			0.1	Cacodylic Acid	75-60-5					1.6E+03	6.6E+03				1.3E+03	
1.8E-03	I			1.0E-04	A	1.0E-05	A				1.36E+09			0.025	Cadmium (Diet)	7440-43-9			2.1E+03	2.1E+03	7.8E+00	8.2E+01	1.4E+04			7.1E+00	
1.8E-03	I			1.0E-04	A	1.0E-05	A				1.36E+09			0.05	Cadmium (Water)	7440-43-9											
				5.0E-01	I	2.2E-03	C				1.36E+09			0.1	Caprolactam	105-60-2						3.9E+04	1.6E+05	3.1E+06			3.1E+04
1.5E-01	C	4.3E-05	C	2.0E-03	I						1.36E+09			0.1	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						1.36E+09			0.1	Caplan	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						1.36E+09			0.1	Carbaryl	63-25-2						7.8E+03	3.3E+04				6.3E+03
				5.0E-03	I						1.36E+09			0.1	Carbofuran	1563-66-2						3.9E+02	1.6E+03				3.2E+02
				1.0E-01	I	7.0E-01	I	V			7.38E+02	1.36E+09	1.17E+03	1		Carbon Disulfide	75-15-0				7.8E+03			8.5E+02	7.7E+02		
7.0E-02	I	6.0E-06	I	4.0E-03	I	1.0E-01	I	V			4.58E+02	1.36E+09	1.49E+03	1		Carbon Tetrachloride	56-23-5	9.9E+00		7.0E-01	6.5E-01	3.1E+02			1.6E+02	1.0E+02	
				1.0E-02	I	1.0E-01	P	V			5.89E+03	1.36E+09	6.46E+02	1		Carbonyl Sulfide	463-58-1				3.1E+02			6.7E+01	6.7E+01		
				1.0E-02	I						1.36E+09			0.1	Carbosulfan	55285-14-8					7.8E+02	3.3E+03				6.3E+02	
				1.0E-01	I						1.36E+09			0.1	Carboxin	5234-68-4					7.8E+03	3.3E+04				6.3E+03	
				1.0E-01	I						1.36E+09			0.1	Ceric oxide	1306-38-3								1.3E+06		1.3E+06	
				1.5E-02	I						1.36E+09	1.45E+05		1	Chloral Hydrate	302-17-0					7.8E+03					7.8E+03	
						9.0E-04	I				1.36E+09			0.1	Chloramben	133-90-4					1.2E+03	4.9E+03				9.5E+02	
4.0E-01	H										1.36E+09			0.1	Chloramines, Organic	E701235											
				5.0E-04	G						1.36E+09	1.49E+06		0.04	Chloranil	118-75-2	1.7E+00	6.1E+00		1.3E+00	3.9E+01	4.1E+02				3.6E+01	
				5.0E-04	G						1.36E+09	1.49E+06		0.04	Chlordane (alpha)	5103-71-9					3.9E+01	4.1E+02				3.6E+01	
3.5E-01	I	1.0E-04	I	5.0E-04	I	7.0E-04	I	V			1.36E+09	1.53E+06		0.04	Chlordane (gamma)	5103-74-2					3.9E+01	4.1E+02				3.6E+01	
1.0E+01	I	4.6E-03	C	3.0E-04	I						1.36E+09	1.53E+06		0.04	Chlordane (technical mixture)	12789-03-6	2.0E+00	1.8E+01	4.3E+01	1.7E+00	3.9E+01	4.1E+02	1.1E+03			3.5E+01	
											1.36E+09			0.1	Chlordecone (Kepone)	143-50-0	7.0E-02	2.5E-01	8.3E+02	5.4E-02	2.3E+01	9.9E+01				1.9E+01	
				7.0E-04	A						1.36E+09			0.1	Chlorfenvinphos	470-90-6					5.5E+01	2.3E+02				4.4E+01	
				9.0E-02	O						1.36E+09			0.1	Chlorimuron, Ethyl-	90982-32-4					7.0E+03	3.0E+04				5.7E+03	
				1.0E-01	I	1.5E-04	A	V			2.78E+03	1.36E+09	1.22E+03	1		Chlorine	7782-50-5				7.8E+03			1.8E-01	1.8E-01		
				3.0E-02	I	2.0E-04	I	V			1.36E+09			1	Chlorine Dioxide	10049-04-4					2.3E+03			2.8E+05		2.3E+03	
				3.0E-02	I						1.36E+09			1	Chlorite (Sodium Salt)	7758-19-2					2.3E+03					2.3E+03	
				5.0E+01	I						1.15E+03	1.36E+09	1.03E+03	1		Chloro-1,1-difluoroethane, 1-	75-68-3							5.4E+04	5.4E+04		
4.6E-01	H	3.0E-04	I	2.0E-02	H	2.0E-02	I	V			7.86E+02	1.36E+09	1.08E+03	1		Chloro-1,3-butadiene, 2- (Chloroprene)	126-99-8			1.0E-02	1.0E-02	1.6E+03			2.2E+01		
1.0E-01	P	7.7E-05	C	3.0E-03	X						1.36E+09			0.1	Chloro-2-methylaniline HCl, 4-	3165-93-3	1.5E+00	5.4E+00			1.2E+00					1.9E+02	
2.7E-01	X										1.36E+09			0.1	Chloro-2-methylaniline, 4-	95-89-2	7.0E+00	2.5E+01	5.0E+04	5.4E+00	2.3E+02	9.9E+02				1.9E+02	
				3.5E-03	C						1.36E+09	1.62E+04		1	Chloroacetaldehyde, 2-	107-20-0	2.6E+00										

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

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 (mg/m ³)	k e y	v o l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS _d	Analyte	CAS No.	Ingestion SL TR=1E-06 (mg/kg)	Dermal SL TR=1E-06 (mg/kg)	Inhalation SL TR=1E-06 (mg/kg)	Carcinogenic SL TR=1E-06 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
				1.0E-02	H						1.36E+09		1	0.1	Chlorpyrifos Methyl	5598-13-0					7.8E+02	3.3E+03		6.3E+02
				5.0E-02	O						1.36E+09		1	0.1	Chlorsulfuron	64902-72-3					3.9E+03	1.6E+04		3.2E+03
				1.0E-02	I						1.36E+09		1	0.1	Chlorthal-dimethyl	1861-32-1					7.8E+02	3.3E+03		6.3E+02
				8.0E-04	H						1.36E+09		1	0.1	Chlorthiophos	60238-56-4					6.3E+01	2.6E+02		5.1E+01
				1.5E+00	I						1.36E+09		0.013		Chromium(III), Insoluble Salts	18065-83-1					1.2E+05			1.2E+05
5.0E-01	C	8.4E-02	G	3.0E-03	I	1.0E-04	I		M		1.36E+09		0.025		Chromium(VI)	18540-29-9	3.1E-01	1.6E+01	3.0E-01		2.3E+02		1.4E+05	2.3E+02
				1.3E-02	I						1.36E+09		1	0.1	Chromium, Total	7440-47-3					1.0E+03	4.3E+03		8.2E+02
		9.0E-03	P	3.0E-04	P	6.0E-06	P		V	M	1.36E+09		1		Cobalt	7440-48-4		4.2E+02	4.2E+02		2.3E+01		8.5E+03	2.3E+01
		6.2E-04	I	4.0E-02	H						1.36E+09		1		Coke Oven Emissions	E649830								
				5.0E-02	I	6.0E-01	C				1.36E+09		1	0.1	Copper	7440-50-8					3.1E+03			3.1E+03
				5.0E-02	I	6.0E-01	C				1.36E+09		1	0.1	Cresol, m-	108-39-4					3.9E+03	1.6E+04	8.5E+08	3.2E+03
				2.0E-02	P	6.0E-01	C				1.36E+09		1	0.1	Cresol, o-	95-48-7					1.6E+03	6.6E+03	8.5E+08	1.3E+03
				1.0E-01	A						1.36E+09		1	0.1	Cresol, p-	106-44-5					7.8E+03	3.3E+04		6.3E+03
				1.0E-01	A	6.0E-01	C				1.36E+09		1	0.1	Cresols	1319-77-3					7.8E+03	3.3E+04	8.5E+08	6.3E+03
1.9E+00	H			1.0E-03	P				V	1.66E+04	1.36E+09	1.89E+04	1	0.1	Crotonaldehyde, trans-	123-73-9	3.7E-01		3.7E-01		7.8E+01			7.8E+01
				1.0E-01	I	4.0E-01	I	V		2.68E+02	1.36E+09	6.21E+03	1		Cumene	98-82-8					7.8E+03		2.6E+03	1.9E+03
2.2E-01	C	6.3E-05	C	2.0E-03	H						1.36E+09		1	0.1	Cupferron	135-20-6	3.2E+00	1.1E+01	6.1E+04	2.5E+00	1.6E+02	6.6E+02		1.3E+02
8.4E-01	H			2.0E-03	H						1.36E+09		1	0.1	Cyanazine	21725-46-2	8.3E-01	2.9E+00		6.5E-01				
				1.0E-03	I	9.0E-03	C				1.36E+09		1		Cyanides						7.8E+01		1.3E+07	7.8E+01
				5.0E-03	I						1.36E+09		1		~Calcium Cyanide	592-01-8					3.9E+02			3.9E+02
				6.0E-04	I	8.0E-04	G	V		9.54E+05	1.36E+09	5.33E+04	1		~Cyanide (CN-)	57-12-5				4.7E+01		4.4E+01	2.3E+01	
				1.0E-03	I						1.36E+09		1		~Cyanogen	460-19-5					7.8E+01			7.8E+01
				9.0E-02	I						1.36E+09		1		~Cyanogen Bromide	506-68-3					7.0E+03			7.0E+03
				5.0E-02	I						1.36E+09		1		~Cyanogen Chloride	506-77-4					3.9E+03			3.9E+03
				6.0E-04	I	8.0E-04	I	V		1.00E+07	1.36E+09	5.22E+04	1		~Hydrogen Cyanide	74-90-8				4.7E+01		4.4E+01	2.3E+01	
				2.0E-03	I	9.0E-03	C				1.36E+09		1		~Potassium Cyanide	151-50-8				1.6E+02		1.3E+07		1.6E+02
				5.0E-03	I						1.36E+09		0.04		~Potassium Silver Cyanide	506-61-6				3.9E+02				3.9E+02
				1.0E-01	I						1.36E+09		0.04		~Silver Cyanide	506-64-9				7.8E+03				7.8E+03
				1.0E-03	I	9.0E-03	C				1.36E+09		1		~Sodium Cyanide	143-33-9				7.8E+01		1.3E+07		7.8E+01
				2.0E-04	P						1.36E+09		1		~Thiocyanates	E1790665					1.6E+01			1.6E+01
				2.0E-04	X						1.36E+09		1		~Thiocyanic Acid	463-56-9				1.6E+01				1.6E+01
				5.0E-02	I						1.36E+09		1		~Zinc Cyanide	557-21-1				3.9E+03				3.9E+03
				6.0E+00	I	V				1.17E+02	1.36E+09	1.04E+03	1	0.1	Cyclohexane	110-82-7				3.5E+01	1.2E+02		2.7E+01	1.6E+03
2.0E-02	X			2.0E-02	X						1.36E+09		1	0.1	Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	3.5E+01	1.2E+02			1.6E+03	6.6E+03		1.3E+03
				5.0E+00	I	7.0E-01	P	V		5.11E+03	1.36E+09	4.17E+04	1		Cyclohexanone	108-94-1				2.7E+01		3.0E+04		2.8E+04
				5.0E-03	P	1.0E+00	X	V		2.83E+02	1.36E+09	1.46E+03	1		Cyclohexene	110-83-8				3.9E+02		1.5E+03		3.1E+02
				2.0E-01	I					2.93E+05	1.36E+09	7.46E+04	1		Cyclohexylamine	108-91-8				1.6E+04				1.6E+04
				2.5E-02	I						1.36E+09		1	0.1	Cyfluthrin	68359-37-5					2.0E+03	8.2E+03		1.6E+03
				5.0E-01	O						1.36E+09		1	0.1	Cyromazine	66215-27-8					3.9E+04	1.6E+05		3.2E+04
2.4E-01	I	6.9E-05	C	5.0E-04	A						1.36E+09		1	0.1	DDD, p,p'- (DDD)	72-54-8	2.9E+00	1.0E+01	5.5E+04	2.3E+00	3.9E+01	1.6E+02		3.2E+01
3.4E-01	I	9.7E-05	C	5.0E-04	A				V		1.36E+09	2.10E+06	1		DDE, p,p'-	72-55-9	2.0E+00	2.0E+00	6.1E+01	2.0E+00	3.9E+01			3.9E+01
3.4E-01	I	9.7E-05	I	5.0E-04	I						1.36E+09		1	0.03	DDT	50-29-3	2.0E+00	2.4E+01	3.9E+04	1.9E+00	3.9E+01	5.5E+02		3.7E+01
				3.0E-02	I						1.36E+09		1	0.1	Dalapon	75-99-0					2.3E+03	9.9E+03		1.9E+03
1.8E-02	C	5.1E-06	C	1.5E-01	I						1.36E+09		1	0.1	Daminozide	1698-84-5	3.9E+01	1.4E+02	7.5E+05	3.0E+01	1.2E+04	4.9E+04		9.5E+03
7.0E-04	I			7.0E-03	I						1.36E+09		1	0.1	Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'-(BDE-209)	1163-19-5	9.9E+02	3.5E+03		7.8E+02	5.5E+02	2.3E+03		4.4E+02
				4.0E-05	I						1.36E+09		1	0.1	Demeton	8065-48-3					3.1E+00	1.3E+01		2.5E+00
1.2E-03	I			6.0E-01	I						1.36E+09		1	0.1	Di(2-ethylhexyl)adipate	103-23-1	5.8E+02	2.1E+03		4.5E+02	4.7E+04	2.0E+05		3.8E+04
6.1E-02	H			7.0E-04	A						1.36E+09		1	0.1	Diallate	2303-16-4	1.1E+01	4.1E+01		8.9E+00				
8.0E-01	P	6.0E-03	P	2.0E-04	P	2.0E-04	I	V	M	9.79E+02	1.36E+09	3.20E+04	1		Diazinon	333-41-5					5.5E+01	2.3E+02		4.4E+01
2.5E-01	C			3.0E-04	C						1.36E+09		1	0.1	Dibromo-3-chloropropane, 1,2-	96-12-8	1.9E-01	9.9E+00	5.4E-03	5.3E-03	1.6E+01		6.7E+00	4.7E+00
				4.0E-04	X					1.59E+02	1.36E+09	1.93E+04	1		Dibromoacetic acid	631-64-1	2.8E+00	9.9E+00		2.2E+00	2.3E+01	9.9E+01		1.9E+01
				1.0E-02	I						1.36E+09	2.20E+04	1		Dibromobenzene, 1,3-	108-36-1					3.1E+01			3.1E+01
				2.0E-02	I						1.36E+09	7.95E+03	1		Dibromobenzene, 1,4-	106-37-6					7.8E+02			7.8E+02
8.4E-02	I			2.0E-02	I					8.02E+02	1.36E+09	7.95E+03	1		Dibromochloromethane	124-48-1	8.3E+00			8.3E+00				1.6E+03
2.0E+00	I	6.0E-04	I	9.0E-03	I	9.0E-03	I	V		1.34E+03	1.36E+09	8.64E+03	1		Dibromoethane, 1,2-	106-93-4	3.5E-01		4.0E-02	3.6E-02	7.0E+02		8.1E+01	7.3E+01
				3.0E-04	P	4.0E-03	X	V																

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

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1				
SFO	k _e	IUR	k _e	RfD _o	k _e	RfC	k _e	v	mutagen	C _{sat}	PEF	VF	GIABS	A _{BS_d}	Analyte	CAS No.	Ingestion SL TR=1E-06 (mg/kg)	Dermal SL TR=1E-06 (mg/kg)	Inhalation SL TR=1E-06 (mg/kg)	Carcinogenic SL TR=1E-06 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
3.7E-02	P	3.7E-06	P	1.0E-02 4.0E-02 2.0E-02	I P	4.0E-03	I V			1.36E+09 1.36E+09 1.49E+03	1.36E+09 3.79E+03 6.76E+03	1 1 1		0.05	Dichlorophenoxy Acetic Acid, 2,4-Dichloropropane, 1,2-Dichloropropane, 1,3-Dichloropropanol, 2,3-Dichloropropene, 1,3-Dichlorovos	94-75-7 78-87-5 142-28-9 616-23-9 542-75-6 62-73-7	1.9E+01		2.9E+00	2.5E+00	7.8E+02 3.1E+03 1.6E+03	6.6E+03 1.6E+01	1.6E+01	7.0E+02 1.6E+01 1.6E+03
1.0E-01 2.9E-01	I I	4.0E-06 8.3E-05	I C	3.0E-02 5.0E-04	I I	2.0E-02 5.0E-04	I I			1.57E+03 1.36E+09	1.36E+09 3.55E+03	1 1		0.1 0.1	Dichloropropanol, 2,3-Dichloropropene, 1,3-Dichlorovos	141-86-2 77-73-6 60-57-1	7.0E+00 2.4E+00	8.5E+00	2.5E+00 4.6E+04	1.8E+00 1.9E+00	2.3E+02 3.9E+01	9.9E+02 1.6E+02	7.4E+01 7.1E+05	7.2E+01 3.2E+01
1.6E+01	I	4.6E-03 3.0E-04	I C	3.0E-05 8.0E-02 5.0E-05	O P	3.0E-04 X V	X V			2.56E+02 1.36E+09 1.36E+09	4.11E+03 1.36E+09 1.36E+09	1 1 1		0.1 0.1 0.1	Dicropthos Dicyclopentadiene Diieldrin Diesel Engine Exhaust Diethanolamine	141-86-2 77-73-6 60-57-1 E17136615 111-42-2	4.3E-02	1.5E-01	8.3E+02	3.4E-02	2.3E+00 6.3E+03 3.9E+00	9.9E+00 1.6E+01	1.3E+00	1.9E+00 1.3E+00 3.2E+00
3.5E+02	C	1.0E-01	C	2.0E-03 3.0E-02 6.0E-02 1.0E-03	P P P	2.0E-04 1.0E-04 3.0E-04	P P V			1.12E+05 1.36E+09 1.36E+09	1.39E+05 1.36E+09 1.36E+09	1 1 1		0.1 0.1 0.1	Diethylene Glycol Monobutyl Ether Diethylene Glycol Monoethyl Ether Diethylformamide Diethylstilbestrol	112-34-5 111-90-0 617-84-5 56-53-1	2.0E-03	7.1E-03	3.8E+01	1.6E-03	4.7E+03 4.7E+03 7.8E+01	2.0E+04 4.3E+05	4.3E+05	3.8E+03 3.8E+03 7.8E+01
4.4E-02	C	1.3E-05	C	8.3E-02 2.0E-02	O I	4.0E+01 3.0E+01	X X			1.43E+03 6.91E+02	1.15E+03 7.58E+02	1 1		0.1 0.1	Difenzoquat Diflubenzuron Difluoroethane, 1,1-Difluoropropane, 2,2-Dihydrosafrole	43222-48-6 35367-38-5 75-37-6 420-45-1	1.6E+01		2.7E+01	9.9E+00	6.5E+03 1.6E+03	2.7E+04 6.6E+03	4.8E+04 2.4E+04	5.2E+03 1.3E+03 4.8E+04
1.6E+00 1.7E-03 4.6E+00	P P C	1.4E-01 3.1E-03 1.3E-03	C C C	8.0E-02 2.2E-02 2.2E-03	I O O	7.0E-01 P V	X V			5.30E+02 1.36E+09 1.36E+09	3.81E+04 1.36E+09 1.36E+09	1 1 1		0.1 0.1 0.1	Diisopropyl Ether Diisopropyl Methylphosphonate Dimethipin Dimethoate	108-20-3 1445-75-6 55290-64-7 60-51-5					6.3E+03 1.7E+03 1.7E+02	7.2E+03 7.3E+02	2.2E+03	2.2E+03 1.4E+03 1.4E+02
5.8E-01 2.0E-01 2.7E-02	H P P	2.0E-03 2.0E-03	X X	1.0E-01 1.0E-04	P X	3.0E-02 2.0E-06	I X			1.06E+05 1.72E+05	1.28E+05 2.77E+04	1 1		0.1 0.1	Dimethylbenzidine, 3,3'-Dimethyl methylphosphonate Dimethylamino azobenzene [p-]	119-90-4 756-79-6 60-11-7	4.3E-01 4.1E+02 1.5E-01	1.5E+00 1.5E+03 5.4E-01	2.7E+01 2.9E+03	3.3E-01 3.2E+02 1.2E-01	4.7E+03	2.0E+04	3.8E+03	
1.1E+01	P			1.0E-01 1.0E-04	P X	3.0E-02 2.0E-06	I X			1.06E+05 1.72E+05	1.28E+05 2.77E+04	1 1		0.1 0.1	Dimethylbenzidine, 3,3'-Dimethylformamide Dimethylhydrazine, 1,1-Dimethylhydrazine, 1,2-Dimethylhydrazine, 1,2-Dimethylphenol, 2,4-Dimethylphenol, 3,4-Dimethylvinylchloride	119-93-7 68-12-2 57-14-7 540-73-8 105-67-9 576-26-1 95-65-8	6.3E-02	2.2E-01		4.9E-02	7.8E+03 7.8E+00	4.0E+03 5.8E-02	2.6E+03 5.7E-02	
5.5E+02	C	1.6E-01	C	2.0E-02 6.0E-04	I I					1.89E+05 1.36E+09 1.36E+09	1.68E+05 1.36E+09 1.36E+09	1 1 1		0.1 0.1 0.1	Dimethylhydrazine, 1,2-Dimethylphenol, 2,4-Dimethylphenol, 3,4-Dimethylvinylchloride	540-73-8 105-67-9 576-26-1 95-65-8	1.3E-03	2.9E-03	8.8E-04	1.6E+03 4.7E+01	6.6E+03 2.0E+02	3.3E+02	1.3E+03 3.8E+01	
4.5E-02	C	1.3E-05	C	1.0E-03 8.0E-05	I X					4.73E+02 1.36E+09	5.48E+03 1.36E+09	1 1		0.1 0.1	Dinitro- <i>o</i> -cresol, 4,6-Dinitro- <i>o</i> -cyclohexyl Phenol, 4,6-Dinitroaniline, 3,5-Dinitrobenzene, 1,2-Dinitrobenzene, 1,3-Dinitrobenzene, 1,4-Dinitrophenol, 2,4-Dinitrotoluene Mixture, 2,4,2,6-Dinitrotoluene, 2,4-Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade	534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5	1.5E+01	1.2E+00	1.1E+00	6.3E+00 1.6E+02	2.6E+01 6.6E+02	5.1E+00		
6.8E-01 3.1E-01 1.5E+00	I C P	8.9E-05 3.8E-01	C C	1.0E-03 1.0E-04 1.0E-04	X X P	2.0E-03 3.0E-04	I X X			1.36E+09 1.36E+09 1.36E+09	1.36E+09 1.012 0.099	1 1 1		0.1 0.1 0.1	Dinitrobenzene, 1,3-Dinitrobenzene, 1,4-Dinitrophenol, 2,4-Dinitrotoluene Mixture, 2,4,2,6-Dinitrotoluene, 2,4-Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade	E1615210 121-14-2 606-20-2 35572-78-2 19406-51-0 25321-14-6	1.0E+00 2.2E+00 4.6E-01	3.6E+00 7.8E+00 1.7E+00	4.3E+04 3.6E-01	8.0E-01 1.7E+00 3.6E-01	1.6E+02 2.3E+01	6.5E+02 1.0E+02	1.3E+02 1.9E+01	
4.5E-01	X			1.0E-03 1.0E-04	X X					1.36E+09 1.36E+09	1.36E+09 1.006 0.009	1 1 1		0.1 0.1 0.1	Dinitrobenzene, 1,3-Dinitrobenzene, 1,4-Dinitrophenol, 2,4-Dinitrotoluene Mixture, 2,4,2,6-Dinitrotoluene, 2,4-Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade	88-85-7 123-91-1	1.5E+00	5.5E+00		1.2E+00	7.8E+00 7.0E+01	3.3E+01 3.0E+02	6.3E+00 5.7E+01	
1.0E-01	I	5.0E-06	I	1.0E-03 3.0E-02	I I	3.0E-02 I V	I V			1.16E+05 1.36E+09	3.96E+04 1.36E+09	1 1		0.1	Dinoseb Dioxane, 1,4-Dioxins	88-85-7 123-91-1	7.0E+00		2.2E+01	5.3E+00	7.8E+01 2.3E+03	3.3E+02 1.2E+03	6.3E+01 8.1E+02	
6.2E+03 1.3E+05	I C	1.3E+00 3.8E+01	I C	7.0E-10 3.0E-02	I I	4.0E-08 C V	C V			1.36E+09 1.36E+09 1.36E+09	1.96E+06 1.36E+09 1.36E+09	1 1 1		0.03 0.03 0.1	~Hexachlorodibenzo-p-dioxin, Mixture ~TCDD, 2,3,7,8-Diphenamid	34465-46-8 1746-01-6 957-51-7	1.1E-04 5.3E-06	1.3E-03 6.3E-05	2.9E+00 1.4E-04	1.0E-04 4.8E-06	5.5E-05 2.3E+03	7.7E-04 9.9E+03	8.2E-02 3.4E+01	5.1E-05 1.9E+03
8.0E-01	I	2.2E-04	I	8.0E-04 1.0E-01	X O	4.0E-04 X V	X V			1.36E+09 1.36E+09 1.36E+09	8.06E+04 1.36E+09 1.36E+09	1 1 1		0.1 0.1 0.1	Diphenyl Ether Diphenyl Sulfone Diphenylamine	101-84-8 127-63-9 122-39-4	8.7E-01	3.1E+00	1.7E+04	6.8E-01	6.3E+01 7.8E+03	2.6E+02 3.3E+04	3.4E+01 6.3E+03	
7.4E+00 7.4E+00 6.7E+00	C C C	2.1E-03 2.1E-03 1.9E-03	C C C	2.2E-03 4.0E-05	I I					1.36E+09 1.36E+09 1.36E+09	1.36E+09 1.36E+09 1.36E+09	1 1 1		0.1 0.1 0.1	Diphenylhydrazine, 1,2-Diquat Direct Black 38 Direct Blue 6 Direct Brown 95 Disulfoton	122-66-7 2764-72-9 1937-57-7 2602-46-2 16071-86-6 298-04-4	9.4E-02 9.4E-02 1.0E-01	3.3E-01 3.3E-01	1.8E+03 2.0E+03	7.3E-02 7.3E-02 8.1E-02	1.7E+02 3.1E+01	7.3E+02	1.4E+02 2.5E+00	
9.9E-03	I	1.2E-06	I	2.0E-02 3.0E-04	I I	1.0E-03 P I V	I V			1.05E+04 1.36E+09	1.89E+04 7.66E+03	1 1		0.1 0.1	Endothall Endrin Epichlorohydrin	145-73-3 72-20-8 106-89-8	7.0E+01	4.4E+01	2.7E+01	3.9E+03 4.7E+02 4.7E+02	2.0E+03 6.6E+03	1.3E+03 3.8E+02	1.3E+03 1.9E+01 1.9E+01	
				4.0E-02	P	2.0E-02	I V			1.53E+04 1.36E+09	7.66E+03 1.36E+09	1 1		0.1	Epoxbutane, 1,2-Ethanol, 2-(2-methoxyethoxy)-	106-88-7 111-77-3					3.1E+03 1.3E+04	1.6E+02	1.6E+02 2.5E+03	

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; c = cancer; n = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; m = ceiling limit exceeded; s = Csat exceeded; V = volatile; M = mutagen.															Contaminant				Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1			
Toxicity and Chemical-specific Information															Analyte		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 _o (mg/m ³)	k _e y ⁻¹	o	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS _d	CAS No.	Ingestion SL TR=1E-06 (mg/kg)	Dermal SL TR=1E-06 (mg/kg)	Inhalation SL TR=1E-06 (mg/kg)	Carcinogenic SL TR=1E-06 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THQ=1 (mg/kg)			
				5.0E-03	I						1.36E+09		1	0.1	Ethephon	16672-87-0					3.9E+02	1.6E+03		3.2E+02		
				5.0E-04	I						1.36E+09		1	0.1	Ethion	563-12-2					3.9E+01	1.6E+02		3.2E+01		
				1.0E-01	P	6.0E-02	P	V		2.38E+04	1.36E+09	6.15E+04	1		Ethoxvethanol Acetate, 2-	111-15-9					7.8E+03		3.8E+03	2.6E+03		
				9.0E-02	P	4.0E-02	P	V		1.06E+05	1.36E+09	9.84E+04	1		Ethoxvethanol, 2-	110-80-5					7.0E+03		4.1E+03	2.6E+03		
				7.0E-01	P	7.0E-02	P	V		1.08E+04	1.36E+09	8.62E+03	1		Ethyl Acetate	141-78-6					5.5E+04		6.3E+02	6.2E+02		
				5.0E-03	P	8.0E-03	P	V		2.50E+03	1.36E+09	6.34E+03	1		Ethyl Acrylate	140-88-5					3.9E+02		5.3E+01	4.7E+01		
										4.0E+00	P	V			Ethyl Chloride (Chloroethane)	75-00-3							5.4E+03			
				2.0E-01	I					1.01E+04	1.36E+09	3.12E+03	1		Ethyl Ether	60-29-7					1.6E+04			1.6E+04		
										1.10E+03	1.36E+09	5.77E+03	1		Ethyl Methacrylate	97-63-2					7.8E+04		1.8E+03	1.8E+03		
				8.0E-08	I	1.0E+00	I	4.0E+01	I	V	2.87E+03	1.36E+09	3.67E+03	1		Ethyl Tertiary Butyl Ether (ETBE)	637-92-3	1.3E+02	1.3E+02			1.5E+05	5.2E+04			
				1.1E-02	C	2.5E-06	C			1.0E+00	I	V		4.80E+02	Ethyl-p-nitrophenyl Phosphonate	2104-64-5	6.3E+01		6.4E+00	5.8E+00	7.8E-01	3.3E+00		6.3E-01		
				5.0E-02	P	1.0E+00	I	V		1.36E+09			1	0.1	Ethylbenzene	100-41-4					3.9E+03			2.4E+03		
				7.0E-02	P					1.36E+09			1	0.1	Ethylene Cyanohydrin	109-78-4					5.5E+03	2.3E+04	5.9E+03	4.4E+03		
				9.0E-02	P					1.89E+05	1.36E+09	1.80E+05	1		Ethylene Diamine	107-15-3					7.0E+03			7.0E+03		
				8.0E-01	A	4.0E-01	C			1.36E+09			1	0.1	Ethylene Glycol	107-21-1					6.3E+04	2.6E+05	5.7E+08	5.1E+04		
				1.0E-01	I	1.6E+00	I			1.36E+09			1	0.1	Ethylene Glycol Monobutyl Ether	111-76-2					7.8E+03	3.3E+04	2.3E+09	6.3E+03		
				3.1E-01	C	3.0E-03	I			3.0E-02	C	V	M	1.21E+05	Ethylene Oxide	75-21-8	4.9E-01		2.1E-03	2.0E-03	6.3E+00	2.6E+01	1.9E+02	1.9E+02		
				4.5E-02	C	1.3E-05	C			8.0E-05	I			1.36E+09	Ethylene Thiourea	96-45-7	1.5E+01	5.5E+01	2.9E+05	1.2E+01				5.1E+00		
				6.5E+01	C	1.9E-02	C				V			1.54E+05	Ethyleneimine	151-56-4	1.1E-02		3.5E-03	2.7E-03						
				3.0E+00	I					1.36E+09			1	0.1	Ethylphthalyl Ethyl Glycolate	84-72-0					2.3E+05	9.9E+05		1.9E+05		
				2.5E-04	I					1.36E+09			1	0.1	Fenamiphos	22224-92-6					2.0E+01	8.2E+01		1.6E+01		
				2.5E-02	I					1.36E+09			1	0.1	Fenpropathrin	39515-41-8					2.0E+03	8.2E+03		1.6E+03		
				2.5E-02	I					1.36E+09			1	0.1	Fenvalerate	51630-58-1					2.0E+03	8.2E+03		1.6E+03		
				1.3E-02	I					1.36E+09			1	0.1	Fluometuron	2164-17-2					1.0E+03	4.3E+03		8.2E+02		
				4.0E-02	C	1.3E-02	C			1.36E+09			1		Fluoride	16984-48-8					3.1E+03		1.8E+07	3.1E+03		
				6.0E-02	I	1.3E-02	C			1.36E+09			1		Fluorine (Soluble Fluoride)	7782-41-4					4.7E+03		1.8E+07	4.7E+03		
				8.0E-02	I					1.36E+09			1	0.1	Fluridone	59756-60-4					6.3E+03	2.6E+04		5.1E+03		
				4.0E-02	O					1.36E+09			1	0.1	Flurprimidol	56425-91-3					3.1E+03	1.3E+04		2.5E+03		
				2.0E-03	O					1.36E+09			1	0.1	Flusilazole	85509-19-9					1.6E+02	6.6E+02		1.3E+02		
				5.0E-01	O					1.36E+09			1	0.1	Flutolanil	66332-96-5					3.9E+04	1.6E+05		3.2E+04		
				1.0E-02	I					1.36E+09			1	0.1	Fluralinate	69409-94-5					7.8E+02	3.3E+03		6.3E+02		
				9.0E-02	O					1.36E+09			1	0.1	Folpet	133-07-3					7.0E+03	3.0E+04		5.7E+03		
				1.0E-02	O					1.36E+09			1	0.1	Fomesafen	72178-02-0					7.8E+02	3.3E+03		6.3E+02		
				2.0E-03	I					1.36E+09			1	0.1	Fonofos	944-22-9					1.6E+02	6.6E+02		1.3E+02		
				2.1E-02	C	1.3E-05	I			2.0E-01	I	9.8E-03	A	V	4.24E+04	Formaldehyde	50-00-0	3.3E+01		1.7E+01	1.1E+01	1.6E+04		8.0E+02	7.6E+02	
				9.0E-01	P	3.0E-04	X	V		1.06E+05	1.36E+09	9.30E+04	1		Formic Acid	64-18-6					7.0E+04		2.9E+01	2.9E+01		
				2.5E+00	O					1.36E+09			1	0.1	Fosetyl-AL	39148-24-8					2.0E+05	8.2E+05		1.6E+05		
				1.0E-03	X					V			1		Furans											
				1.0E-03	I					V			1		~Dibenzofuran	132-64-9					7.8E+01			7.8E+01		
				1.0E-03	I					6.22E+03	1.36E+09	2.62E+03	1		~Furan	110-00-9					7.8E+01			7.8E+01		
				3.8E+00	H					9.0E-01	I	2.0E+00	I	V	1.65E+05	~Tetrahydrofuran	109-99-9					7.0E+04	2.5E+04		1.8E+04	
										1.36E+09			1	0.1	Furazolidone	67-45-8	1.8E-01	6.5E-01								
				3.0E-03	I	5.0E-02	H	V		1.01E+04	1.36E+09	4.86E+04	1		Furfural	98-01-1					2.3E+02		2.5E+03	2.1E+02		
				1.5E+00	C	4.3E-04	C				1.36E+09		1	0.1	Furium	531-82-8	4.6E-01	1.6E+00	8.9E+03	3.6E-01						
				3.0E-02	I	8.6E-06	C				1.36E+09		1	0.1	Furmecycloz	60568-05-0	2.3E+01	8.2E+01	4.4E+05	1.8E+01						
				6.0E-03	O					1.36E+09			1	0.1	Glufosinate, Ammonium	77182-82-2					4.7E+02	2.0E+03		3.8E+02		
				1.0E-01	A	8.0E-05	C			1.36E+09			1	0.1	Glutaraldehyde	111-30-8					7.8E+03	3.3E+04	1.1E+05	6.0E+03		
				4.0E-04	I	1.0E-03	X	V		1.06E+05	1.36E+09	8.43E+04	1		Glycidaldehyde	765-34-4					3.1E+01		8.8E+01	2.3E+01		
				1.0E-01	I					1.36E+09			1	0.1	Glyphosate	1071-83-6					7.8E+03	3.3E+04		6.3E+03		
				1.0E-02	X					1.36E+09	1.45E+05		1		Guanidine	113-00-8					7.8E+02			7.8E+02		
				2.0E-02	P					1.36E+09			1	0.1	Guanidine Chloride	50-01-1					1.6E+03	6.6E+03		1.3E+03		
				3.0E-02	X					1.36E+09			1	0.1	Guanidine Nitrate	506-93-4					2.3E+03	9.9E+03		1.9E+03		
				4.5E+00	I	1.3E-03	I			5.0E-05	I			1	0.1	Haloxypol, Methyl	69806-40-2					3.9E+00	1.6E+01		3.2E+00	
				9.1E+00	I	2.6E-03	I	1.3E-05	I	1.0E-04	A		V	1.36E+09	Heptachlor	76-44-8	1.5E-01		1.0E+00	1.3E-01	7.8E+00			7.8E+00		
										1.36E+09	8.43E+05		1		Heptachlor Epoxide	1024-57-3	7.6E-02		9.1E-01	7.0E-02	1.0E+00			1.0E+00		
				3.0E-04	X	4.0E-01	P	V		2.09E+02	1.36E+09	7.80E+03	1		Heptanal, n-	111-71-7					2.3E+01		2.4E+01	2.4E+01		
				2.0E-03	I					5.79E+01	1.36E+09	8.95E+02	1		Heptane, N-	142-82-5					2.3E+01		3.7E+02	2.2E+01		
				1.6E+00	I	4.6E-04	I	1.0E-05	P		1.36E+09	3.80E+05	1	0.1	Hexabromobenzene	87-82-1					1.6E+02			1.6E+02		
				7.8E-02	I	2.2E-05	I	1.0E-03	P		1.36E+09	1.08E+04	1		Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BDE-153)	68631-49-2	4.3E-01		4.1E-01	2.1E-01	1.6E+01	6.6E+01		1.3E+01		
				1.8E+00	I	5.3E-04	I			1.36E+09	6.80E+04		1		Hexachlorobutadiene	118-74-1	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		1.36E+09		1	0.1	Hexachlorocyclohexane, Alpha-	319-84-6	1.1E-01	3.9E-01	2.1E+03	8.6E-02						
										1.36E+09			1	0.1	Hexachlorocyclohexane, Beta-	319-85-7					3.9E-01	1.4E+00	7.2E+03	3.0E-01		
										1.36E+09			1	0.1	Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	6.3E-01	5.6E+00	1.2E+04	5.7E-01						
				1.8E+00	I	5.1E-04	I			1.36E+09			1	0.1	Hexachlorocyclohexane, Technical	608-7										

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

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 _o (mg/m ³)	k e y	v o l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS		ABS _d	Analyte	CAS No.	Ingestion SL TR=1E-06 (mg/kg)	Dermal SL TR=1E-06 (mg/kg)	Inhalation SL TR=1E-06 (mg/kg)	Carcinogenic SL TR=1E-06 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)
				5.0E-03	I	3.0E-02	I	V		3.28E+03	1.36E+09	1.33E+04	1		Hexanone, 2-Hexazinone	591-78-6 51235-04-2					3.9E+02		4.2E+02	2.0E+02
				3.3E-02	I						1.36E+09		1	0.1							2.6E+03	1.1E+04		2.1E+03
				2.5E-02	I						1.36E+09		1	0.1	Hexythiazox	78587-05-0					2.0E+03	2.8E+03		1.6E+03
				1.7E-02	O						1.36E+09		1	0.1	Hydramethylnon	67485-29-4					1.3E+03	5.6E+03		1.1E+03
3.0E+00	I	4.9E-03	I			3.0E-05	P	V		1.12E+05	1.36E+09	6.52E+04	1		Hydrazine	302-01-2	2.3E-01		3.7E-02	3.2E-02			2.0E+00	2.0E+00
3.0E+00	I	4.9E-03	I								1.36E+09		1		Hydrazine Sulfate	10034-93-2	2.3E-01		7.8E+02	2.3E-01			2.8E+07	2.8E+07
				4.0E-02	C	1.4E-02	C	V			1.36E+09		1		Hydrogen Chloride	7647-01-0					3.1E+03		2.0E+07	3.1E+03
				2.0E-02	I	2.0E-03	I	V			1.36E+09		1		Hydrogen Fluoride	7664-39-3							2.8E+06	2.8E+06
6.0E-02	P			4.0E-02	P						1.36E+09		1	0.1	Hydrogen Sulfide	7783-06-4	1.2E+01	4.1E+01		9.0E+00	3.1E+03	1.3E+04		2.8E+06
6.1E-02	O			1.1E-01	O						1.36E+09		1	0.1	Hydroquinone	123-31-9	1.1E+01	4.0E+01		8.9E+00	8.4E+03	3.6E+04		2.5E+03
				2.5E-01	I						1.36E+09		1	0.1	Imazalil	35554-44-0					2.0E+04	8.2E+04		1.6E+04
				2.5E+00	I						1.36E+09		1	0.1	Imazequin	81335-37-7					2.0E+05	8.2E+05		1.6E+05
				1.0E-02	A						1.36E+09		1		Imazethapyr	81335-77-5					7.8E+02			7.8E+02
				4.0E-02	I						1.36E+09		1	0.1	Iprodione	36734-19-7					3.1E+03	1.3E+04		2.5E+03
				7.0E-01	P						1.36E+09		1		Iron	7439-89-6					5.5E+04			5.5E+04
				3.0E-01	I	4.0E-01	X	V		1.00E+04	1.36E+09	2.81E+04	1		Isobutyl Alcohol	78-83-1					2.3E+04		1.2E+04	7.8E+03
9.5E-04	I			2.0E-01	I	2.0E+00	C				1.36E+09		1	0.1	Isophorone	78-59-1	7.3E+02	2.6E+03		5.7E+02	1.6E+04	6.6E+04	2.8E+09	1.3E+04
				1.5E-02	I			V			1.36E+09	4.20E+05	1		Isopropalin	33820-53-0					1.2E+03			1.2E+03
				2.0E+00	P	2.0E-01	P	V		1.09E+05	1.36E+09	2.77E+04	1		Isopropanol	67-63-0					1.6E+05		5.8E+03	5.6E+03
				1.0E-01	I						1.36E+09		1	0.1	Isopropyl Methyl Phosphonic Acid	1832-54-8					7.8E+03	3.3E+04		6.3E+03
				5.0E-02	I						1.36E+09		1	0.1	Isoxaben	82558-50-7					3.9E+03	1.6E+04		3.2E+03
				3.0E-01	A	V					1.36E+09		1		JP-7	E1737665							4.3E+08	4.3E+08
				8.0E-03	O						1.36E+09		1	0.1	Lactofen	77501-63-4					6.3E+02	2.6E+03		5.1E+02
				2.0E-04	X						1.36E+09		1	0.1	Lactonitrile	78-97-7					1.6E+01	6.6E+01		1.3E+01
				5.0E-05	P						1.36E+09		1		Lanthanum	7439-91-0					3.9E+00			3.9E+00
				2.1E-05	P						1.36E+09		1	0.1	Lanthanum Acetate Hydrate	100587-90-4					1.6E+00	6.9E+00		1.3E+00
				1.9E-05	P						1.36E+09		1		Lanthanum Chloride Heptahydrate	10025-84-0					1.5E+00			1.5E+00
				2.8E-05	P						1.36E+09		1		Lanthanum Chloride, Anhydrous	10099-58-8					2.2E+00			2.2E+00
				1.6E-05	P						1.36E+09		1		Lanthanum Nitrate Hexahydrate	10277-43-7					1.3E+00			1.3E+00
8.5E-03	C	1.2E-05	C								1.36E+09		1		Lead Compounds									
2.1E-01	C	8.0E-05	C								1.36E+09		1	0.1	~Lead Phosphate	7446-27-7	8.2E+01		3.2E+05	8.2E+01				
											1.36E+09		1		~Lead acetate	301-04-2	3.3E+00	1.2E+01	4.8E+04	2.6E+00				
3.8E-02	C	1.1E-05	C								1.36E+09		1	0.1	~Lead and Compounds	7439-92-1								4.0E+02
											1.36E+09		1	0.1	~Lead subacetate	1335-32-6	1.8E+01	6.5E+01	3.5E+05	1.4E+01				
				1.0E-07	I			V		2.43E+00	1.36E+09	1.91E+03	1		~Tetraethyl Lead	78-00-2					7.8E-03			7.8E-03
				5.0E-06	P			V		3.83E+02	1.36E+09	2.55E+04	1		Lewisite	541-25-3					3.9E-01			3.9E-01
				7.7E-03	O						1.36E+09		1	0.1	Linuron	330-55-2					6.0E+02	2.5E+03		4.9E+02
				2.0E-03	P						1.36E+09		1		Lithium	7439-93-2					1.6E+02			1.6E+02
				5.0E-04	I						1.36E+09		1	0.1	MCPA	94-74-6					3.9E+01	1.6E+02		3.2E+01
				4.4E-02	O						1.36E+09		1	0.1	MCPB	94-81-5					3.4E+03	1.5E+04		2.8E+03
				1.0E-03	I						1.36E+09		1	0.1	MCPP	93-65-2					7.8E+01	3.3E+02		6.3E+01
				2.0E-02	I						1.36E+09		1	0.1	Malathion	121-75-5					1.6E+03	6.6E+03		1.3E+03
				1.0E-01	I	7.0E-04	C				1.36E+09		1	0.1	Maleic Anhydride	108-31-6					7.8E+03	3.3E+04	9.9E+05	6.3E+03
				5.0E-01	I						1.36E+09		1	0.1	Maleic Hydrazide	123-33-1					3.9E+04	1.6E+05		3.2E+04
				1.0E-04	P						1.36E+09		1	0.1	Malononitrile	109-77-3					7.8E+00	3.3E+01		6.3E+00
				3.0E-02	H						1.36E+09		1	0.1	Mancozeb	8018-01-7					2.3E+03	9.8E+03		1.9E+03
				5.0E-03	I						1.36E+09		1	0.1	Maneb	12427-38-2					3.9E+02	1.6E+03		3.2E+02
				1.4E-01	I	5.0E-05	I				1.36E+09		1		Manganese (Diet)	7439-96-5								
				2.4E-02	G	5.0E-05	I				1.36E+09	0.04			Manganese (Non-diet)	7439-96-5					1.9E+03		7.1E+04	1.8E+03
				9.0E-05	H						1.36E+09		1	0.1	Mephsfolan	950-10-7					7.0E+00	3.0E+01		5.7E+00
				3.0E-02	I						1.36E+09		1	0.1	Mepiquat Chloride	24307-26-4					2.3E+03	9.9E+03		1.9E+03
1.1E-02	P			4.0E-03	P						1.36E+09		1	0.1	Mercaptobenzothiazole, 2-Mercuro Compounds	149-30-4	6.3E+01	2.2E+02		4.9E+01	3.1E+02	1.3E+03		2.5E+02
				3.0E-04	I	3.0E-04	G				1.36E+09	0.07			~Mercuric Chloride (and other Mercury salts)	7487-94-7					2.3E+01		4.3E+05	2.3E+01
				3.0E-04	I	V				3.13E+00	1.36E+09	3.47E+04	1		~Mercury (elemental)	7439-97-6						1.1E+01		1.1E+01
				1.0E-04	I						1.36E+09		1		~Methyl Mercury	22967-92-6					7.8E+00			7.8E+00
				8.0E-05	I						1.36E+09		1	0.1	~Phenylmercuric Acetate	62-38-4					6.3E+00	2.6E+01		5.1E+00
				3.0E-05	I						1.36E+09	1.94E+06	1		Merphos	150-50-5					2.3E+00			2.3E+00
				6.0E-02	I						1.36E+09		1	0.1	Metalaxyl	57837-19-1					4.7E+03	2.0E+04		3.9E+03
				1.0E-04	I	3.0E-02	P	V																

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 (ug/m ³) ⁻¹	IUR (ug/m ³) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _o (mg/m ³)	k _e y	v _o l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS _d	Analyte	CAS No.	Ingestion SL TR=1E-06 (mg/kg)	Dermal SL TR=1E-06 (mg/kg)	Inhalation SL TR=1E-06 (mg/kg)	Carcinogenic SL TR=1E-06 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THQ=1 (mg/kg)		
7.8E-03	O			1.9E-01	O						1.36E+09			1	0.1	Oryzalin	19044-88-3	8.9E+01	3.2E+02		7.0E+01	1.5E+04	6.3E+04		1.2E+04	
				5.0E-03	I						1.36E+09			1	0.1	Oxadiazon	19666-30-9				3.9E+02	1.6E+03			3.2E+02	
7.3E-02	O			2.5E-02	I						1.36E+09			1	0.1	Oxamyl	23135-22-0				2.0E+03	8.2E+03			1.6E+03	
				4.0E-02	O						1.36E+09			1	0.1	Oxyfluorfen	42874-03-3	9.5E+00	3.4E+01		7.4E+00	3.1E+03	1.3E+04		2.5E+03	
				1.3E-02	I						1.36E+09			1	0.1	Paclobutrazol	76738-62-0				1.0E+03	4.3E+03			8.2E+02	
				4.5E-03	I						1.36E+09			1	0.1	Paraquat Dichloride	1910-42-5				3.5E+02	1.5E+03			2.8E+02	
				6.0E-03	H						1.36E+09			1	0.1	Parathion	56-38-2				4.7E+02	2.0E+03			3.8E+02	
				5.0E-02	H		V				1.36E+09	4.49E+04		1		Pebulate	1114-71-2				3.9E+03				3.9E+03	
				3.0E-01	O						1.36E+09			1	0.1	Pendimethalin	40487-42-1				2.3E+04	9.9E+04			1.9E+04	
				2.0E-03	I					3.12E-01	1.36E+09	5.13E+05		1		Pentabromodiphenyl Ether	32534-81-9				1.6E+02				1.6E+02	
				1.0E-04	I		V				1.36E+09			1	0.1	Pentabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-99)	60348-60-9					7.8E+00	3.3E+01			6.3E+00
9.0E-02	P			8.0E-04	I		V			4.57E+02	1.36E+09	9.65E+03		1		Pentachlorobenzene	608-93-5				7.7E+00				6.3E+01	
				3.0E-03	I		V				1.36E+09	4.32E+05		1		Pentachloronitrobenzene	82-68-8				2.7E+00				2.3E+02	
2.6E-01	H			5.0E-03	I		V				1.36E+09			0.25		Pentachlorophenol	87-86-5	1.7E+00	2.5E+00	7.5E+05	1.0E+00	3.9E+02	6.6E+02		2.5E+02	
4.0E-01	I	5.1E-06	C	9.0E-03	P						1.36E+09			1	0.1	Pentaerythritol tetranitrate (PETN)	78-11-5	1.6E+02	5.7E+02		7.0E+02	3.0E+03			5.7E+02	
4.3E-03	X			1.0E-04	X						1.36E+09			1	0.1	Pentamethylphosphoramide (PMPA)	10159-46-3					7.8E+00	3.3E+01			6.3E+00
						1.0E+00	P	V			3.88E+02	7.79E+02		1		Pentane, n-	109-66-0						8.1E+02		8.1E+02	
											1.36E+09			1	0.1	Per- and Polyfluoroalkyl Substances (PFAS)										
				3.0E-06	D						1.36E+09			1	0.1	~Ammonium perfluoro-2-methyl-3-oxahexanoate	62037-80-3					2.3E-01	9.9E-01			1.9E-01
				1.0E-03	I						1.36E+09			1	0.1	~Ammonium perfluorobutanoate	10495-86-0					7.8E+01	3.3E+02			6.3E+01
				5.0E-04	I		V				1.36E+09	1.45E+06		1		~Ammonium perfluorohexanoate	21615-47-4					3.9E+01				3.9E+01
				3.0E-06	D		V				1.36E+09			1		~Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6					2.3E-01				2.3E-01
				3.0E-04	P						1.36E+09			1	0.1	~Perfluorobutanesulfonate	45187-15-3					2.3E+01	9.9E+01			1.9E+01
				3.0E-04	P						1.36E+09			1	0.1	~Perfluorobutanesulfonic acid (PFBS)	375-73-5					2.3E+01	9.9E+01			1.9E+01
				1.0E-03	I		V			9.26E+04	1.36E+09	4.97E+04		1		~Perfluorobutanoate	45048-62-2					7.8E+01				7.8E+01
				1.0E-03	I		V			2.57E+03	1.36E+09	3.82E+04		1		~Perfluorobutanoic acid (PFBA)	375-22-4					7.8E+01				7.8E+01
				2.0E-05	A						1.36E+09			1	0.1	~Perfluorohexanesulfonate	108427-53-8					1.6E+00	6.6E+00			1.3E+00
				2.0E-05	A						1.36E+09			1	0.1	~Perfluorohexanesulfonic acid (PFHxS)	355-46-4					1.6E+00	6.6E+00			1.3E+00
				5.0E-04	I		V			6.84E+03	1.36E+09	1.43E+06		1		~Perfluorohexanoate	92612-52-7					3.9E+01				3.9E+01
				5.0E-04	I						1.36E+09			1	0.1	~Perfluorohexanoic acid (PFHxA)	307-24-4					3.9E+01	1.6E+02			3.2E+01
				3.0E-06	A						1.36E+09			1	0.1	~Perfluorononanoate	72007-68-2					2.3E-01	9.9E-01			1.9E-01
				3.0E-06	A						1.36E+09			1	0.1	~Perfluorononanoic acid (PFNA)	375-95-1					2.3E-01	9.9E-01			1.9E-01
				2.0E-06	A						1.36E+09			1	0.1	~Perfluorooctanesulfonate	45298-90-6					1.6E-01	6.6E-01			1.3E-01
				2.0E-06	A						1.36E+09			1	0.1	~Perfluorooctanesulfonic acid (PFOS)	1763-23-1					1.6E-01	6.6E-01			1.3E-01
7.0E-02	D			3.0E-06	A						1.36E+09			1	0.1	~Perfluorooctanoate	45285-51-6	9.9E+00	3.5E+01		7.8E+00	2.3E-01	9.9E-01			1.9E-01
7.0E-02	D			3.0E-06	A						1.36E+09			1	0.1	~Perfluorooctanoic acid (PFOA)	335-67-1	9.9E+00	3.5E+01		7.8E+00	2.3E-01	9.9E-01			1.9E-01
				2.0E-03	I		V			1.10E+05	1.36E+09	5.20E+04		1		~Potassium heptafluorobutanoate	2966-54-3					1.6E+02				1.6E+02
				3.0E-04	P						1.36E+09			1	0.1	~Potassium perfluorobutanesulfonate	29420-49-3					2.3E+01	9.9E+01			1.9E+01
				2.0E-06	A						1.36E+09			1	0.1	~Potassium perfluorooctanesulfonate	2795-39-3					1.6E-01	6.6E-01			1.3E-01
				1.0E-03	I		V			1.03E+05	1.36E+09	5.11E+04		1		~Sodium perfluorobutanoate	2218-54-4					7.8E+01				7.8E+01
				5.0E-04	I		V				1.36E+09	1.47E+06		1		~Sodium perfluorohexanoate	2923-26-4					3.9E+01				3.9E+01
				7.0E-04	I						1.36E+09			1		Perchlorates										
				7.0E-04	I						1.36E+09			1		~Ammonium Perchlorate	7790-98-9					5.5E+01				5.5E+01
				7.0E-04	I						1.36E+09			1		~Lithium Perchlorate	7791-03-9					5.5E+01				5.5E+01
				7.0E-04	I						1.36E+09			1		~Perchlorate and Perchlorate Salts	14797-73-0					5.5E+01				5.5E+01
				7.0E-04	I						1.36E+09			1		~Potassium Perchlorate	7778-74-7					5.5E+01				5.5E+01
				7.0E-04	I						1.36E+09			1		~Sodium Perchlorate	7601-89-0					5.5E+01				5.5E+01
				5.0E-02	I						1.36E+09			1	0.1	Permethrin	52645-53-1					3.9E+03	1.6E+04			3.2E+03
2.2E-03	C	6.3E-07	C								1.36E+09			1	0.1	Phenacetin	62-44-2	3.2E+02	1.1E+03	6.1E+06	2.5E+02					
				2.4E-01	O						1.36E+09			1	0.1	Phenmedipham	13684-63-4					1.9E+04	7.9E+04			1.5E+04
				3.0E-01	I	2.0E-01	C				1.36E+09			1	0.1	Phenol	108-95-2					2.3E+04	9.9E+04	2.8E+08		1.9E+04
				4.0E-03	I						1.36E+09			1	0.1	Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1					3.1E+02	1.3E+03			2.5E+02
				5.0E-04	X						1.36E+09			1	0.1	Phenothiazine	92-84-2					3.9E+01	1.6E+02			3.2E+01
				2.0E-04	X		V			1.29E+02	1.36E+09	7.06E+03		1		Phenyl Isothiocyanate	103-72-0					1.6E+01				1.6E+01
1.2E-01	P			6.0E-03	I						1.36E+09			1	0.1	Phenylenediamine, m-	108-45-2					4.7E+02	2.0E+03			3.8E+02
				4.0E-03	P						1.36E+09			1	0.1	Phenylenediamine, o-	95-54-5	5.8E+00	2.1E+01							

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; c = cancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; m = ceiling limit exceeded; s = Csat exceeded; V = volatile; M = mutagen.														Contaminant				Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1			
Toxicity and Chemical-specific Information														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 _o (mg/m ³)	k _e y ⁻¹	v _o l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS _o	Analyte	CAS No.	Ingestion SL TR=1E-06 (mg/kg)	Dermal SL TR=1E-06 (mg/kg)	Inhalation SL TR=1E-06 (mg/kg)	Carcinogenic SL TR=1E-06 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL TH1=1 (mg/kg)	
				1.0E+00	P						1.36E+09		1		~Tetrasodium pyrophosphate	7722-88-5					7.8E+04			7.8E+04	
				1.0E+00	P						1.36E+09		1		~Tripotassium phosphate	7778-53-2					7.8E+04			7.8E+04	
				1.0E+00	P						1.36E+09		1		~Trisodium phosphate	7601-54-9					7.8E+04			7.8E+04	
				3.0E-04	I	3.0E-04	I	V			1.36E+09		1		Phosphine	7803-51-2					2.3E+01		4.3E+05	2.3E+01	
				1.0E+00	P	1.0E-02	I				1.36E+09		1		Phosphoric Acid	7664-38-2					7.8E+04		1.4E+07	7.8E+04	
				2.0E-05	I			V			1.36E+09	6.92E+03	1		Phosphorus, White	7723-14-0					1.6E+00			1.6E+00	
1.4E-02	I	2.4E-06	C	2.0E-02	I						1.36E+09		1	0.1	~Bis(2-ethylhexyl)phthalate	117-81-7	5.0E+01	1.8E+02	1.6E+06	3.9E+01	1.6E+03	6.6E+03		1.3E+03	
1.9E-03	P			2.0E-01	I						1.36E+09		1	0.1	~Butyl Benzyl Phthalate	85-68-7	3.7E+02	1.3E+03		2.9E+02	1.6E+04	6.6E+04		1.3E+04	
				1.0E+00	I						1.36E+09		1	0.1	~Butylphthalyl Butylglycolate	85-70-1					7.8E+04	3.3E+05		6.3E+04	
				1.0E-01	I						1.36E+09		1	0.1	~Diethyl Phthalate	84-74-2					7.8E+03	3.3E+04		6.3E+03	
				8.0E-01	I						1.36E+09		1	0.1	~Diethyl Phthalate	84-66-2					6.3E+04	2.6E+05		5.1E+04	
				1.0E-01	I			V			1.36E+09	2.13E+04	1		~Dimethylterephthalate	120-61-6					7.8E+03			7.8E+03	
				1.0E-02	P						1.36E+09		1	0.1	~Octyl Phthalate, di-N-	117-84-0					7.8E+02	3.3E+03		6.3E+02	
				5.0E-01	X						1.36E+09		1	0.1	~Phthalic Acid, p-	100-21-0					3.9E+04	1.8E+05		3.2E+04	
				2.0E+00	I	2.0E-02	C				1.36E+09		1	0.1	~Phthalic Anhydride	85-44-9					1.6E+05	6.6E+05	2.8E+07	1.3E+05	
				7.0E-02	X						1.36E+09		1	0.1	Picloram	1918-02-1					5.5E+03	2.3E+04		4.4E+03	
				1.0E-04	X						1.36E+09		1	0.1	Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3					7.8E+00	3.9E+01		6.3E+00	
				2.0E-03	X						1.36E+09		1	0.1	Picric Acid (2,4,6-Trinitrophenol)	88-89-1					1.6E+02	6.6E+02		1.3E+02	
				7.3E-04	O						1.36E+09		1	0.1	Primingphos, Methyl	29232-93-7					5.7E+01	2.4E+02		4.6E+01	
3.0E+01	C	8.6E-03	C	7.0E-06	H						1.36E+09		1	0.1	Polybrominated Biphenyls	36355-01-8	2.3E-02	8.2E-02	4.4E+02	1.8E-02	5.5E-01	2.3E+00		4.4E-01	
											1.36E+09	5.86E+05	1	0.14	Polychlorinated Biphenyls (PCBs)										
7.0E-02	G	2.0E-05	G	7.0E-05	I			V			1.36E+09		1	0.14	~Aroclor 1016	12674-11-2	9.9E+00	2.5E+01	8.2E+01	6.6E+00	5.5E+00	1.6E+01		4.1E+00	
2.0E+00	G	5.7E-04	G					V			1.36E+09	2.04E+05	1	0.14	~Aroclor 1221	11104-28-2	3.5E-01	8.8E-01	1.0E+00	2.0E-01					
2.0E+00	G	5.7E-04	G					V			1.36E+09	1.12E+05	1	0.14	~Aroclor 1232	11141-16-5	3.5E-01	8.8E-01	5.5E-01	1.7E-01					
2.0E+00	G	5.7E-04	G					V			1.36E+09	5.91E+05	1	0.14	~Aroclor 1242	53469-21-9	3.5E-01	8.8E-01	2.9E+00	2.3E-01					
2.0E+00	G	5.7E-04	G					V			1.36E+09	5.14E+05	1	0.14	~Aroclor 1248	12672-29-6	3.5E-01	8.8E-01	2.5E+00	2.3E-01					
2.0E+00	G	5.7E-04	G	2.0E-05	I			V			1.36E+09	8.43E+05	1	0.14	~Aroclor 1254	11097-69-1	3.5E-01	8.8E-01	4.1E+00	2.4E-01	1.6E+00	4.7E+00		1.2E+00	
2.0E+00	G	5.7E-04	G					V			1.36E+09	1.31E+06	1	0.14	~Aroclor 1260	11096-82-5	3.5E-01	8.8E-01	6.5E+00	2.4E-01					
				6.0E-04	X			V			1.36E+09	1.15E+06	1	0.14	~Aroclor 5460	11126-42-4					4.7E+01	1.4E+02		3.5E+01	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	2.43E+06	1	0.14	~Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	1.8E-01	4.5E-01	6.0E+00	1.3E-01	1.8E+00	5.5E+00	3.4E+03	1.4E+00	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.58E+06	1	0.14	~Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167)	52663-72-6	1.8E-01	4.5E-01	3.9E+00	1.2E-01	1.8E+00	5.5E+00	2.2E+03	1.4E+00	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.04E+06	1	0.14	~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	1.8E-01	4.5E-01	2.6E+00	1.2E-01	1.8E+00	5.5E+00	1.4E+03	1.4E+00	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.11E+06	1	0.14	~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 156)	38380-08-4	1.8E-01	4.5E-01	2.7E+00	1.2E-01	1.8E+00	5.5E+00	1.5E+03	1.4E+00	
3.9E+03	W	1.1E+00	W	2.3E-08	W	1.3E-06	W	V			1.36E+09	1.58E+06	1	0.14	~Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	1.8E-04	4.5E-04	3.9E-03	1.2E-04	1.8E-03	5.5E-03	2.2E+00	1.4E-03	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	7.33E+05	1	0.14	~Pentachlorobiphenyl, 2,3,4,4',5'- (PCB 123)	65510-44-3	1.8E-01	4.5E-01	1.8E+00	1.2E-01	1.8E+00	5.5E+00	1.0E+03	1.4E+00	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	5.90E+05	1	0.14	~Pentachlorobiphenyl, 2,3,4,4',5'- (PCB 118)	31508-00-6	1.8E-01	4.5E-01	1.5E+00	1.2E-01	1.8E+00	5.5E+00	8.2E+02	1.4E+00	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	6.01E+05	1	0.14	~Tetrachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	1.8E-01	4.5E-01	1.5E+00	1.2E-01	1.8E+00	5.5E+00	8.4E+02	1.4E+00	
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	W	V			1.36E+09	1.05E+06	1	0.14	~Pentachlorobiphenyl, 2,3,4,4',5'- (PCB 114)	74472-37-0	1.8E-01	4.5E-01	2.6E+00	1.2E-01	1.8E+00	5.5E+00	1.5E+03	1.4E+00	
1.3E+04	W	3.8E+00	W	7.0E-09	W	4.0E-07	W	V			1.36E+09	7.26E+05	1	0.14	~Pentachlorobiphenyl, 3,3',4,4',5'- (PCB 126)	57465-28-8	5.3E-05	1.4E-04	5.4E-04	3.6E-05	5.5E-04	1.6E-03	3.0E-01	4.1E-04	
2.0E+00	I	5.7E-04	I					V			1.36E+09	5.32E+05	1	0.14	~Polychlorinated Biphenyls (high risk)	1336-36-3	3.5E-01	8.8E-01	2.6E+00	2.3E-01					
4.0E-01	I	1.0E-04	I					V			1.36E+09		1	0.14	~Polychlorinated Biphenyls (low risk)	1336-36-3									
7.0E-02	I	2.0E-05	I					V			1.36E+09		1	0.14	~Polychlorinated Biphenyls (lowest risk)	1336-36-3									
1.0E-01	W	3.8E-03	W	7.0E-06	W	4.0E-04	W	V			1.36E+09		1	0.14	~Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	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	W	1.1E-02	W	2.3E-06	W	1.3E-04	W	V			1.36E+09	5.09E+05	1	0.14	~Tetrachlorobiphenyl, 3,4,4',5'- (PCB 81)	70362-50-4	1.8E-02	4.5E-02	1.3E-01	1.2E-02	1.8E-01	5.5E-01	7.1E-01	1.4E-01	
				6.0E-04	I			V			1.36E+09		1	0.1	~Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9					8.5E+05			8.5E+05	
				6.0E-02	I			V			1.36E+09	1.41E+05	1	0.13	Polynuclear Aromatic Hydrocarbons (PAHs)						4.7E+03	1.5E+04		3.6E+03	
				3.0E-01	I			V			1.36E+09	5.23E+05	1	0.13	~Acenaphthene	120-12-7					2.3E+04	7.6E+04		1.8E+04	
1.0E-01	E	6.0E-05	E					V	M		1.36E+09	4.41E+06	1	0.13	~Benz[a]anthracene	56-55-3	1.5E+00	4.6E+00	7.4E+01	1.1E+00					
				9.0E-05	X	2.0E-06	X				1.36E+09		1	0.1	~Benzo[e]pyrene	192-97-2					7.0E+00	3.0E+01	2.8E+03	5.7E+00	
1.2E+00	C	1.1E-04	C					V			1.36E+09		1	0.13	~Benzo[j]fluoranthene	205-82-3	5.8E-01	1.6E+00	3						

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

Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child 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	vo l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS _d	Analyte	CAS No.	Ingestion SL TR=1E-06 (mg/kg)	Dermal SL TR=1E-06 (mg/kg)	Inhalation SL TR=1E-06 (mg/kg)	Carcinogenic SL TR=1E-06 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THQ=1 (mg/kg)	
2.0E-03	I									1.11E+05	1.36E+09	6.27E+04	1		Propargyl Alcohol	107-19-7					1.6E+02				1.6E+02
2.0E-02	I										1.36E+09		1	0.1	Propazine	139-40-2					1.6E+03	6.6E+03			1.3E+03
2.0E-02	I										1.36E+09		1	0.1	Propham	122-42-9					1.6E+03	6.6E+03			1.3E+03
1.0E-01	O										1.36E+09		1	0.1	Propiconazole	60207-90-1					7.8E+03	3.3E+04			6.3E+03
1.0E-01	X									3.26E+04	1.36E+09	8.94E+03	1		Propionaldehyde	123-38-6							7.5E+01		7.5E+01
1.0E-01	X									1.0E+00	1.36E+09	6.99E+03	1		Propyl benzene	103-65-1					7.8E+03		7.3E+03		3.8E+03
2.0E+01	P									3.0E+00	1.36E+09	7.04E+02	1		Propylene	115-07-1							2.2E+03		2.2E+03
										2.7E-04	1.36E+09		1	0.1	Propylene Glycol	57-55-6					1.6E+06	6.6E+06			1.3E+06
										7.0E-01	1.36E+09	7.83E+04	1		Propylene Glycol Dinitrate	6423-43-4							3.9E+05		3.9E+05
2.4E-01	I	3.7E-06	I							1.06E+05	1.36E+09	1.03E+04	1		Propylene Glycol Monomethyl Ether	107-98-2	2.9E+00		7.8E+00	2.1E+00	5.5E+04		1.6E+05		4.1E+04
										1.0E-03	1.36E+09	5.54E+04	1		Propylene Oxide	75-56-9							3.2E+02		3.2E+02
										5.0E-04	1.36E+09		1	0.1	Pyridine	110-86-1					7.8E+01				7.8E+01
3.0E+00	I										1.36E+09		1	0.1	Quinalphos	13593-03-8	2.3E-01	8.2E-01		1.8E-01	3.9E+01	1.6E+02			3.2E+01
											1.36E+09		1	0.1	Quinoline	91-22-5									
											1.36E+09		1	0.1	Quizalofop-ethyl	76578-14-8					7.0E+02	3.0E+03			5.7E+02
											1.36E+09		1		Refractory Ceramic Fibers (units in fibers)	E715557									
											1.36E+09		1	0.1	Resmethrin	10453-86-8					2.3E+03	9.9E+03			1.9E+03
											1.36E+09	4.65E+05	1		Ronnel	299-84-3					3.9E+03				3.9E+03
2.2E-01	C	6.3E-05	C								1.36E+09		1	0.1	Rotenone	83-79-4					3.1E+02	1.3E+03			2.5E+02
											1.36E+09		1	0.1	Safrole	94-59-7	7.0E-01	2.7E+00	2.2E+04	5.5E-01					
											1.36E+09		1	0.1	Selenious Acid	7783-00-8					3.9E+02				3.9E+02
											1.36E+09		1		Selenium	7782-49-2					3.9E+02		2.8E+07		3.9E+02
											1.36E+09		1		Selenium Sulfide	7446-34-6					3.9E+02		2.8E+07		3.9E+02
											1.36E+09		1	0.1	Sethoxydim	74051-80-2					1.1E+04	4.6E+04			8.8E+03
											1.36E+09		1		Silica (crystalline, respirable)	7631-86-9							4.3E+06		4.3E+06
1.2E-01	H										1.36E+09	0.04			Silver	7440-22-4					3.9E+02				3.9E+02
											1.36E+09		1	0.1	Simazine	122-34-9	5.8E+00	2.1E+01		4.5E+00	3.9E+02	1.6E+03			3.2E+02
											1.36E+09		1	0.1	Sodium Acifluorfen	62476-59-9					1.0E+03	4.3E+03			8.2E+02
											1.36E+09		1		Sodium Azide	26628-22-8					3.1E+02				3.1E+02
2.7E-01	H										1.36E+09		1	0.1	Sodium Diethyldithiocarbamate	148-18-5	2.6E+00	9.2E+00		2.0E+00	2.3E+03	9.9E+03			1.9E+03
											1.36E+09		1		Sodium Fluoride	7681-49-4					3.9E+03		2.0E+07		3.9E+03
											1.36E+09		1	0.1	Sodium Fluoroacetate	62-74-8					1.6E+00	6.6E+00			1.3E+00
											1.36E+09		1		Sodium Metavanadate	13718-26-8					7.8E+01				7.8E+01
											1.36E+09		1		Sodium Tungstate	13472-45-2					6.3E+01				6.3E+01
2.4E-02	H										1.36E+09		1	0.1	Sodium Tungstate Dihydrate	10213-10-2					6.3E+01				6.3E+01
											1.36E+09		1		Stirofos (Tetrachlorovinphos)	961-11-5	2.9E+01	1.0E+02		2.3E+01	2.3E+03	9.9E+03			1.9E+03
											1.36E+09		1		Strontium, Stable	7440-24-6					4.7E+04				4.7E+04
											1.36E+09		1	0.1	Strychnine	57-24-9					2.3E+01	9.9E+01			1.9E+01
											1.36E+09	9.35E+03	1		Styrene	100-42-5					1.6E+04		9.7E+03		6.0E+03
											1.36E+09		1	0.1	Styrene-Acrylonitrile (SAN) Trimer (THNA isomer)	57964-39-3					2.3E+02	9.9E+02			1.9E+02
											1.36E+09		1	0.1	Styrene-Acrylonitrile (SAN) Trimer (THNP isomer)	57964-40-6					2.3E+02	9.9E+02			1.9E+02
											1.36E+09		1	0.1	Sulfolane	126-33-0					7.8E+01	3.3E+02	2.8E+06		6.3E+01
											1.36E+09		1	0.1	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9					6.3E+01	2.6E+02			5.1E+01
											1.36E+09		1		Sulfur Trioxide	7446-11-9							1.4E+06		1.4E+06
											1.36E+09		1		Sulfuric Acid	7664-93-9							1.4E+06		1.4E+06
2.5E-02	I	7.1E-06	I								1.36E+09		1	0.1	Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl	140-57-8	2.8E+01	9.9E+01	5.4E+05	2.2E+01	3.9E+03	1.6E+04			3.2E+03
											1.36E+09		1	0.1	TCMTB	21564-17-0					2.3E+03	9.9E+03			1.9E+03
											1.36E+09		1	0.1	Tebuthiuron	34014-18-1					5.5E+03	2.3E+04			4.4E+03
											1.36E+09		1	0.1	Temephos	3383-96-8					1.6E+03	6.6E+03			1.3E+03
											1.36E+09		1	0.1	Terbacil	5902-51-2					1.0E+03	4.3E+03			8.2E+02
											1.36E+09	2.64E+05	1		Terbufos	13071-79-9					2.0E+00				2.0E+00
5.0E-03	C	1.3E-06	C								1.36E+09		1	0.1	Terbutryn	886-50-0					7.8E+01	3.3E+02			6.3E+01
											1.36E+09	3.99E+03	1		Tert-Butyl Acetate	540-88-5	1.4E+02		8.6E+00	8.1E+00					
											1.36E+09		1	0.1	Tetrabromodiphenyl ether, 2,2',4,4'-(BDE-47)	5436-43-1					7.8E+00	3.3E+01			6.3E+00
											1.36E+09	5.07E+04	1		Tetrachlorobenzene, 1,2,4,5-	95-94-3					2.3E+00				2.3E+00
2.6E-02	I	7.4E-06	I								1.36E+09	5.68E+03	1		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								1.90E+03	1.51E+04	1		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								1.66E+02	2.35E+03	1		Tetrachloroethylene	127-18-4	3.3E+02		2.5E+01	2.4E+01	4.7E+02		9.8E+01		8.1E+01
											1.36E+09		1	0.1	Tetrachlorophenol, 2,3,4,6-	58-90-2									

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; W = TEF applied; E = RPF applied; G = see user's guide; U = user provided; c = cancer; n = noncancer; * = where: nc SL < 100X ca SL; ** = where nc SL < 10X ca SL; SSL values are based on DAF=1; m = ceiling limit exceeded; s = Csat exceeded; V = volatile; M = mutagen.															Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Child Hazard Index (HI) = 1				
Toxicity and Chemical-specific Information															Analyte	CAS No.	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 _o (mg/m ³)	k _e y ⁻¹	Vol	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS _d			Ingestion SL TR=1E-06 (mg/kg)	Dermal SL TR=1E-06 (mg/kg)	Inhalation SL TR=1E-06 (mg/kg)	Carcinogenic SL TR=1E-06 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)	
				1.5E-02	O						1.36E+09			1	0.1	Thiram	137-26-8					1.2E+03	4.9E+03	9.5E+02	
				6.0E-01	H						1.36E+09			1		Tin	7440-31-5					4.7E+04		4.7E+04	
				8.0E-02	I	1.0E-04 A V				8.18E+02	1.36E+09	4.29E+03		1		Titanium Tetrachloride	7550-45-0					1.4E+05	1.4E+05	4.7E+04	
3.9E-02	C	1.1E-05 C		2.0E-04	X	5.0E+00 I V					1.36E+09	7.62E+05		1		Toluene	108-88-3					6.3E+03	2.2E+04	4.9E+03	
1.8E-01	X			8.0E-06	C V						1.36E+09			1		Toluene-2,4-disocyanate	584-84-9	1.8E+01	1.9E+02	1.6E+01		6.4E+00	6.4E+00	6.4E+00	
3.9E-02	C	1.1E-05 C		2.0E-04	X	8.0E-06 C V				1.71E+03	1.36E+09	6.32E+05		1	0.1	Toluene-2,5-diamine	95-70-5	3.9E+00	1.4E+01	3.0E+00		1.6E+01	6.6E+01	1.3E+01	
				1.0E-04	X						1.36E+09			1	0.1	Toluene-2,6-disocyanate	91-08-7	1.8E+01	1.6E+02	1.6E+01		5.3E+00	5.3E+00	5.3E+00	
				1.0E-04	X						1.36E+09			1	0.1	Toluenediamine, 2,3-	2687-25-4					7.8E+00	3.3E+01	6.3E+00	
				1.0E-04	X						1.36E+09			1	0.1	Toluenediamine, 3,4-	496-72-0					7.8E+00	3.3E+01	6.3E+00	
				5.0E-03	P						1.36E+09			1	0.1	Toxic Acid, p-	99-34-5					3.9E+02	1.6E+03	3.2E+02	
1.6E-02	P	5.1E-05 C		4.0E-03	X						1.36E+09			1	0.1	Toluidine, o- (Methylaniline, 2-)	95-34-4	4.3E+01	1.5E+02	7.5E+04	3.4E+01	7.8E+00	3.3E+01	6.3E+00	
3.0E-02	P			3.0E+00	P					3.42E-01	1.36E+09	1.38E+03		1	0.1	Toluidine, p-	106-49-0	2.3E+01	8.2E+01	1.8E+01		3.9E+02	1.3E+03	2.5E+02	
				5.0E-03	P	4.0E-01 P V					1.12E+02	1.36E+09	1.65E+03		1		Total Petroleum Hydrocarbons (Aliphatic High)	E1790670					2.3E+05	2.3E+05	2.3E+05
				1.0E-02	X	1.0E-01 P V				6.86E+00	1.36E+09	1.04E+03		1		Total Petroleum Hydrocarbons (Aliphatic Low)	E1790666					3.9E+02	6.9E+02	2.5E+02	
				3.0E-04	P	2.0E-06 P M					1.36E+09			1	0.13	Total Petroleum Hydrocarbons (Aliphatic Medium)	E1790668					7.8E+02	1.1E+02	9.6E+01	
				1.0E-02	P	6.0E-02 P V				2.31E+02	1.36E+09	7.75E+03		1		Total Petroleum Hydrocarbons (Aromatic High)	E1790676					2.3E+01	7.6E+01	1.8E+01	
1.1E+00	I	3.2E-04 I		9.0E-05	P						1.36E+09			1	0.1	Total Petroleum Hydrocarbons (Aromatic Medium)	E1790674					7.8E+02	4.8E+02	3.0E+02	
				3.0E-05	X						1.36E+09			1	0.1	Toxaphene	8001-35-2	6.3E-01	2.2E+00	1.2E+04	4.9E-01	7.0E+00	3.0E+01	5.7E+00	
				8.0E-01	X						1.36E+09			1	0.1	Toxaphene, Weathered	E1841606					2.3E+00	9.9E+00	1.9E+00	
				3.4E-02	O						1.36E+09			1	0.1	Tralometrin	68841-25-6					5.9E+02	2.5E+03	4.7E+02	
7.2E-02	O			2.5E-02	O						1.36E+09	3.62E+05		1	0.1	Tri-n-butyltin	688-73-3					2.3E+01	2.5E+03	2.3E+01	
				1.0E-02	I						1.36E+09			1	0.1	Triacetin	102-76-1					6.3E+06	2.6E+07	5.1E+06	
				8.0E-03	I						1.36E+09			1	0.1	Triadimefon	43121-43-3					2.7E+03	1.1E+04	2.1E+03	
				5.0E-03	I						1.36E+09	4.83E+04		1	0.1	Triallate	2303-17-5	9.7E+00		9.7E+00		2.0E+03		2.0E+03	
				9.0E-03	X						1.36E+09			1	0.1	Triasulfuron	82097-50-5					7.8E+02	3.3E+03	6.3E+02	
				2.0E-04	P						1.36E+09			1	0.1	Tribenuron-methyl	101200-48-0					6.3E+02	2.6E+03	5.1E+02	
				1.0E-02	P						1.36E+09			1	0.1	Tribromobenzene, 1,2,4-	615-54-3					3.9E+02		3.9E+02	
9.0E-03	P			3.0E-04	P						1.36E+09			1	0.1	Tribromophenol, 2,4,6-	118-79-6					7.0E+02	3.0E+03	5.7E+02	
				3.0E-04	P						1.36E+09			1	0.1	Tributyl Phosphate	78-48-8					1.6E+01	6.6E+01	1.3E+01	
				3.0E-04	P						1.36E+09			1	0.1	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						1.36E+09			1	0.1	Tributyltin Compounds	E1790679					2.3E+01	9.9E+01	1.9E+01	
				3.0E-04	P						1.36E+09			1	0.1	Tributyltin Oxide	56-35-9					2.3E+01	9.9E+01	1.9E+01	
				3.0E+01	I	5.0E+00 P V				9.10E+02	1.36E+09	1.29E+03		1		Trichloramine	10025-85-1					2.3E+06		6.7E+03	
7.0E-02	I			2.0E-02	I						1.36E+09			1	0.1	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1					2.3E+06		6.7E+03	
2.9E-02	H			3.0E-05	X						1.36E+09			1	0.1	Trichloroacetic Acid	76-03-9					1.6E+03	6.6E+03	1.3E+03	
7.0E-03	X			8.0E-04	X						1.36E+09	3.22E+04		1	0.1	Trichloroaniline HCl, 2,4,6-	33663-50-2	9.9E+00	3.5E+01	7.8E+00		1.6E+03	6.6E+03	1.3E+03	
2.9E-02	P			2.0E+00	I	2.0E-03 P V				4.04E+02	1.36E+09	2.99E+04		1		Trichloroaniline, 2,4,6-	634-93-5	2.4E+01				9.9E+01	8.5E+01	1.9E+01	
				5.7E-02	I	1.6E-05 I				2.16E+03	1.36E+09	7.22E+03		1		Trichlorobenzene, 1,2,3-	87-61-6	9.9E+01				2.3E+00	9.9E+00	1.9E+00	
4.6E-02	I	4.1E-06 I		3.0E-01	I	2.0E-03 I V				6.92E+02	1.36E+09	2.21E+03		1		Trichlorobenzene, 1,2,4-	120-82-1					6.3E+01	6.6E+01	6.3E+01	
				1.0E-01	I	2.0E-03 I V				1.23E+03	1.36E+09	1.04E+03		1	0.1	Trichlorobenzene, 1,1,1-	71-55-6	2.4E+01		2.4E+01		1.6E+05	6.2E+01	8.1E+03	
1.1E-02	I	3.1E-06 I		1.0E-03	I						1.36E+09			1	0.1	Trichloroethane, 1,1,2-	79-00-5					3.1E+02	1.5E+00	1.5E+00	
				1.0E-02	I						1.36E+09			1	0.1	Trichloroethylene	79-01-6	8.8E+00	1.3E+00	1.1E+00		3.9E+01	4.6E+00	4.1E+00	
				5.0E-03	I						1.36E+09			1	0.1	Trichlorofluoromethane	75-69-4					2.3E+04		2.3E+04	
				8.0E-03	I						1.36E+09			1	0.1	Trichlorophenol, 2,4,5-	95-95-4					7.8E+03	3.3E+04	6.3E+03	
				5.0E-03	I						1.36E+09			1	0.1	Trichlorophenol, 2,4,6-	88-06-2	6.3E+01	2.2E+02	1.2E+06	4.9E+01	7.8E+01	3.3E+02	6.3E+01	
				3.0E+01	I						1.36E+09			1	0.1	Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5					7.8E+02	3.3E+03	6.3E+02	
				4.0E-03	I						1.36E+09			1	0.1	Trichlorophenoxypropionic acid, -2,4,5	93-72-1					6.3E+02	2.6E+03	5.1E+02	
				3.0E-03	X	3.0E-04 P V				1.28E+03	1.36E+09	1.50E+04		1		Trichloropropane, 1,1,2-	598-77-6					3.9E+02		3.9E+02	
				2.0E-02	A					1.40E+03	1.36E+09	1.57E+04		1		Trichloropropane, 1,2,3-	96-18-4	5.1E-03		5.1E-03		3.1E+02	4.9E+00	4.8E+00	
				3.0E-03	I						1.36E+09			1	0.1	Trichloropropene, 1,2,3-	96-19-5					2.3E+02		7.3E-01	
				2.0E-02	A						1.36E+09			1	0.1	Tricresyl Phosphate (TCP)	1330-78-5					1.6E+03	6.6E+03	1.3E+03	
				2.0E+00	P	7.0E-03 I V				2.79E+04	1.36E+09	1.58E+04		1	0.1	Tridiphane	58138-08-2					2.3E+02	9.9E+02	1.9E+02	
				7.7E-03	I						1.36E+09	5.13E+05		1	0.1	Triethylamine	121-44-8					1.6E+05	6.6E+05	1.2E+02	
2.0E-02	P			1.0E-02	P	2.0E+01 P V				4.81E+03	1.36E+09	7.12E+02		1		Triethylene Glycol	112-27-6					1.6E+05	6.6E+05	1.3E+05	
				1.0E-02	I	6.0E-02 I V				2.93E+02	1.36E+09	9.44E+03		1		Trifluoroethane, 1,1,1-	420-46-2					1.6E+05	6.6E+05	1.5E+04	
				1.0E-02	I	6.0E-02 I V				2.19E+02	1.36E+09	7.91E+03		1		Trifluralin	1582-09-8	9.0E+01		9.0E+01		5.9E+02		5.9E+02	
				1.0E-02	X					1.82E+02	1.36E+09	6.61E+03		1		Trimethyl Phosphate	512-56-1	3.5E+01	1.2E+02	2.7E+01		7.8E+02	3.3E+03	6.3E+02	
				3.0E-02	I						1.36E+09			1	0.019	Trimethylbenzene, 1,2,3-	526-73-8					7.8E+02		5.9E+02	
				5.0E-04	I						1.36E+09			1	0.032	Trimethylbenzene, 1,2,4-	95-63-6					7.8E+02		3.0E+02	
				2.0E-02	P						1.36E+09			1		Trimethylbenzene, 1,3,5-	108-67-8					7.8E+02		2.7E+02	
				1.0E-02	X					2.96E+01	1.36E+09	1.00E+03		1		Trimethylbenzene, 2,4,4-	25167-70-8					7.8E+02		7	

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

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)	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS _d	Analyte	CAS No.	Ingestion SL TR=1E-06 (mg/kg)	Dermal SL TR=1E-06 (mg/kg)	Inhalation SL TR=1E-06 (mg/kg)	Carcinogenic SL TR=1E-06 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)	
7.2E-01	I	1.5E-05	P	3.0E-03	I	1.0E-01	I	V	M	2.47E+03	1.36E+09	1.37E+03	1		Vinyl Bromide	593-60-2			2.6E-01	2.6E-01			4.3E+00		4.3E+00
		4.4E-06	I	3.0E-03	I	1.0E-01	I	V		3.92E+03	1.36E+09	9.56E+02	1		Vinyl Chloride	75-01-4	9.4E-02		1.6E-01	5.9E-02	2.3E+02		1.0E+02	7.0E+01	
				3.0E-04	I						1.36E+09		1	0.1	Warfarin	81-81-2					2.3E+01	9.9E+01		1.9E+01	
				2.0E-01	G	1.0E-01	G	V		3.88E+02	1.36E+09	5.47E+03	1		Xylene, m-	108-38-3					1.6E+04		5.7E+02	5.5E+02	
				2.0E-01	G	1.0E-01	G	V		4.34E+02	1.36E+09	6.45E+03	1		Xylene, o-	95-47-6					1.6E+04		6.7E+02	6.4E+02	
				2.0E-01	G	1.0E-01	G	V		3.90E+02	1.36E+09	5.58E+03	1		Xylene, p-	106-42-3					1.6E+04		5.8E+02	5.6E+02	
				2.0E-01	I	1.0E-01	I	V		2.60E+02	1.36E+09	5.74E+03	1		Xylenes	1330-20-7					1.6E+04		6.0E+02	5.8E+02	
				3.0E-04	I						1.36E+09		1		Zinc Phosphide	1314-84-7					2.3E+01			2.3E+01	
				3.0E-01	I						1.36E+09		1		Zinc and Compounds	7440-66-6					2.3E+04			2.3E+04	
				5.0E-02	I						1.36E+09		1	0.1	Zinc	12122-67-7					3.9E+03	1.6E+04		3.2E+03	
		8.0E-05	X								1.36E+09		1		Zirconium	7440-67-7					6.3E+00			6.3E+00	

TR=1E-06
THQ=1.0