

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; T = ATSDR DRAFT; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; R = ORD; N = WI; W = TEF applied; E = RPF applied; G = see user's guide; 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	RD ₅₀ (mg/kg-day)	k e y	RIC ₁ (mg/m ³)	k e y	v o	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS ₀	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)		
2.20E-06	I	3.00E-04	O	9.00E-03	I	V				1.07E+05	1.36E+09	8.72E+03	1	0.1	Acetophenone	30660-19-1			1.1E+01	1.1E+01	2.3E+01	9.9E+01			1.9E+01	
2.00E-02	I	9.00E-01	V	1.14E+05						1.36E+09	1.37E+04	1	0.1	Acetaldehyde	75-07-0									8.2E+01	8.2E+01	
9.00E-01	I	2.00E-03	X	1.28E+05						1.36E+09	1.30E+04	1	0.1	Acetochlor	34256-82-1					1.6E+03	6.6E+03			1.3E+03	1.3E+03	
3.80E+00	C	1.30E-03	C	1.00E-01						2.52E+03	5.97E+04	1		Acetone	67-64-1					7.0E+04				7.0E+04	7.0E+04	
5.00E-01	I	1.00E-04	I	2.00E-03	I	V				1.36E+09	1.30E+04	1	0.1	Acetone Cyanohydrin	75-86-5									2.8E+06	2.8E+06	
5.40E-01	I	6.80E-05	I	9.00E-05	T	2.00E-03	I	V		1.13E+04	1.36E+09	7.69E+03	1		Acetonitrile	75-05-8							8.1E+02	8.1E+02	8.1E+02	
5.60E-02	C	1.00E-02	I	1.00E-03						1.36E+09	5.97E+04	1		Acetophenone	98-86-2	1.8E-01	6.5E-01	2.9E+03	1.4E-01	7.8E+03				7.8E+03	7.8E+03	
1.70E+01	I	4.90E-03	I	3.00E-05						1.36E+09	1.72E+06	1		Acetylaminofluorene, 2-	53-96-3									1.4E-01	1.4E-01	
2.10E-02	C	6.00E-06	C	4.00E-03	P	1.00E-04	X	V		1.11E+05	1.36E+09	3.42E+04	1		Acrolein	107-02-8	1.8E-01	6.5E-01	2.9E+03	1.4E-01	7.8E+03				7.8E+03	
4.00E-04	I	1.00E+00	P	5.00E-03	P					1.42E+03	1.58E+03	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	1.3E+02	
2.10E+01	C	6.00E-03	C	4.00E-04	I	9.00E-03	I			1.36E+09	1.36E+09	6.91E+03	1		Acrylic Acid	79-10-7	3.1E-01	1.2E+00	1.4E+04	2.4E-01	1.6E+02	6.6E+02	8.5E+06		1.3E+02	1.3E+02
8.00E-02	P	4.00E-03	X	2.00E-02						1.36E+09	1.36E+09	1	0.1	Acrylonitrile	107-13-1	1.3E+00		3.2E-01	2.5E-01	7.0E+00			1.6E+01	4.9E+00		
4.00E-03	X	2.00E-02	P	2.50E-03	I					1.36E+09	1.36E+09	1	0.1	Adiponitrile	111-69-3									8.5E+06	8.5E+06	
2.00E-02	P	2.00E-02	P	2.00E-02						1.36E+09	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	6.3E+02	
2.50E-03	I	5.00E-01	I	2.00E-03	I	V				1.36E+09	1.36E+09	1	0.1	Aldicarb	116-06-3									6.3E+01	6.3E+01	
2.00E-03	X	2.00E-01	I	3.00E-03	X	V				1.36E+09	1.36E+09	1	0.1	Aldicarb Sulfone	1646-88-4					7.8E+01	3.3E+02			6.3E+01	6.3E+01	
5.70E-03	I	1.60E-06	C	7.00E-03	P	1.00E-03	I	V		1.37E+04	2.62E+04	1	0.1	Aldicarb sulfoxide	1646-87-3									3.3E+02	3.3E+02	
4.00E-02	P	4.00E-04	I	3.00E-04	A					1.36E+09	1.36E+09	1	0.1	Aldrin	309-00-2	4.1E-02		9.8E-01	3.9E-02	2.3E+00				2.3E+00	2.3E+00	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P			1.36E+09	1.36E+09	1	0.1	Allyl Alcohol	107-18-6	3.3E+01		7.4E-01	7.2E-01	3.1E+02				3.6E+00	3.6E+00	
2.30E-01	C	3.00E-03	A	3.00E-03	X	V				1.36E+09	1.36E+09	2.62E+04	1		Allyl Chloride	107-05-1					1.6E+04				1.6E+04	1.6E+04
8.80E-01	C	2.50E-04	C	4.00E-04	I					1.36E+09	1.36E+09	1	0.1	Ammonia	7773-06-0	1.2E+02	4.3E+02	2.4E+06	9.5E+01	5.5E+02	2.3E+03	8.2E+01	1.4E+06	4.4E+02	4.4E+02	
1.10E-01	I	3.10E-05	I	3.00E-04	A	1.00E-02	A			1.36E+09	5.23E+05	1	0.1	Ammonium Picrate	131-74-8	1.7E+01	6.2E+01	2.4E+06	1.4E+01	1.6E+02	6.6E+02			1.3E+02	1.3E+02	
2.30E-01	C	3.00E-03	A	3.00E-03	X	V				1.36E+09	1.36E+09	1	0.1	Ammonium Sulfamate	7773-06-0					1.6E+04				1.6E+04	1.6E+04	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P			1.36E+09	1.36E+09	1	0.1	Amyl Alcohol, tert-	75-85-4	1.2E+02	4.3E+02	2.4E+06	9.5E+01	5.5E+02	2.3E+03	8.2E+01	1.4E+06	4.4E+02	4.4E+02	
2.30E-01	C	3.00E-03	A	3.00E-03	X	V				1.36E+09	1.36E+09	1	0.1	Aniline	62-53-3					1.6E+04				1.6E+04	1.6E+04	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P			1.36E+09	1.36E+09	1	0.1	Anthraquinone, 9,10-	84-65-1	1.7E+01	6.2E+01	2.4E+06	1.4E+01	1.6E+02	6.6E+02			1.3E+02	1.3E+02	
2.30E-01	C	3.00E-03	A	3.00E-03	X	V				1.36E+09	1.36E+09	1	0.1	Antimony (metallic)	7440-36-0					3.1E+01				3.1E+01	3.1E+01	
1.50E+00	I	4.30E-03	I	3.00E-04	I	1.50E-05	C			1.36E+09	1.36E+09	0.15	0.03	Antimony Pentoxide	1314-60-9	7.7E-01	5.5E+00	8.9E+02	6.8E-01	1.6E+03	6.6E+03			1.3E+03	1.3E+03	
2.30E-01	C	3.00E-03	A	3.00E-03	X	V				1.36E+09	1.36E+09	1	0.1	Antimony Trioxide	1332-81-6					3.1E+01				3.1E+01	3.1E+01	
8.80E-01	C	2.50E-04	C	4.00E-04	I					1.36E+09	1.36E+09	1	0.1	Antimony Trioxide	1309-64-4					2.8E+05				2.8E+05	2.8E+05	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P			1.36E+09	1.36E+09	1	0.1	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	3.5E+01	3.5E+01	
2.30E-01	C	3.00E-03	A	3.00E-03	X	V				1.36E+09	1.36E+09	1	0.1	Arsine	7784-42-1					2.7E-01				2.7E-01	2.7E-01	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P			1.36E+09	1.36E+09	1	0.1	Asbestos (units in fibers)	1332-21-4					2.8E+04	1.2E+05			2.3E+04	2.3E+04	
2.30E-01	C	3.00E-03	A	3.00E-03	X	V				1.36E+09	1.36E+09	1	0.1	Asulam	3337-71-1	3.0E+00	1.1E+01		2.4E+00	2.3E+02	9.9E+02			1.9E+02	1.9E+02	
8.80E-01	C	2.50E-04	C	4.00E-04	I					1.36E+09	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			1.9E+02	1.9E+02	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P			1.36E+09	1.36E+09	1	0.1	Auramine	492-60-8	7.9E-01	2.8E+00	1.5E+04	6.2E-01	3.1E+01	1.3E+02			2.5E+01	2.5E+01	
2.30E-01	C	3.00E-03	A	3.00E-03	X	V				1.36E+09	1.36E+09	1	0.1	Avermectin B1	85-195-55-3					3.1E+01	1.3E+02			2.5E+01	2.5E+01	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P			1.36E+09	1.36E+09	1	0.1	Azobenzene	86-50-0	6.3E+00	4.7E+01	5.6E+00		2.3E+02	9.9E+02			1.4E+07	1.9E+02	
2.30E-01	C	3.00E-03	A	3.00E-03	X	V				1.36E+09	1.36E+09	1	0.1	Azodicarbonamide	103-33-3					7.8E+04	3.3E+05	9.9E+03		8.6E+03	8.6E+03	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P			1.36E+09	1.36E+09	1	0.1	Barium	7440-39-3					1.6E+04				1.5E+04	1.5E+04	
2.30E-01	C	3.00E-03	A	3.00E-03	X	V				1.36E+09	1.36E+09	1	0.1	Benfluralin	1861-40-1					3.9E+02				3.9E+02	3.9E+02	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P			1.36E+09	1.36E+09	1	0.1	Benomyl	17804-35-2					3.9E+03	1.6E+04			3.2E+03	3.2E+03	
2.30E-01	C	3.00E-03	A	3.00E-03	X	V				1.36E+09	1.36E+09	1	0.1	Bensulfuron-methyl	83055-99-6					1.6E+04	6.6E+04			1.3E+04	1.3E+04	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P			1.36E+09	1.36E+09	1	0.1	Benzenesulfonamide	25057-89-0					2.3E+03	9.9E+03			1.9E+03	1.9E+03	
5.50E-02	I	7.80E-06	I	4.00E-03	I	3.00E-02	I	V		1.82E+03	1.36E+09	3.54E+03	1		Benzaldehyde	100-52-7	1.7E+02			1.7E+02	7.8E+03			7.8E+03	7.8E+03	
1.00E-01	X	3.00E-04	X	4.00E-03	C	V				1.82E+02	1.36E+09	1.23E+04	1		Benzene	71-43-2	1.3E+01		1.3E+00	1.2E+00	3.1E+02			1.1E+02	8.2E+01	
2.30E-01	C	3																								

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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	key	IUR (ug/m ³) ¹	key	RD ₅₀ (mg/kg-day)	key	RIC ₁ (mg/m ³)	key	vo	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS ₀	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)	
7.90E-03	I	1.10E-06		2.00E-02		V				9.15E+02	1.36E+09	9.70E+03	1		Bromoform	75-25-2	8.8E+01		2.5E+01	1.9E+01	1.6E+03				1.6E+03
				1.40E-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	
				5.00E-03	H	V				1.36E+09	1.24E+05	1		Bromophos	2104-96-3						3.9E+02			3.9E+02	
		3.70E-06	C	1.00E-01	A	V				9.66E+02	1.36E+09	2.14E+03	1		Bromopropane, 1-	106-94-5			1.6E+00	1.6E+00			2.2E+02	2.2E+02	
1.03E-01	O			1.50E-02	O					1.36E+09	4.74E+05	1	0.1	Bromoxynil	1689-84-5	6.7E+00	2.4E+01		5.3E+00	1.2E+03	4.9E+03			9.5E+02	
1.03E-01	O			1.50E-02	O					1.36E+09	8.66E+02	1		Bromoxynil Octanoate	1689-99-2	6.7E+00			6.7E+00	1.2E+03				1.2E+03	
6.00E-01	C	3.00E-05	I	2.00E-03	I	V				6.67E+02	1.36E+09	6.66E+02	1		Butadiene, 1,3-	106-99-0	1.2E+00		8.1E-02	7.6E-02				1.8E+00	
				1.00E-01	I	V				7.64E+03	1.36E+09	3.00E+04	1		Butanol, n-	71-36-3								7.8E+03	
5.00E-04	I			4.00E-01	I	V				1.36E+09	2.87E+04	1		Butyl Alcohol, 1-	75-65-0	1.4E+03			1.4E+03	7.8E+03			3.1E+04	1.5E+05	
				2.00E+00	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	
				5.00E-02	I	V				1.36E+09	8.63E+04	1		Butylate	2008-41-5					3.9E+03				3.9E+03	
2.00E-04	C	5.70E-08	C	3.00E-01	P					1.36E+09		1	0.1	Butylated hydroxyanisole	25013-16-5	3.5E+03	1.2E+04	6.7E+07	2.7E+03					1.9E+04	
3.60E-03	P			5.00E-02	P	V				1.08E+02	1.36E+09	8.14E+03	1		Butylated hydroxytoluene	128-37-0	1.9E+02	6.9E+02	1.5E+02			2.3E+04	9.9E+04	1.9E+04	
				1.00E-01	X	V				1.45E+02	1.36E+09	7.35E+03	1		Butylbenzene, n-	104-51-8					3.9E+03			3.9E+03	
				1.00E-01	X	V				1.83E+02	1.36E+09	7.36E+03	1		Butylbenzene, sec-	135-98-8					7.8E+03			7.8E+03	
				1.00E-01	X	V				1.83E+02	1.36E+09	7.36E+03	1		Butylbenzene, tert-	98-06-6					7.8E+03			7.8E+03	
1.80E-03	I	1.00E-04	A	1.00E-05	A					1.36E+09		1	0.1	Caecidic Acid	75-60-5					1.6E+03	6.6E+03		1.4E+04	1.3E+03	
1.80E-03	I	1.00E-04	A	1.00E-05	A					1.36E+09	0.025	0.001		Cadmium (Diet)	7440-43-9			2.1E+03	2.1E+03	7.8E+00	8.2E+01			7.1E+00	
				1.80E-03	I	2.20E-03	C			1.36E+09	0.05	0.001		Cadmium (Water)	7440-43-9										
				5.00E-01	I	V				1.36E+09		1	0.1	Caprolactam	105-60-2					3.9E+04	1.6E+05		3.1E+06	3.1E+04	
1.50E-01	C	4.30E-05	C	2.00E-03	I	V				1.36E+09		1	0.1	Captafol	2425-06-1	4.6E+00	1.6E+01	8.9E+04	3.6E+00	1.6E+02	6.6E+02		4.3E+04	1.3E+02	
2.30E-03	C	6.00E-07	C	1.30E-01	I	V				1.36E+09		1	0.1	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	
				1.00E-01	I	V				1.36E+09		1	0.1	Carbaryl	63-25-2					7.8E+03	3.3E+04			6.3E+03	
				5.00E-03	I	V				1.36E+09		1	0.1	Carbafuran	1563-66-2					3.9E+02	1.6E+03			3.2E+02	
				1.00E-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	
7.00E-02	I	6.00E-06	I	4.00E-03	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	
				1.00E-02	I	V				5.89E+03	1.36E+09	6.46E+02	1		Carbonyl Sulfide	463-58-1					7.8E+03		6.7E+01	6.3E+02	
				1.00E-01	I	V				1.36E+09		1	0.1	Carbosulfan	55285-14-8					7.8E+02	3.3E+03			6.3E+02	
				1.00E-01	I	V				1.36E+09		1	0.1	Carboxin	5234-68-4					7.8E+03	3.3E+04			6.3E+03	
				1.00E-01	I	V				1.36E+09	1.45E+05	1		Ceric oxide	1306-38-3					7.8E+03		1.3E+06		1.3E+06	
				1.50E-02	I	V				1.36E+09		1	0.1	Chloral Hydrate	302-17-0					7.8E+03	4.9E+03			9.8E+03	
				5.00E-01	I	V				1.36E+09		1	0.1	Chloramben	133-90-4					1.2E+03		4.9E+03		7.5E+02	
4.03E-01	H			5.00E-04	G	V				1.36E+09		1	0.1	Chloramines, Organic	E701235										
				5.00E-04	G	V				1.36E+09		1	0.1	Chloranil	118-75-2	1.7E+00	6.1E+00		1.3E+00						
				5.00E-04	G	V				1.36E+09	1.49E+06	1	0.04	Chlordane (alpha)	5103-71-9					3.9E+01	4.1E+02			3.6E+01	
				5.00E-04	G	V				1.36E+09	1.49E+06	1	0.04	Chlordane (gamma)	5103-74-2					3.9E+01	4.1E+02			3.6E+01	
3.50E-01	I	1.00E-04	I	7.00E-04	I	V				1.36E+09	1.53E+06	1	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	
1.00E+01	I	4.60E-03	C	3.00E-04	I	V				1.36E+09		1	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	
				7.00E-04	A	V				1.36E+09		1	0.1	Chlorfenvinphos	470-90-6					5.5E+01	2.3E+02			4.4E+01	
				9.00E-02	O	V				1.36E+09		1	0.1	Chlorimuron, Ethyl-	90982-32-4					7.0E+03	3.0E+04			5.7E+03	
				1.00E-01	I	V				2.78E+03	1.36E+09	1.22E+03	1		Chlorine	7782-50-5					7.8E+03		1.8E-01	1.8E-01	
				3.00E-02	I	V				1.36E+09		1		Chlorine Dioxide	10049-04-4					2.3E+03	2.8E+05			2.3E+03	
				3.00E-02	I	V				1.36E+09		1		Chlorite (Sodium Salt)	7758-19-2					2.3E+03				2.3E+03	
4.60E-01	H	3.00E-04	I	2.00E-02	H	V				1.15E+03	1.36E+09	1.03E+03	1		Chloro-1,1-difluoroethane, 1-	75-68-3					1.6E+03		5.4E+04	5.4E+04	
1.00E-01	P	7.70E-05	C	3.00E-03	X					7.86E+02	1.36E+09	1.08E+03	1	0.1	Chloro-1,3-butadiene, 2- (Chloroprene)	128-99-8	1.5E+00	5.4E+00	3.6E-03	3.6E-03	1.6E+03		5.4E+04		
2.70E-01	X			3.50E-03	C	V				1.36E+09		1	0.1	Chloro-2-methylaniline HCl, 4-	3165-93-3	7.0E+00	2.5E+01	5.0E+04	5.4E+00	2.3E+02	9.9E+02			1.9E+02	
				3.50E-03	C	V				1.36E+09	1.62E+04	1		Chloroacetalddehyde, 2-	107-20-0	2.6E+00			2.6E+00	2.3E+02	9.9E+02			1.9E+02	
				3.00E-05	I	V				1.36E+09		1	0.1	Chloroacetic Acid	79-11-8					2.7E+02	1.2E+03			2.2E+02	
2.00E-01	P			5.00E-04	P	V				1.36E+09		1	0.1	Chloroacetophenone, 2-	532-27-4	3.5E+00	1.2E+01		2.7E+00	3.9E+01	1.6E+02		4.3E+04	4.3E+04	
				2.00E-02	I	V				1.36E+09	6.45E+03	1		Chloroaniline, p-	106-47-8					1.6E+03				2.8E+02	
				1.00E-01	X	V				1.36E+09		1	0.1	Chlorobenzene	108-90-7					7.8E+03	3.3E+04			6.3E+03	
1.10E-01	C	3.10E-05	C	2.00E-02	I	V				1.36E+09		1	0.1	Chlorobenzene sulfonic acid, p-	98-66-8	6.3E+00	2.2E+01	1.2E+05	4.9E+00	1.6E+03	6.6E+03			1.3E+03	
				3.00E-02	X	V				1.36E+09		1	0.1	Chlorobenzoate	510-15-6					2.3E+03	9.9E+03			1.9E+03	
				8.60E-06	C	3.00E-01	P	V		2.90E+02	1.36E+09	6.76E+0													

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	RD ₅₀ (mg/kg-day)	k e y	RIC ₁ (mg/m ³)	k e y	v o l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS ₀	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 THQ=1 (mg/kg)	
				4.00E-02	H						1.36E+09			1	Copper	7440-50-8					3.1E+03			3.1E+03	
				5.00E-02	I	6.00E-01	C				1.36E+09		0.1	Cresol, m-	108-39-4					3.9E+03	1.6E+04	8.5E+08	3.2E+03		
				5.00E-02	I	6.00E-01	C				1.36E+09		0.1	Cresol, o-	95-48-7					3.9E+03	1.6E+04	8.5E+08	3.2E+03		
				2.00E-02	P	6.00E-01	C				1.36E+09		0.1	Cresol, p-	106-44-5					1.6E+03	6.6E+03	8.5E+08	1.3E+03		
				1.00E-01	A						1.36E+09		0.1	Cresol, p-chloro-m-	59-50-7					7.8E+03	3.3E+04	8.5E+08	6.3E+03		
1.90E+00	H			1.00E-01	A	6.00E-01	C				1.36E+09		0.1	Cresols	1319-77-3				3.7E-01	7.8E+03	3.3E+04	8.5E+08	6.3E+03		
				1.00E-03	P				1.66E+04	1.36E+09	1.89E+04		1	Crotonaldehyde, trans-	123-73-9				3.7E-01	7.8E+03			7.8E+01	7.8E+01	
				1.00E-01	I	4.00E-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		
2.20E-01	C	6.30E-05	C								1.36E+09		0.1	Cupferron	135-20-6	3.2E+00	1.1E+01	6.1E+04	2.9E+00						
8.40E-01	H			2.00E-03	H						1.36E+09		0.1	Cyanazine	21725-46-2	8.3E-01	2.9E+00		6.5E-01	1.6E+02	6.6E+02		1.3E+02		
											1.36E+09		1	Cyanides											
				1.00E-03	I	9.00E-03	C				1.36E+09		1	-Calcium Cyanide	592-01-8					7.8E+01		1.3E+07	7.8E+01		
				5.00E-03	I						1.36E+09		1	-Copper Cyanide	544-92-3					3.9E+02			3.9E+02		
				6.00E-04	I	8.00E-04	G	V	9.54E+05	1.36E+09	5.89E+04		1	-Cyanide (CN-)	57-12-5					4.7E+01		4.9E+01	2.4E+01		
				1.00E-03	I						1.36E+09		1	-Cyanogen	460-19-5					7.8E+01			7.8E+01		
				9.00E-02	I						1.36E+09		1	-Cyanogen Bromide	506-68-3					7.0E+03			7.0E+03		
				5.00E-02	I						1.36E+09		1	-Cyanogen Chloride	506-77-4					3.9E+03			3.9E+03		
				6.00E-04	I	8.00E-04	I	V	1.00E+07	1.36E+09	5.22E+04		1	-Hydrocyan Cyanide	74-90-8					4.7E+01		4.4E+01	2.3E+01		
				2.00E-03	I	9.00E-03	C				1.36E+09		0.04	-Potassium Cyanide	151-50-8					1.6E+02		1.3E+07	1.6E+02		
				5.00E-03	I						1.36E+09		0.04	-Potassium Silver Cyanide	506-61-6					3.9E+02			3.9E+02		
				1.00E-01	I						1.36E+09		1	-Silver Cyanide	506-64-9					7.8E+03			7.8E+03		
				1.00E-03	I	9.00E-03	C				1.36E+09		1	-Sodium Cyanide	143-33-9					7.8E+01		1.3E+07	7.8E+01		
				5.00E-02	I						1.36E+09		1	-Zinc Cyanide	557-21-1					3.9E+03			3.9E+03		
2.00E-02	X			2.00E-02	X	6.00E+00	I	V	1.17E+02	1.36E+09	1.04E+03		0.1	Cyclohexane	110-82-7				2.7E+01	1.6E+03	6.6E+03	6.5E+03	6.5E+03		
				5.00E+00	I	7.00E-01	P	V	5.11E+03	1.36E+09	4.17E+04		1	Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	3.5E+01	1.2E+02			3.9E+05	6.6E+03	3.0E+04	1.3E+03		
											1.36E+09		1	Cyclohexanone	108-94-1								2.8E+04		
				5.00E-03	P	1.00E+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		
				2.00E-01	I				2.93E+05	1.36E+09	7.46E+04		1	Cyclohexylamine	108-91-8					1.6E+04			1.6E+04		
				2.50E-02	I						1.36E+09		0.1	Cyfluthrin	68359-37-5					2.0E+03	8.2E+03		1.6E+03		
				5.00E-01	O						1.36E+09		0.1	Cymazine	66215-27-8					3.9E+04	1.6E+05		3.2E+04		
				3.00E-02	I						1.36E+09		0.1	Dalapon	75-99-0					2.3E+03	9.9E+03		1.9E+03		
1.80E-02	C	5.10E-06	C								1.36E+09		0.1	Daminozide	1596-84-5	3.9E+01	1.4E+02	7.5E+05	3.0E+01	1.2E+04	4.9E+04		9.5E+03		
7.00E-04	I			7.00E-03	I						1.36E+09		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		
				4.00E-05	I						1.36E+09		0.1	Demeton	8065-48-3					3.1E+00		1.3E+01	2.5E+00		
1.20E-03	I			6.00E-01	I						1.36E+09		0.1	Di(2-ethylhexyl)adipate	103-23-1				5.8E+02	4.7E+04	2.0E+05		3.8E+04		
6.10E-02	H			7.00E-04	A						1.36E+09		0.1	Diallate	2303-16-4	1.1E+01	4.1E+01		8.9E+00						
				3.00E-04	P				9.79E+02	1.36E+09	3.20E+04		1	Diazinon	333-41-5					5.5E+01	2.3E+02		4.4E+01		
8.00E-01	P	6.00E-03	P	2.00E-04	P	2.00E-04	I	V	M	9.79E+02	1.36E+09	3.20E+04		1	Dibromo-3-chloropropane, 1,2-	96-12-8	1.9E-01		5.4E-03	5.3E-03	1.6E+01		6.7E+00	4.7E+00	
2.50E-01	C			3.00E-04	C						1.36E+09		0.1	Dibromoacetic acid	631-64-1	2.8E+00	9.9E+00		2.2E+00	2.3E+01	9.9E+01		1.9E+01		
				4.00E-04	X			V	1.59E+02	1.36E+09	1.93E+04		1	Dibromobenzene, 1,3-	108-36-1					3.1E+01			3.1E+01		
				1.00E-02	I						1.36E+09		1	Dibromobenzene, 1,4-	106-37-6					7.8E+02			7.8E+02		
8.40E-02	I			2.00E-02	I			V	8.02E+02	1.36E+09	7.95E+03		1	Dibromochloromethane	124-48-1	8.3E+00			8.3E+00	1.6E+03			1.6E+03		
2.00E+00	I	6.00E-04	I	9.00E-03	I	9.00E-03	I	V	1.34E+03	1.36E+09	8.64E+03		1	Dibromomethane, 1,2-	106-93-4	3.5E-01		4.0E-02	3.6E-02	7.0E+02		8.1E+01	7.3E+01		
				1.00E-02	I	4.00E-03	X	V	2.82E+03	1.36E+09	5.64E+03		1	Dibromomethane (Methylene Bromide)	74-95-3					1.6E+03		2.4E+01	2.4E+01		
				3.00E-04	P						1.36E+09		0.1	Dibutyltin Compounds	E1790661					2.3E+01	9.9E+01		1.9E+01		
				3.00E-02	I						1.36E+09		0.1	Dicamba	1918-00-9					2.3E+03	9.9E+03		1.9E+03		
											1.36E+09		0.1	Dichloramine	3400-09-7										
4.20E-03	P							V	5.54E+02	1.36E+09	3.21E+03		1	Dichloro-2-butene, 1,4-	764-41-0			2.1E-03	2.1E-03						
4.20E-03	P							V	5.19E+02	1.36E+09	1.11E+04		1	Dichloro-2-butene, cis-1,4-	1476-11-5			7.4E-03	7.4E-03						
4.20E-03	P							V	7.60E+02	1.36E+09	1.11E+04		1	Dichloro-2-butene, trans-1,4-	110-57-6			7.4E-03	7.4E-03						
5.00E-02	I			4.00E-03	I						1.36E+09		0.1	Dichloroacetic Acid	79-43-6	1.4E+01	4.9E+01		1.1E+01	3.1E+02	1.3E+03		2.5E+02		
				9.00E-02	I	2.00E-01	H	V	3.76E+02	1.36E+09	1.17E+04		1	Dichlorobenzene, 1,2-	95-50-1					7.0E+03		2.4E+03	1.8E+03		
5.40E-03	C	1.10E-05	C	7.00E-02	A	8.00E-01	I	V		1.36E+09	1.04E+04		1	Dichlorobenzene, 1,4-	106-46-7	1.3E+02		2.7E+00	2.6E+00	5.5E+03		8.7E+03	3.4E+03		
4.50E-01	I	3.40E-04	C								1.36E+09		0.1	Dichlorobenzidine, 3,3'-	91-94-1	1.5E+00	5.5E+00	1.1E+04	1.2E+00						
				9.00E-03	X						1.36E+09		0.1	Dichlorobenzophenone, 4,4'-	90-98-2					7.0E+02	3.0E+03		5.7E+02		
				2.00E-01	I	1.00E-01	X	V	8.45E+02	1.36E+09	8.41E+02		1	Dichlorodifluoromethane	75-71-8					1.6E+04		8.8E+01	8.7E+01		
2.40E-01	I	6.90E-05	C	5.00E-04	A						1.36E+09		0.1	Dichlorodiphenyldichloroethane, 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		
3.40E-01	I																								

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; T = ATSDR DRAFT; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; R = ORD; N = WI; W = TEF applied; E = RPF applied; G = see user's guide; 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	RD ₅₀ (mg/kg-day)	k e y	RIC ₁ (mg/m ³) ¹	k e y	v o l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS ₀	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)	
4.40E-02	C	1.30E-05	C			3.00E+01	X	V		6.91E+02	1.36E+09	7.58E+02	1		Diffluoropropane, 2,2-Dihydroasafrole Diisopropyl Ether	420-45-1 94-58-6 108-20-3	1.6E+01		2.7E+01	9.9E+00			2.4E+04	2.4E+04	
				8.00E-02	I	7.00E-01	P	V		2.26E+03	1.36E+09	3.81E+04	1		Diisopropyl Methylphosphonate	1445-75-6					6.3E+03			6.3E+03	
				2.18E-02	O					5.30E+02	1.36E+09	3.06E+03	1	0.1	Dimethipin	55290-64-7					1.7E+03	7.2E+03		1.4E+03	
				2.20E-03	O						1.36E+09		1	0.1	Dimethoate	60-51-5					1.7E+02	7.3E+02		1.4E+02	
1.60E+00	P			2.00E-02	X	2.00E-04	X	V	M	5.35E+03	1.36E+09	2.97E+03	1	0.1	Dimethoxybenzidine, 3,3'-Dimethyl Sulfide	119-90-4 75-18-3	9.6E-02	3.7E-01		7.6E-02	1.6E+03		6.2E-01	6.2E-01	
1.70E-03	P			6.00E-02	P						1.36E+09		1	0.1	Dimethyl methylphosphonate	756-79-6	4.1E+02	1.5E+03	3.2E+02		4.7E+03	2.0E+04		3.8E+03	
4.60E+00	C	1.30E-03	C								1.36E+09		1	0.1	Dimethylamino azobenzene [p-]	60-11-7	1.5E-01	5.4E-01	2.9E+03	1.2E-01					
5.80E-01	H										1.36E+09		1	0.1	Dimethylamine HCl, 2,4-Dimethylaniline, 2,4-Dimethylaniline, 2,4-Dimethylaniline, N,N-Dimethylbenzidine, 3,3'-Dimethylformamide	121-69-7 119-93-7 68-12-2	2.6E+01	1.2E+01	2.6E+01	1.1E-02	4.7E+03		4.0E+03	2.6E+03	
2.00E-01	P			2.00E-03	X						1.36E+09		1	0.1	Dimethylhydrazine, 1,1-Dimethylhydrazine, 1,2-Dimethylphenol, 2,4-Dimethylphenol, 2,5-Dimethylphenol, 3,4-Dimethylvinylchloride	95-68-1 540-73-8 105-67-9 576-26-1	3.5E+00	1.2E+01	2.7E+00		1.6E+02	6.6E+02		1.3E+02	
2.70E-02	P			2.00E-03	I			V	M	8.30E+02	1.36E+09	3.13E+04	1	0.1	Dinitro-p-cresol, 4,6-Dinitro-o-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, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	2.6E+01	1.2E+01	2.6E+01	1.1E-02	7.8E+00		5.8E-02	1.6E+02	
1.10E+01	P			1.00E-01	P	3.00E-02	I	V		1.06E+05	1.36E+09	1.28E+05	1	0.1	Dimethylhydrazine, 1,1-Dimethylhydrazine, 1,2-Dimethylphenol, 2,4-Dimethylphenol, 2,5-Dimethylphenol, 3,4-Dimethylvinylchloride	57-14-7 540-73-8 105-67-9 576-26-1	1.3E-03		2.9E-03	8.8E-04	7.8E+00		5.8E-02	5.7E-02	
5.50E+02	C	1.60E-01	C	2.00E-02	I						1.36E+09		1	0.1	Dimethylhydrazine, 1,1-Dimethylhydrazine, 1,2-Dimethylphenol, 2,4-Dimethylphenol, 2,5-Dimethylphenol, 3,4-Dimethylvinylchloride	57-14-7 540-73-8 105-67-9 576-26-1	1.3E-03		2.9E-03	8.8E-04	1.6E+03	6.6E+03		3.8E+01	
				6.00E-04	I						1.36E+09		1	0.1	Dimethylphenol, 2,5-Dimethylphenol, 3,4-Dimethylvinylchloride	95-65-9	4.7E+01				4.7E+01	2.0E+02		1.3E+01	
4.50E-02	C	1.30E-05	C	1.00E-03	I			V		4.73E+02	1.36E+09	5.48E+03	1	0.1	Dinitrochlorobenzene, 1,2-Dinitro-p-cresol, 4,6-Dinitro-o-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, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	1.5E+01		1.2E+00	1.1E+00	7.8E+00		5.8E-02	1.6E+02	
				8.00E-05	X						1.36E+09		1	0.1	Dinitrochlorobenzene, 1,2-Dinitro-p-cresol, 4,6-Dinitro-o-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, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	1.5E+01		1.2E+00	1.1E+00	6.3E+00	2.6E+01		5.1E+00	
				2.00E-03	I						1.36E+09		1	0.1	Dinitrochlorobenzene, 1,2-Dinitro-p-cresol, 4,6-Dinitro-o-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, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	1.5E+01		1.2E+00	1.1E+00	1.6E+02	6.6E+02		2.8E+06	
				4.00E-04	X	2.00E-03	X				1.36E+09		1	0.1	Dinitrochlorobenzene, 1,2-Dinitro-p-cresol, 4,6-Dinitro-o-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, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	1.5E+01		1.2E+00	1.1E+00	3.1E+01	1.3E+02	2.8E+06	2.5E+01	
				1.00E-04	P						1.36E+09		1	0.1	Dinitrochlorobenzene, 1,2-Dinitro-p-cresol, 4,6-Dinitro-o-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, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	1.5E+01		1.2E+00	1.1E+00	7.8E+00	3.3E+01		6.3E+00	
				1.00E-04	I						1.36E+09		1	0.1	Dinitrochlorobenzene, 1,2-Dinitro-p-cresol, 4,6-Dinitro-o-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, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	1.5E+01		1.2E+00	1.1E+00	7.8E+00	3.3E+01		6.3E+00	
				1.00E-04	P						1.36E+09		1	0.1	Dinitrochlorobenzene, 1,2-Dinitro-p-cresol, 4,6-Dinitro-o-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, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	1.5E+01		1.2E+00	1.1E+00	7.8E+00	3.3E+01		6.3E+00	
				2.00E-03	I						1.36E+09		1	0.1	Dinitrochlorobenzene, 1,2-Dinitro-p-cresol, 4,6-Dinitro-o-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, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	1.5E+01		1.2E+00	1.1E+00	1.6E+02	6.6E+02		1.3E+02	
6.80E-01	I			3.10E-01	C	8.90E-05	C			2.00E-03	I				Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	1.0E+00	3.6E+00		8.0E-01	1.6E+02	6.5E+02		1.3E+02	
1.50E+00	P			3.00E-04	X						1.36E+09		1	0.099	Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	4.6E-01	1.7E+00		3.6E-01	2.3E+01	1.0E+02		1.9E+01	
				1.00E-04	X						1.36E+09		1	0.006	Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	4.6E-01	1.7E+00		3.6E-01	7.8E+00	5.5E+02		7.7E+00	
				1.00E-04	X						1.36E+09		1	0.009	Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	4.6E-01	1.7E+00		3.6E-01	7.8E+00	3.7E+02		7.7E+00	
4.50E-01	X			9.00E-04	X						1.36E+09		1	0.1	Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	1.5E+00	5.5E+00		1.2E+00	7.0E+01	3.0E+02		5.7E+01	
				1.00E-03	I						1.36E+09		1	0.1	Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 3,4-Dinitrotoluene, 4-Amino-2,6-Dinitrotoluene, Technical grade Dinoseb	513-37-1 534-52-1 131-89-5 618-87-1 528-29-0 99-65-0 100-25-4 51-28-5 E1615210 121-14-2	1.5E+00	5.5E+00		1.2E+00	7.8E+01	3.3E+02		6.3E+01	
1.00E-01	I	5.00E-06	I	3.00E-02	I	3.00E-02	I	V		1.16E+05	1.36E+09	3.96E+04	1		Dioxane, 1,4-Dioxins	123-91-1	7.0E+00		2.2E+01	5.3E+00	2.3E+03		1.2E+03	8.1E+02	
6.20E+03	I	1.30E+00	I								1.36E+09		1	0.03	-Hexachlorodibenzo-p-dioxin, Mixture	34465-46-8	1.1E-04	1.3E-03	2.9E+00	1.0E-04	5.5E-05	7.7E-04	8.2E-02	5.1E-05	
1.30E+05	C	3.80E+01	C	7.00E-10	I	4.00E-08	C	V			1.36E+09	1.96E+06	1	0.03	-TCDD, 2,3,7,8-Diphenamid	1746-01-6 957-51-7	5.3E-06	6.3E-05	1.4E-04	4.8E-06	2.3E+03	9.9E+03		1.9E+03	
				3.00E-02	I						1.36E+09		1	0.1	Diphenamid	957-51-7					2.3E+03	9.9E+03		1.9E+03	
						4.00E-04	X	V			1.36E+09	8.06E+04	1		Diphenyl Ether	101-84-8							3.4E+01	3.4E+01	
				8.00E-04	X						1.36E+09		1	0.1	Diphenyl Sulfone	127-63-9					6.3E+01	2.6E+02		5.1E+01	
8.00E-01	I	2.20E-04	I	1.00E-01	O						1.36E+09		1	0.1	Diphenylamine	122-39-4	8.7E-01	3.1E+00	1.7E+04	6.8E-01	7.8E+03	3.3E+04		6.3E+03	
				2.20E-03	I						1.36E+09		1	0.1	Diphenylhydrazine, 1,2-Dioxal	2764-72-9					7.8E+03	3.3E+04		6.3E+03	
7.40E+00	C	2.10E-03	C								1.36E+09														

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	RD ₅₀ (mg/kg-day)	k e y	RIC ₁ (mg/m ³)	k e y	v o l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS ₀	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)		
				1.30E-02	I						1.36E+09			1	0.1	Fluometuron	2164-17-2					1.0E+03	4.3E+03		8.2E+02	
				4.00E-02	C	1.30E-02	C				1.36E+09			1		Fluoride	16984-48-8				3.1E+03		1.8E+07	3.1E+03		
				6.00E-02	I	1.30E-02	C				1.36E+09			1		Fluorine (Soluble Fluoride)	7782-41-4				4.7E+03		1.8E+07	4.7E+03		
				8.00E-02	I						1.36E+09			1	0.1	Fluoride	59756-60-4				6.3E+03	2.6E+04		5.1E+03		
				4.00E-02	O						1.36E+09			1	0.1	Flurprimidol	56425-91-3				3.1E+03	1.3E+04		2.5E+03		
				2.00E-03	O						1.36E+09			1	0.1	Flusilazole	85509-19-9				1.6E+02	6.6E+02		1.3E+02		
				5.00E-01	O						1.36E+09			1	0.1	Flutolanil	96332-98-5				3.9E+04	1.6E+05		3.2E+04		
				1.00E-02	I						1.36E+09			1	0.1	Fluvalinate	89409-94-5				7.8E+02	3.3E+03		6.3E+02		
				9.00E-02	O						1.36E+09			1	0.1	Folpet	133-07-3				7.0E+03	3.0E+04		5.7E+03		
				1.00E-02	O						1.36E+09			1	0.1	Fomesafen	72178-02-0				7.8E+02	3.3E+03		6.3E+02		
				2.00E-03	I						1.36E+09			1	0.1	Fonofos	944-22-9				1.6E+02	6.6E+02		1.3E+02		
	2.10E-02	C	7.40E-06	I	2.00E-01	I	7.00E-03	I	V	M	4.24E+04	1.36E+09	7.77E+04	1		Formaldehyde	50-00-0	7.3E+00		1.1E+01	4.3E+00	1.6E+04		5.5E+02		
				9.00E-01	P	3.00E-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		
				2.50E+00	O						1.36E+09			1	0.1	Fosetyl-AL	39148-24-8				2.0E+05	8.2E+05		1.6E+05		
				1.00E-03	X			V			1.36E+09	1.56E+05	1		-Dibenzofuran	132-64-9					7.8E+01			7.8E+01		
				1.00E-03	I			V		6.22E+03	1.36E+09	2.62E+03	1		-Furan	110-00-9					7.8E+01			7.8E+01		
	3.80E+00	H		9.00E-01	I	2.00E+00	I	V		1.65E+05	1.36E+09	1.20E+04	1		-Tetrahydrofuran	109-99-9					7.0E+04		2.5E+04	1.8E+04		
				3.00E-03	I	5.00E-02	H	V			1.36E+09	4.86E+04	1		Furazolidone	67-45-8	1.8E-01	6.5E-01		1.4E-01					2.1E+02	
				1.50E+00	C	4.30E-04	C				1.01E+04	1.36E+09	4.86E+04	1	0.1	Furfural	98-01-1					2.3E+02			2.1E+02	
	3.00E-02	I	8.60E-06	C							1.36E+09			1	0.1	Furium	531-82-8	4.6E-01	1.6E+00	8.9E+03	3.6E-01				2.1E+02	
				6.00E-03	O						1.36E+09			1	0.1	Furmecyclo	80568-05-0	2.3E-01	8.2E+01	4.4E+05	1.8E+01				2.1E+02	
				1.00E-01	A	8.00E-05	C				1.36E+09			1	0.1	Glufosinate, Ammonium	77182-82-2					4.7E+02	2.0E+03		3.8E+02	
				4.00E-04	A	1.00E-03	X	V		1.06E+05	1.36E+09	8.43E+04	1		Glutaraldehyde	111-30-8					7.8E+03	3.3E+04	1.1E+05	6.0E+03		
				1.00E-01	I						1.36E+09			1	0.1	Glycidaldehyde	765-34-4				3.1E+01		8.8E+01	2.3E+01		
				1.00E-02	X			V			1.36E+09	1.45E+05	1		Guanidine	113-00-8					7.8E+02			7.8E+02		
				2.00E-02	P						1.36E+09			1	0.1	Guanidine Chloride	50-01-1				1.6E+03	6.6E+03		1.3E+03		
				3.00E-02	X						1.36E+09			1	0.1	Guanidine Nitrate	506-93-4				2.3E+03	9.9E+03		1.9E+03		
	4.50E+00	I	1.30E-03	I	1.00E-04	A		V			1.36E+09	4.79E+05	1		Haloxypol, Methyl	69806-40-2	1.5E-01		1.0E+00	1.3E-01	3.9E+00	1.6E+01		3.2E+00		
	9.10E+00	I	2.60E-03	I	1.30E-05	I					1.36E+09	4.79E+05	1		Heptachlor	76-44-8	7.6E-02				7.8E+00			7.8E+00		
				3.00E-04	X	4.00E-01	P	V		5.79E+01	1.36E+09	8.95E+02	1		Heptachlor Epoxide	1024-87-3					1.0E+00			1.0E+00		
				2.00E-03	I			V			1.36E+09	7.80E+03	1		Heptanal, n-	111-71-7			9.1E-01	7.0E-02				2.4E+01		
				2.00E-04	X						1.36E+09	3.80E+05	1		Heptane, n-	142-82-5					2.3E+01			2.2E+01		
	1.60E+00	I	4.60E-04	I	1.00E-05	P		V			1.36E+09	6.80E+04	1	0.1	Hexabromobenzene	87-82-1	4.3E-01		4.1E-01	2.1E-01	1.6E+02				1.6E+02	
	7.80E-02	I	2.20E-05	I	1.00E-03	P		V		1.68E+01	1.36E+09	1.08E+04	1		Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BDE-153)	86631-49-2					1.6E+01	6.6E+01		1.3E+01		
	6.30E+00	I	1.80E-03	I	9.00E-04	A					1.36E+09			1	0.1	Hexachlorobenzene	118-74-1	8.9E+00		1.4E+00	1.2E+00	7.8E-01			7.8E-01	
	1.80E+00	I	5.30E-04	I							1.36E+09			1	0.1	Hexachlorobutadiene	87-88-3	3.9E-01	3.9E-01	2.1E+03	8.6E-02	7.0E+01	3.0E+02		5.7E+01	
				6.00E-08	X						1.36E+09			1	0.1	Hexachlorocyclohexane, Alpha-	319-84-6				3.9E-01	1.4E+00	7.2E+03	3.0E-01		
	1.10E+00	C	3.10E-04	C	8.00E-07	A					1.36E+09			0.04	Hexachlorocyclohexane, Beta-	319-86-8					4.7E+03	2.0E-02		3.8E-03		
	1.80E+00	I	5.10E-04	I							1.36E+09			1	0.1	Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	6.3E-01	5.6E+00	1.2E+04	5.7E-01	6.3E-02	6.6E-01		5.7E-02	
				6.00E-03	I	2.00E-04	I	V		1.57E+01	1.36E+09	8.51E+03	1		Hexachlorocyclohexane, Technical	608-73-1	3.9E-01	1.4E+00	7.5E+03	3.0E-01	6.3E-02	6.6E-01		5.7E-02		
	4.00E-02	I	1.10E-05	C	3.00E-02	I	V				1.36E+09	8.01E+03	1		Hexachlorocyclopentadiene	77-47-4	1.7E+01		2.0E+00	1.8E+00	4.7E+02				1.8E+00	
				3.00E-04	I						1.36E+09			1	0.1	Hexachlorocyclopentadiene	67-72-1				5.5E+01	3.3E+04		4.5E+01		
	8.00E-02	I		4.00E-03	I						1.36E+09			0.015	Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	8.7E+00	2.1E+02		8.3E+00	2.3E+01	9.9E+01		1.9E+01		
				1.00E-05	I	V				3.39E+03	1.36E+09	3.00E+05	1		0.1	Hexamethylene Diisocyanate, 1,6-	822-06-0				3.1E+02	8.8E+03		3.1E+00		
				4.00E-04	C						1.36E+09			1	0.1	Hexamethylene diisocyanate biuret	4035-89-6							5.7E+05		
				4.00E-04	P						1.36E+09			1	0.1	Hexamethylene diisocyanate isocyanurate	3779-63-3							5.7E+05		
				2.00E-07	X	6.00E-01	P	V		1.41E+02	1.36E+09	8.29E+02	1		Hexamethylphosphoramide	680-31-9					3.1E+01	1.3E+02		2.5E+01		
				7.00E-01	I	V				1.41E+02	1.36E+09	8.29E+02	1		Hexane, Commercial	E5241997			1.2E+01	1.2E+01			5.2E+02		5.2E+02	
				2.00E+00	P						1.36E+09			1	0.1	Hexane, N-	110-54-3				1.6E+05	6.6E+05		6.1E+02		
	9.50E-03	P		7.00E-02	P	4.00E-04	P	V		2.74E+02	1.36E+09	3.62E+04	1		Hexanedioic Acid	124-04-9	7.3E+01			7.3E+01	5.5E+03		1.5E+01	1.5E+01		
				5.00E-03	I	3.00E-02	I	V		3.28E+03	1.36E+09	1.33E+04	1		Hexanone, 2-	591-78-6				3.9E+02		4.2E+02		2.0E+02		
				3.30E-02	I						1.36E+09			1	0.1	Hexazinone	51235-04-2				2.6E+03	1.1E+04		2.1E+03		
				2.50E-02	I						1.36E+09			1	0.1	Hexythiazox	78587-05-0				2.0E+03	8.2E+03		1.6E+03		
	3.00E+00	I	4.90E-03	I							1.36E+09			1	0.1	Hydramethylnon	67485-29-4				1.3E+03	5.6E+03		1.1E+03		
	3.00E+00	I	4.90E-03	I		3.00E-05	P	V		1.12E+05	1.36E+09	6.52E+04	1		Hydrazine	302-01-2	2.3									

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; T = ATSDR DRAFT; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; R = ORD; N = WI; W = TEF applied; E = RPF applied; G = see user's guide; 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	RD ₅₀ (mg/kg-day)	k _e y ⁻¹	RIC ₁ (mg/m ³)	k _e y ⁻¹	v ₀ l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS ₀	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)	
				1.87E-05	P						1.36E+09		1	Lanthanum Chloride Heptahydrate	10025-84-0					1.5E+00				1.5E+00	
				2.83E-05	P						1.36E+09		1	Lanthanum Chloride, Anhydrous	10099-58-8					2.2E+00				2.2E+00	
				1.60E-05	P						1.36E+09		1	Lanthanum Nitrate Hexahydrate	10277-43-7					1.3E+00				1.3E+00	
8.50E-03	C	1.20E-05	C								1.36E+09		1	Lead Compounds	7446-27-7	8.2E+01		3.2E+05	8.2E+01						
2.10E-01	C	8.00E-05	C								1.36E+09		1	Lead acetate	301-04-2	3.3E+00	1.2E+01	4.8E+04	2.6E+00						2.0E+02
											1.36E+09		1	Lead and Compounds	7439-92-1										1.0E+02
3.80E-02	C	1.10E-05	C								1.36E+09		1	Lead and Compounds (with other sources of lead present, see Guidance)	7439-92-1										
				1.00E-07	I		V			2.43E+00	1.36E+09	1.91E+03	1	Lead subacetate	1335-32-6	1.8E+01	6.5E+01	3.5E+05	1.4E+01						
				5.00E-06	P		V			3.83E+02	1.36E+09	2.55E+04	1	Tetraethyl Lead	78-00-2							7.8E-03			7.8E-03
				7.70E-03	O						1.36E+09		1	Lewisite	541-25-3							3.9E-01			3.9E-01
				2.00E-03	P						1.36E+09		1	Linuron	330-55-2							6.0E+02	2.5E+03		4.9E+02
				5.00E-04	I						1.36E+09		1	Lithium	7439-93-2							1.6E+02			1.6E+02
				4.40E-02	O						1.36E+09		1	MCPPA	94-74-6							3.9E+01	1.6E+02		3.2E+01
				1.00E-03	I						1.36E+09		1	MCPB	94-81-5							3.4E+03	1.5E+04		2.8E+03
				2.00E-02	I						1.36E+09		1	MCPP	93-65-2							7.8E+01	3.3E+02		6.3E+01
				1.00E-01	I	7.00E-04	C				1.36E+09		1	Malathion	121-75-5							1.6E+03	6.6E+03		1.3E+03
				5.00E-01	I						1.36E+09		1	Maleic Anhydride	108-31-6							7.8E+03	3.3E+04	9.9E+05	6.3E+03
				1.00E-04	P						1.36E+09		1	Maleic Hydrazide	123-33-1							3.9E+04	1.6E+05		3.2E+04
				3.00E-02	H						1.36E+09		1	Malonitrile	109-77-3							7.8E+00	3.3E+01		6.3E+00
				5.00E-03	I						1.36E+09		1	Mancozeb	8018-01-7							2.3E+03	9.9E+03		1.9E+03
				1.40E-01	I	5.00E-05	I				1.36E+09		1	Maneb	12427-38-2							3.9E+02	1.6E+03		3.2E+02
				2.40E-02	G	5.00E-05	I		0.04		1.36E+09		1	Manganese (Diet)	7439-96-5							1.9E+03		7.1E+04	1.8E+03
				9.00E-05	H						1.36E+09		1	Manganese (Non-diet)	7439-96-5							7.0E+00	3.0E+01		5.7E+00
				3.00E-02	I						1.36E+09		1	Mepiquat Chloride	24307-26-4							2.3E+03	9.9E+03		1.9E+03
1.10E-02	P			4.00E-03	P						1.36E+09		1	Mercaptobenzothiazole, 2-Mercury Compounds	149-30-4	6.3E+01	2.2E+02		4.9E+01		3.1E+02	1.3E+03			2.5E+02
				3.00E-04	I	3.00E-04	G		0.07		1.36E+09		1	Mercury Chloride (and other Mercury salts)	7487-94-7							2.3E+01	4.3E+05		2.3E+01
				1.00E-04	I		V			3.13E+00	1.36E+09	2.27E+04	1	Mercury (elemental)	7439-97-6							7.1E+00			7.1E+00
				8.00E-05	I						1.36E+09		1	Methyl Mercury	22967-92-6							7.8E+00			7.8E+00
				3.00E-05	I		V				1.36E+09	1.94E+06	1	Phenylmercuric Acetate	62-36-4							6.3E+00	2.6E+01		5.1E+00
				6.00E-02	I						1.36E+09		1	Mephos	159-50-5							2.3E+00			2.3E+00
				1.00E-04	I	3.00E-02	P V			4.58E+03	1.36E+09	6.79E+03	1	Metallaxyl	57837-19-1							4.7E+03	2.0E+04		3.8E+03
				5.00E-05	I						1.36E+09		1	Methacrylonitrile	126-98-7							7.8E+00	2.1E+02		7.5E+00
				2.00E+00	I	2.00E+01	I V			1.06E+05	1.36E+09	2.90E+04	1	Methamidophos	10265-92-6							3.9E+00	1.6E+01		3.2E+00
				1.50E-03	O						1.36E+09		1	Methanol	67-56-1							1.6E+05		6.1E+05	1.2E+05
				2.50E-02	I						1.36E+09		1	Methidathion	950-37-8							1.2E+02	4.9E+02		9.5E+01
4.90E-02	C			5.00E-03	I						1.36E+09		1	Methomyl	16752-77-5	1.4E+01	5.0E+01		1.1E+01		2.0E+03	8.2E+03			1.6E+03
				8.00E-03	P	1.00E-03	P V			1.15E+05	1.36E+09	1.24E+05	1	Methoxyethanol Acetate, 2-Methoxyethanol, 2-Methyl Acetate	110-49-6						6.3E+02		1.3E+02		1.1E+02
				5.00E-03	P	7.00E-03	P V			1.06E+05	1.36E+09	1.01E+05	1	Methyl Acrylate	109-86-4							3.9E+02	7.4E+02		2.6E+02
				1.00E+00	X					2.90E+04	1.36E+09	8.12E+03	1	Methyl Acrylate	79-20-9							7.8E+04			7.8E+04
				6.00E-01	I	5.00E+00	I V			6.75E+03	1.36E+09	6.97E+03	1	Methyl Ethyl Ketone (2-Butanone)	96-33-3							7.8E+00			1.5E+02
				1.00E-03	X	1.00E-03	P	2.00E-05	X V	1.80E+05	1.36E+09	5.04E+04	1	Methyl Hydrazine	78-93-3							4.7E+04			6.4E+04
				3.00E+00	I V					3.36E+03	1.36E+09	1.06E+04	1	Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1			1.4E-01	1.4E-01		7.8E+01			1.1E+00	1.0E+00
				1.00E-03	C V					1.01E+04	1.36E+09	4.42E+03	1	Methyl Isocyanate	624-83-9							2.0E+01	8.2E+01		1.6E+01
				1.40E+00	I	7.00E-01	I V			2.36E+03	1.36E+09	6.33E+03	1	Methyl Methacrylate	80-62-6							1.1E+05			4.6E+03
				2.50E-04	I						1.36E+09		1	Methyl Parathion	298-00-0							2.0E+01	8.2E+01		1.6E+01
				6.00E-02	X						1.36E+09		1	Methyl Phosphonic Acid	993-13-5							4.7E+03	2.0E+04		3.8E+03
				6.00E-03	H	4.00E-02	H V			3.93E+02	1.36E+09	2.43E+04	1	Methyl Styrene (Mixed Isomers)	25013-15-4						4.7E+02			1.0E+03	3.2E+02
9.90E-02	C	2.80E-05	C								1.36E+09		1	Methyl methanesulfonate	66-27-3	7.0E+00	2.5E+01	1.4E+05	5.5E+00						
1.80E-03	C	2.60E-07	C								1.36E+09		1	Methyl tert-Butyl Ether (MTBE)	1634-04-4	3.9E+02		5.3E+01	4.7E+01						
				3.00E-04	X						1.36E+09		1	Methyl-1,4-benzenediamine dihydrochloride, 2-Methyl-2-Pentanol, 4-Methyl-5-Nitroaniline, 2-Methyl-N-nitro-N-nitrosoguanidine, N-Methylbenzene, 1,4-diamine monohydrochloride, 2-Methylbenzene-1,4-diamine sulfate, 2-Methylcyclohexane	615-45-2					2.3E+01	9.9E+01		1.5E+04	1.5E+04	
				2.00E-02	X					2.45E+03	1.36E+09	1.72E+04	1	Methyl-2-Pentanol, 4-Methyl-5-Nitroaniline, 2-Methyl-N-nitro-N-nitrosoguanidine, N-Methylbenzene, 1,4-diamine monohydrochloride, 2-Methylbenzene-1,4-diamine sulfate, 2-Methylcyclohexane	108-11-2	7.7E+01	2.7E+02		6.0E+01	1.6E+03	6.6E+03	5.4E+04	5.4E+04		
9.00E-03	P			2.00E-02	X						1.36E+09		1	Methyl-2-Pentanol, 4-Methyl-5-Nitroaniline, 2-Methyl-N-nitro-N-nitrosoguanidine, N-Methylbenzene, 1,4-diamine monohydrochloride, 2-Methylbenzene-1,4-diamine sulfate, 2-Methylcyclohexane	99-55-8	8.4E-02	3.0E-01	1.6E+03	6.5E-02						
8.30E+00	C	2.40E-03	C								1.36E+09		1	Methyl-2-Pentanol, 4-Methyl-5-Nitroaniline, 2-Methyl-N-nitro-N-nitrosoguanidine, N-Methylbenzene, 1,4-diamine monohydrochloride, 2-Methylbenzene-1,4-diamine sulfate, 2-Methylcyclohexane	70-25-7										
1.30E-01	C	3.70E-05	C								1.36E+09		1	Methyl-2-Pentanol, 4-Methyl-5-Nitroaniline, 2-Methyl-N-nitro-N-nitrosoguanidine, N-Methylbenzene, 1,4-diamine monohydrochloride, 2-Methylbenzene-1,4-diamine sulfate, 2-Methylcyclohexane	636-21-5	5.3E+00	1.9E+01	1.0E+05	4.2E+00						
				1.00E-02	A																				

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; T = ATSDR DRAFT; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; R = ORD; N = WI; W = TEF applied; E = RPF applied; G = see user's guide; 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	RD ₅₀ (mg/kg-day)	k _e y	RIC ₁ (mg/m ³)	k _e y	v _o l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS ₀	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 THQ=1 (mg/kg)
1.80E+00	C	0.00E+00	C	2.00E-03 3.00E-02	I X	1.00E-01	V P				1.36E+09 5.70E+04	1			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
2.60E-04	C	1.10E-02	C	1.20E-01 1.10E-02	O C	1.40E-05	C				1.36E+09 1.36E+09	1	0.1		Napropamide Nickel Acetate Nickel Carbonate	15299-99-7 373-02-4 3333-67-3		1.5E+04	1.5E+04	1.5E+04	8.6E+02 3.6E+03	4.0E+04 3.6E+03	2.0E+04 2.0E+04	7.6E+03 6.7E+02
2.60E-04	C	1.10E-02	C	1.20E-02	C	1.40E-05	C				1.36E+09	1	0.1		Nickel Carbonyl Nickel Hydroxide Nickel Oxide	13463-39-3 12054-49-7 1313-99-1	1.5E+04	1.5E+04	1.5E+04	8.6E+02 8.6E+02	1.5E+04 1.5E+04	2.0E+04 2.0E+04	8.2E+02 8.2E+02	
1.70E+00	C	4.80E-04	I	2.40E-04 2.60E-04	I C	1.10E-02 2.00E-02	C I	1.00E-05	T		1.36E+09 1.36E+09	0.04			Nickel Refinery Dust Nickel Soluble Salts Nickel Sub sulfide	E715532 7440-02-0 12035-72-2	4.1E-01	2.7E+00	8.0E+03	4.1E-01	8.6E+02 8.6E+02	1.6E+03 2.0E+04	1.4E+04 2.0E+04	8.2E+02 8.2E+02
9.10E-01	C	2.60E-04	C	1.10E-02 1.60E+00	C I	1.40E-05	C				1.36E+09 1.36E+09	1	0.1		Nickelocene Nitrate (measured as nitrogen) Nitrate + Nitrite (measured as nitrogen)	1271-28-9 14797-65-8 E701177	7.6E-01	2.7E+00	1.5E+04	6.0E-01	8.6E+02 1.3E+05	3.6E+03	2.0E+04	6.7E+02 1.3E+05
2.00E-02	P	4.00E-05	I	1.00E-01 4.00E-03	X P	5.00E-05	X	5.00E-05	X		1.36E+09 1.36E+09	1	0.1		Nitrite (measured as nitrogen) Nitroaniline, 2- Nitroaniline, 4-	14797-65-0 88-74-4 100-01-6	3.5E+01	1.2E+02		2.7E+01	7.8E+03 3.1E+02	3.3E+03 1.3E+03	7.1E+04 8.5E+06	7.8E+03 2.5E+02
1.30E+00	C	3.70E-04	C	2.00E-03 7.00E-02	I H	9.00E-03	I	V		3.05E+03	7.32E+04	1			Nitrobenzene Nitrocellulose Nitrofurantoin	98-95-3 9004-70-0 67-20-9			5.1E+00	5.1E+00	1.6E+02 2.3E+08	9.9E+08 2.3E+04	6.9E+02 1.9E+08	1.3E+02 4.4E+03
1.70E-02	P	1.00E-04	P	1.00E-01	I						1.36E+09	1	0.1		Nitrofurazone Nitroglycerin Nitroquinidine	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
2.70E+01	C	7.70E-03	C	8.80E-06 5.80E-04	P X	5.00E-03 2.00E-02	P I	V		1.80E+04 4.86E+03	1.69E+04 1.31E+04	1	0.1		Nitromethane Nitropropane, 2- Nitroso-N-ethylurea, N-	75-52-5 79-46-9 759-73-9			5.4E+00 6.4E-02	5.4E+00 6.4E-02			8.8E+01 2.7E+02	
1.20E-02	C	3.40E-02	C	5.40E+00 2.80E+00	I I	1.60E-03	I				1.36E+09	2.43E+05	1	0.1		Nitroso-N-methylurea, N- Nitrosodibutylamine, N- Nitrosodiethanolamine, N-	684-93-5 924-16-3 1116-54-7	5.7E-03	2.2E-02	1.8E+02	4.5E-03			
5.10E+01	I	1.40E-02	I	4.90E-03	I	2.60E-06	I				1.36E+09	8.23E+04	1	0.1		Nitrosodiethylamine, N- Nitrosodiphenylamine, N- Nitrosodipropylamine, N-	55-18-5 62-75-9 86-30-6	1.0E-03 3.0E-03	4.0E-03	3.2E+01	8.1E-04	6.3E-01	3.4E+00	5.3E-01
7.00E+00	I	2.00E-03	C	2.00E-03	H						1.36E+09	1.21E+05	1	0.1		Nitrosodipropylamine, N- Nitrosomethylmethylethylamine, N- Nitrosomorpholine [N]	621-64-7 10595-95-6 59-89-2	9.9E-02	3.5E-01	1.9E+03	7.8E-02			
9.40E+00	C	2.70E-03	C	2.10E+00	I	6.10E-04	I				1.36E+09	1	0.1		Nitrosopiperidine [N]- Nitrosopyrrolidine, N- Nitrotoluene, m-	100-75-4 930-55-2 99-08-1	7.4E-02	2.6E-01	1.4E+03	5.8E-02	7.8E+00	3.3E+01		6.3E+00
2.20E-01	P	9.00E-04	P	1.60E-02	P	3.00E-04	X	2.00E-02	P	V	1.51E+03	1.37E+05	1	0.1		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	1.3E+03	2.5E+02 1.1E+01
7.79E-03	O	2.00E-03	H	1.90E-01 5.00E-03	O I						1.36E+09	1	0.1		Norfurazone Octabromodiphenyl Ether Octahydro-1,3,5,7-tetrahydro-1,3,5,7-tetrazocine (HMX)	27314-13-2 32536-52-0 2691-41-0					1.2E+02 2.3E+02	4.9E+02 9.9E+02		9.5E+01 1.9E+02
7.79E-03	O	2.00E-03	H	1.90E-01 5.00E-03	O I						1.36E+09	1	0.1		Octamethylphosphoramide Oryzalin Oxadiazon	152-18-9 19044-88-3 19666-30-9	8.9E+01	3.2E+02		7.0E+01	1.6E+02 1.5E+04	6.6E+02 6.3E+04	1.3E+02 1.2E+04	3.2E+02
7.32E-02	O	2.50E-02	I	4.00E-02	O						1.36E+09	1	0.1		Oxamyl Oxyflufen Paclobutrazol	23135-22-0 42874-03-3 76738-62-0	9.5E+00	3.4E+01		7.4E+00	2.0E+03 3.1E+03	8.2E+03 1.3E+04	1.6E+03 2.5E+03	
9.00E-02	P	8.00E-04	I	4.50E-03 6.00E-03 5.00E-02	I H H						1.36E+09 1.36E+09 1.36E+09	4.49E+04	1	0.1		Paraquat Dichloride Parathion Pebulate	1910-42-5 56-38-2 1114-71-2				3.5E+02 4.7E+02 3.9E+03	1.5E+03 2.0E+03	2.8E+02 3.8E+02	3.9E+03
2.60E-01	H	3.00E-03	I	2.00E-03	I						1.36E+09	5.13E+05	1	0.1		Pendimethalin Pentabromodiphenyl Ether Pentabromodiphenyl ether, 2,2',4,4',5-(BDE-99)	40487-42-1 32534-81-9 60348-60-9				2.3E+04 1.6E+02	9.9E+04 3.3E+01	1.9E+04 1.6E+02	6.3E+00
4.00E-01	I	5.10E-06	C	4.30E-03	X	1.00E-04	X				1.36E+09 1.36E+09	1	0.25		Pentachloroethane Pentachloronitrobenzene Pentachlorophenol Pentaerythritol tetranitrate (PETN) Pentamethylphosphoramide (PMPA)	76-01-7 82-68-8 87-86-5 78-11-5 10159-46-3	7.7E+00 2.7E+00	2.5E+00	7.5E+05	7.7E+00 2.7E+00	3.9E+02 2.3E+02	6.6E+02 3.0E+03	2.5E+02 5.7E+02	
2.93E+04	D	3.00E-06	D	3.00E-08 3.00E-04 3.00E-06	D R D	1.00E+00	P	V		3.88E+02	7.79E+02	1	0.1		Pentane, n- Per- and Polyfluoroalkyl Substances (PFAS) ~Ammonium perfluoro-2-methyl-3-oxahexanoate	109-66-0 62037-80-3					2.3E-01 7.8E+01	9.9E-01	8.1E+02	1.9E+01
2.93E+04	D	3.00E-06	D	3.00E-08 3.00E-04 3.00E-06	D R D	1.00E+00	P	V		3.88E+02	7.79E+02	1	0.1		~Ammonium perfluorobutanoate ~Ammonium perfluorodecanoate ~Ammonium perfluorooctanoate	10495-86-0 3108-42-7 21615-47-4					7.8E+01 1.6E-04	9.9E-01	8.1E+02	1.9E+01
2.93E+04	D	3.00E-06	D	3.00E-08 3.00E-04 3.00E-06	D R D	1.00E+00	P	V		3.88E+02	7.79E+02	1	0.1		~Ammonium perfluorooctanoate ~Bis(trifluoromethylsulfonyl)amine (TFSI) ~Hexafluoropropylene oxide dimer acid (HFPO-DA)	3825-26-1 82113-65-3 13252-13-6	2.4E-05	8.4E-05		1.9E-05	2.3E-03 2.3E+01	9.9E-03	8.1E+02	1.9E+01
2.93E+04	D	3.00E-06	D	3.00E-08 3.00E-04 3.00E-06	D R D	1.00E+00	P	V		3.88E+02	7.79E+02	1	0.1		~Lithium bis(trifluoromethyl)sulfonylazanide ~Perfluoro(2-propoxypropanoate) ~Perfluorobutanesulfonate	90076-65-6 122499-17-6 45187-15-3					2.3E+01 2.3E-01	9.9E+01	8.1E+02	1.9E+01
2.93E+04	D	3.00E-06	D	3.00E-08 3.00E-04 3.00E-06	D R D	1.00E+00	P	V		3.88E+02	7.79E+02	1	0.1		~Perfluorobutanesulfonic acid (PFBS) ~Perfluorobutanoate ~Perfluorobutanoic acid (PFBA)	375-73-5 45048-62-2 375-22-4					2.3E+01 7.8E+01	9.9E+01	8.1E+02	1.9E+01
2.93E+04	D	3.00E-06	D	3.00E-08 3.00E-04 3.00E-06	D R D	1.00E+00	P	V		3.88E+02	7.79E+02	1	0.1		~Perfluorodecanoate ~Perfluorodecanoic acid (PFDA) ~Perfluorododecanoic acid (PFDoDA)	73829-36-4 335-76-2 307-55-1					1.6E-04 1.6E-04	6.6E-04	8.1E+02	1.9E+01
2.93E+04	D	3.00E-06	D	3.00E-08 3.00E-04 3.00E-06	D R D	1.00E+00	P	V		3.88E+02	7.79E+02	1	0.1		~Perfluorohexanesulfonate ~Perfluorohexanesulfonic acid (PFHxS)	108427-53-8 355-46-4					1.6E+00 1.6E+00	6.6E+00	8.1E+02	1.9E+01

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; T = ATSDR DRAFT; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; R = ORD; N = WI; W = TEF applied; E = RPF applied; G = see user's guide; 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	RD ₅₀ (mg/kg-day)	k _e y	RIC ₁ (mg/m ³)	k _e y	v _o l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS ₀	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)	
				5.00E-04	I						1.36E+09			1	0.1	-Perfluorohexanoate	92612-52-7					3.9E+01	1.6E+02		3.2E+01
				5.00E-04	I						1.36E+09			1	0.1	-Perfluorohexanoic acid (PFHxA)	307-24-4					3.9E+01	1.6E+02		3.2E+01
				3.00E-06	A						1.36E+09			1	0.1	-Perfluorononanoate	72007-68-2					2.3E-01	9.9E-01		1.9E-01
				3.00E-06	A						1.36E+09			1	0.1	-Perfluorononanoic acid (PFNA)	375-95-1					2.3E-01	9.9E-01		1.9E-01
				4.00E-02	N						1.36E+09			1	0.1	-Perfluorooctadecanoic acid (PFODA)	16517-11-6					3.1E+03	3.3E+04		2.5E+03
3.95E+01	D			1.00E-07	D						1.36E+09			1	0.1	-Perfluorooctanesulfonate	45298-90-6	1.8E-02	6.3E-02		1.4E-02	7.8E-03	1.3E+02		6.3E-03
3.95E+01	D			1.00E-07	D						1.36E+09			1	0.1	-Perfluorooctanesulfonic acid (PFOS)	1763-23-1	1.8E-02	6.3E-02		1.4E-02	7.8E-03	3.3E-02		6.3E-03
2.93E+04	D			3.00E-08	D						1.36E+09			1	0.1	-Perfluorooctanoate	45285-51-6	2.4E-05	8.4E-05		1.9E-05	2.3E-03	9.9E-03		1.9E-03
2.93E+04	D			3.00E-08	D						1.36E+09			1	0.1	-Perfluorooctanoic acid (PFOA)	335-67-1	2.4E-05	8.4E-05		1.9E-05	2.3E-03	9.9E-03		1.9E-03
				5.00E-04	R			V		1.36E+04	1.36E+09	2.65E+04		1		-Perfluoropropanoic acid (PFPrA)	422-64-0					3.9E+01			3.9E+01
				1.00E-03	N						1.36E+09			1	0.1	-Perfluorotetradecanoic acid (PFTeDA)	376-06-7					7.8E+01	3.3E+02		6.3E+01
				3.00E-04	N						1.36E+09			1	0.1	-Perfluoroundecanoic acid (PFUDA)	2058-94-8					2.3E+01	9.9E+01		1.9E+01
				3.00E-04	P						1.36E+09			1	0.1	-Potassium perfluorobutanesulfonate	29420-49-3					2.3E+01	9.9E+01		1.9E+01
				2.00E-03	I			V		9.61E+04	1.36E+09	6.13E+04		1		-Potassium perfluorobutanoate	2966-54-3					1.6E+02			1.6E+02
				2.15E-09	I						1.36E+09			1	0.1	-Potassium perfluorodecanoate	51604-85-4					1.7E-04	7.1E-04		1.4E-04
3.95E+01	D			1.00E-07	D						1.36E+09			1	0.1	-Potassium perfluorooctanesulfonate	2795-39-3	1.8E-02	6.3E-02		1.4E-02	7.8E-03	3.3E-02		6.3E-03
				1.00E-03	I			V		8.99E+04	1.36E+09	6.02E+04		1		-Sodium perfluorobutanoate	2218-54-4					7.8E+01			7.8E+01
				2.09E-09	I						1.36E+09			1	0.1	-Sodium perfluorodecanoate	3830-45-3					1.6E-04	6.9E-04		1.3E-04
				5.00E-04	I						1.36E+09			1	0.1	-Sodium perfluorohexanoate	2923-26-4					3.9E+01	1.6E+02		3.2E+01
				7.00E-04	I						1.36E+09			1		-Perchlorates						5.5E+01			5.5E+01
				7.00E-04	I						1.36E+09			1		-Ammonium Perchlorate	7790-98-9					5.5E+01			5.5E+01
				7.00E-04	I						1.36E+09			1		-Lithium Perchlorate	7791-03-9					5.5E+01			5.5E+01
				7.00E-04	I						1.36E+09			1		-Perchlorate and Perchlorate Salts	14797-73-0					5.5E+01			5.5E+01
				7.00E-04	I						1.36E+09			1		-Potassium Perchlorate	7778-74-7					5.5E+01			5.5E+01
				7.00E-04	I						1.36E+09			1		-Sodium Perchlorate	7601-89-0					5.5E+01			5.5E+01
2.20E-03	C	6.30E-07	C	5.00E-02	I						1.36E+09			1	0.1	Permethrin	52645-53-1	3.2E+02	1.1E+03	6.1E+06	2.5E+02	3.9E+03	1.6E+04		3.2E+03
				2.40E-01	O						1.36E+09			1	0.1	Phenacetin	62-44-2					1.9E+04	7.9E+04		1.5E+04
				3.00E-01	I	2.00E-01	C				1.36E+09			1	0.1	Phenmedipham	13684-63-4					2.3E+04	9.9E+04	2.8E+08	1.9E+04
				4.00E-03	I						1.36E+09			1	0.1	Phenol	108-95-2					3.1E+02	1.3E+03		2.5E+02
				5.00E-04	X						1.36E+09			1	0.1	Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1					3.9E+01	1.6E+02		3.2E+01
				2.00E-04	X			V		1.29E+02	1.36E+09	7.06E+03		1		Phenothiazine	92-84-2					3.9E+01	1.6E+02		3.2E+01
1.20E-01	P			6.00E-03	I						1.36E+09			1	0.1	Phenyl isothiocyanate	103-72-0					1.6E+01			1.6E+01
				4.00E-03	P			M			1.36E+09			1	0.1	Phenylenediamine, m-	108-45-2					4.7E+02	2.0E+03		3.8E+02
				4.00E-03	P						1.36E+09			1	0.1	Phenylenediamine, o-	95-54-5	1.3E+00	5.0E+00		1.0E+00	3.1E+02	1.3E+03		2.5E+02
1.94E-03	H			1.00E-03	X						1.36E+09			1	0.1	Phenylenediamine, p-	106-50-3					7.8E+01	3.3E+02		6.3E+01
				2.00E-04	H						1.36E+09			1	0.1	Phenylphenol, 2-	90-43-7	3.6E+02	1.3E+03		2.8E+02	1.6E+01	6.6E+01		1.3E+01
				2.00E-02	I	3.00E-04	I	V		1.61E+03	1.36E+09	9.81E+02		1	0.1	Phorate	298-02-2					1.6E+01	6.6E+03	3.1E-01	1.3E+03
				2.93E+00	X						1.36E+09			1		Phosgene	75-44-5					1.6E+03	6.6E+03		3.1E-01
				3.00E-01	X						1.36E+09			1		Phosmet	732-11-6					1.6E+03	6.6E+03		1.3E+03
				1.00E+00	P						1.36E+09			1		Phosphates, Inorganic									
				1.00E+00	P						1.36E+09			1		-Aluminum metaphosphate	13776-88-0					2.3E+05			2.3E+05
				1.00E+00	P						1.36E+09			1		-Aluminum salts of inorganic phosphates	E524680405					2.3E+04			2.3E+04
				1.00E+00	P						1.36E+09			1		-Dipotassium phosphate	7758-11-4					7.8E+04			7.8E+04
				1.00E+00	P						1.36E+09			1		-Disodium phosphate	7558-79-4					7.8E+04			7.8E+04
				3.54E+00	X						1.36E+09			1	0.1	-Monosodium phosphate	13530-50-2					2.8E+05			2.8E+05
				1.00E+00	P						1.36E+09			1		-Monopotassium phosphate	7778-77-0					7.8E+04			7.8E+04
				1.00E+00	P						1.36E+09			1		-Monosodium phosphate	7558-80-7					7.8E+04			7.8E+04
				1.00E+00	P	1.00E-02	I				1.36E+09			1		-Phosphoric acid	7664-38-2					7.8E+04		1.4E+07	7.8E+04
				1.36E+00	X						1.36E+09			1	0.1	-Phosphoric acid, aluminum salt (1:1) [aluminum phosphate]	7784-30-7					1.1E+05	4.5E+05		8.6E+04
				4.26E+00	X						1.36E+09			1		-Phosphoric acid, aluminum sodium salt (1:X:X) [sodium aluminum phosphate acidic (acidic SALP)]	7785-88-8					3.3E+05			3.3E+05
				1.00E+00	P						1.36E+09			1		-Polyphosphoric acid	8017-16-1					7.8E+04			7.8E+04
				1.00E+00	P						1.36E+09			1		-Potassium salts of inorganic phosphates	E524680403					7.8E+04			7.8E+04
				1.00E+00	P						1.36E+09			1		-Potassium tripolyphosphate	13845-36-8					7.8E+04			7.8E+04
				4.99E+00	X						1.36E+09			1		-Sodium aluminum phosphate (anhydrous)	10279-59-1					3.9E+05			3.9E+05
				3.52E+00	X						1.36E+09			1		-Sodium aluminum phosphate (tetrahydrate)	10305-76-7					2.8E+05			2.8E+05
				1.00E+00	P						1.36E+09			1		-Sodium hexametaphosphate	10124-56-8					7.8E+04			7.8E+04
				1.00E+00	P																				

Key: I = IRIS; P = PPRVT; O = OPP; A = ATSDR; T = ATSDR DRAFT; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; R = ORD; N = WI; W = TEF applied; E = RPF applied; G = see user's guide; 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	RD ₅₀ (mg/kg-day)	k e y	RIC ₁ (mg/m ³)	k e y	v o	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS ₀	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 THQ=1 (mg/kg)
3.00E+01	C	8.60E-03	C	7.30E-04	O	7.00E-06	H				1.36E+09		1	0.1	Pirimiphos, Methyl	29232-93-7	2.3E-02	8.2E-02	4.4E+02	1.8E-02	5.7E+01	2.4E+02		4.6E+01
											1.36E+09		1	0.1	Polybrominated Biphenyls	36355-01-8	5.5E-01				5.5E-01	2.3E+00		4.4E-01
											1.36E+09	5.86E+05	1	0.14	Polychlorinated Biphenyls (PCBs)									
7.00E-02	G	2.00E-05	G	7.00E-05	I			V			1.36E+09	2.04E+05	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
2.00E+00	G	5.71E-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				
2.00E+00	G	5.71E-04	G					V			1.36E+09	1.12E+05	1	0.14	~Aroclor 1232	11141-16-5	3.5E-01	8.8E-01	1.0E+00	2.0E-01				
2.00E+00	G	5.71E-04	G					V			1.36E+09	5.91E+05	1	0.14	~Aroclor 1242	53409-21-9	3.5E-01	8.8E-01	2.9E+00	2.3E-01				
2.00E+00	G	5.71E-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				
2.00E+00	G	5.71E-04	G	2.00E-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
2.00E+00	G	5.71E-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				
				6.00E-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
3.90E+00	W	1.14E-03	W	2.33E-05	W	1.33E-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
3.90E+00	W	1.14E-03	W	2.33E-05	W	1.33E-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
3.90E+00	W	1.14E-03	W	2.33E-05	W	1.33E-03	W	V			1.36E+09	1.04E+06	1	0.14	~Hexachlorobiphenyl, 2,3,3',4,4',5,5'-(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
3.90E+00	W	1.14E-03	W	2.33E-05	W	1.33E-03	W	V			1.36E+09	1.11E+06	1	0.14	~Hexachlorobiphenyl, 2,3,3',4,4',5,5'-(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
3.90E+03	W	1.14E+00	W	2.33E-08	W	1.33E-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
3.90E+00	W	1.14E-03	W	2.33E-05	W	1.33E-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
3.90E+00	W	1.14E-03	W	2.33E-05	W	1.33E-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
3.90E+00	W	1.14E-03	W	2.33E-05	W	1.33E-03	W	V			1.36E+09	6.01E+05	1	0.14	~Tetrachlorobiphenyl, 2,3,3',4,4'-(PCB 105)	32599-14-4	1.8E-01	4.5E-01	1.5E+00	1.2E-01	1.8E+00	5.5E+00	8.4E+02	1.4E+00
3.90E+00	W	1.14E-03	W	2.33E-05	W	1.33E-03	W	V			1.36E+09	1.05E+06	1	0.14	~Pentachlorobiphenyl, 2,3,4,4',5-(PCB 114)	74472-31-0	1.8E-01	4.5E-01	2.6E+00	1.2E-01	1.8E+00	5.5E+00	1.5E+03	1.4E+00
1.30E+04	W	3.80E+00	W	7.00E-09	W	4.00E-07	W	V			1.36E+09	7.26E+05	1	0.14	~Pentachlorobiphenyl, 3,3',4,4',5-(PCB 128)	57465-29-8	5.3E-05	1.4E-04	5.4E-04	3.6E-05	5.5E-04	1.6E-03	3.0E-01	4.1E-04
2.00E+00	I	5.71E-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				
4.00E-01	I	1.00E-04	I					V			1.36E+09		1	0.14	~Polychlorinated Biphenyls (low risk)	1336-36-3								
7.00E-02	I	2.00E-05	I					V			1.36E+09		1	0.14	~Polychlorinated Biphenyls (lowest risk)	1336-36-3								
1.30E+01	W	3.80E-03	W	7.00E-06	W	4.00E-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
3.90E+01	W	1.14E-02	W	2.33E-06	W	1.33E-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
				6.00E-04	I			V			1.36E+09		1	0.1	Polymeric Methylenediphenyl Diisocyanate (PMDI)	9016-87-9							8.5E+05	8.5E+05
				6.00E-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
				3.00E-01	I			V			1.36E+09	5.23E+05	1	0.13	~Anthracene	120-12-7					2.3E+04	7.6E+04		1.8E+04
1.00E-01	E	6.00E-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				
1.00E+00	I	6.00E-04	I	3.00E-04	I	2.00E-06	I	M			1.36E+09		1	0.13	~Benz[a]pyrene	50-32-8	1.5E-01	4.6E-01	2.3E+03	1.1E-01	2.3E+01	7.6E+01	2.8E+03	1.8E+01
1.00E-01	E	6.00E-05	E					V	M		1.36E+09		1	0.13	~Benz[b]fluoranthene	205-99-2	1.5E+00	4.6E+00	2.3E+04	1.1E+00				
				9.00E-05	X	2.00E-06	X				1.36E+09		1	0.1	~Benz[e]pyrene	192-97-2					7.0E+00	3.0E+01	2.8E+03	5.7E+00
1.20E+00	C	1.10E-04	C					V			1.36E+09		1	0.13	~Benzofluoranthene	205-82-3	5.8E-01	1.6E+00	3.5E+04	4.2E-01				
1.00E-02	E	6.00E-06	E					V	M		1.36E+09		1	0.13	~Benzofluoranthene	207-08-9	1.5E+01	4.6E+01	2.3E+05	1.1E+01				
				5.00E-03	X			V			1.36E+09		1	0.1	~Benzofluorene, 2,3-	243-17-4					3.9E+02	1.6E+03		3.2E+02
				8.00E-02	I			V			1.36E+09	7.99E+04	1	0.13	~Chloronaphthalene, Beta-	91-58-7					6.3E+03	2.0E+04		4.8E+03
1.00E-03	E	6.00E-07	E					V	M		1.36E+09		1	0.13	~Chrysene	218-01-9	1.5E+02	4.6E+02	2.3E+06	1.1E+02				
1.00E+00	E	6.00E-04	E					V	M		1.36E+09		1	0.13	~Dibenz[a,h]anthracene	53-70-3	1.5E-01	4.6E-01	2.3E+03	1.1E-01				
1.20E+01	C	1.10E-03	C					V			1.36E+09		1	0.13	~Dibenz[a,e]pyrene	192-65-4	5.8E-02	1.6E-01	3.5E+03	4.2E-02				
2.50E+02	C	7.10E-02	C					V	M		1.36E+09		1	0.13	~Dimethylbenz[a]anthracene, 7,12-	57-97-6	6.1E-04	1.8E-03	1.9E+01	4.6E-04				
				4.00E-02	I			V			1.36E+09		1	0.13	~Fluoranthene	206-44-0					3.1E+03	1.0E+04		2.4E+03
				4.00E-02	I			V			1.36E+09	2.81E+05	1	0.13	~Fluorene	86-73-7					3.1E+03	1.0E+04		2.4E+03
1.00E-01	E	6.00E-05	E					V	M		1.36E+09		1	0.13	~Indeno[1,2,3-cd]pyrene	193-39-5	1.5E+00	4.6E+00	2.3E+04	1.1E+00				
5.10E-02	X	7.00E-02	T	3.00E-06	P	V		V		3.94E+02	1.36E+09	5.86E+04	1	0.13	~Methylnaphthalene, 1-	90-12-0	1.4E+01	3.7E+01		1.0E+01	5.5E+03	1.8E+04	1.8E-01	1.8E-01
				4.00E-03	I			V			1.36E+09	5.80E+04	1	0.13	~Methylnaphthalene, 2-	91-57-6					3.1E+02	1.0E+03		2.4E+02
1.20E-01	C	3.40E-05	C	2.00E-02	I	3.00E-03	I	V			1.36E+09	4.63E+04	1	0.13	~Naphthalene	91-20-3	5.8E+00	1.6E+01	3.8E+00	2.0E+00	1.6E+03	5.1E+03	1.4E+02	1.3E+02
1.20E+00	C	1.10E-04	C					V			1.36E+0													

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; T = ATSDR DRAFT; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; R = ORD; N = WI; W = TEF applied; E = RPF applied; G = see user's guide; 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 = Caat 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	key	IUR (ug/m ³) ¹	key	RD ₅₀ (mg/kg-day)	key	RIC ₁ (mg/m ³)	key	vo	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS ₀	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)	
				5.00E-03	I	3.00E-03	C				1.36E+09		1		Silica (crystalline, respirable)	7631-86-9							4.3E+06	4.3E+06	
1.20E-01	H			5.00E-03	I						1.36E+09		0.04		Silver	7440-22-4	5.8E+00	2.1E+01		4.5E+00	3.9E+02	1.6E+03	3.9E+02	3.9E+02	
				1.30E-02	I						1.36E+09		1	0.1	Simazine	122-34-9					3.9E+02	4.3E+03	4.3E+03	3.2E+02	
2.70E-01	H			4.00E-03	I						1.36E+09		1		Sodium Acifluorfen	62476-59-9	2.6E+00	9.2E+00		2.0E+00	3.1E+02	9.9E+03	9.9E+03	3.1E+02	
				3.00E-02	I						1.36E+09		1	0.1	Sodium Azide	26628-22-8					3.1E+02	9.9E+03	9.9E+03	1.9E+03	
				5.00E-02	A	1.40E-02	C				1.36E+09		1		Sodium Diethyldithiocarbamate	148-18-5					3.9E+03	6.6E+00	2.0E+07	3.9E+03	
				2.00E-05	I						1.36E+09		1	0.1	Sodium Fluoroacetate	7681-49-4					1.6E+00	6.6E+00	2.0E+07	1.3E+00	
				1.00E-03	H						1.36E+09		1		Sodium Melavanadate	13718-26-8					7.8E+01			7.8E+01	
2.40E-02	H			8.00E-04	P						1.36E+09		1		Sodium Tungstate	13472-45-2					6.3E+01			6.3E+01	
				8.00E-04	P						1.36E+09		1		Sodium Tungstate Dihydrate	10213-10-2					6.3E+01			6.3E+01	
				3.00E-02	I						1.36E+09		1	0.1	Stirofos (Tetrachlorovinphos)	961-11-5	2.9E+01	1.0E+02		2.3E+01	2.3E+03	9.9E+03	9.9E+03	1.9E+03	
				6.00E-01	I						1.36E+09		1		Strontium, Stable	7440-24-6					4.7E+04			4.7E+04	
				3.00E-04	I						1.36E+09		1	0.1	Strychnine	57-24-9					2.3E+01	9.9E+01	9.9E+01	1.9E+01	
				2.00E-01	I	1.00E+00	V			8.67E+02	1.36E+09	9.35E+03	1		Styrene	100-42-5					1.6E+04		9.7E+03	6.0E+03	
				3.00E-03	P						1.36E+09		1	0.1	Styrene-Acrylonitrile (SAN) Trimer (THNA isomer)	57964-39-3					2.3E+02	9.9E+02	9.9E+02	1.9E+02	
				3.00E-03	P						1.36E+09		1	0.1	Styrene-Acrylonitrile (SAN) Trimer (THNP isomer)	57964-40-6					2.3E+02	9.9E+02	9.9E+02	1.9E+02	
				1.00E-03	P	2.00E-03	X				1.36E+09		1	0.1	Sulffolane	126-33-0					7.8E+01	3.3E+02	2.8E+06	6.3E+01	
				8.00E-04	P						1.36E+09		1	0.1	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9					6.3E+01	2.6E+02	2.6E+02	5.1E+01	
						1.00E-03	C	V			1.36E+09		1		Sulfur Trioxide	7446-11-9					6.3E+01			1.4E+06	
						1.00E-03	C				1.36E+09		1		Sulfuric Acid	7664-93-9					6.3E+01			1.4E+06	
2.50E-02	I	7.10E-06	I	5.00E-02	H						1.36E+09		1	0.1	Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylphenyl)phenoxy]-1-methylethyl ester	140-57-8	2.8E+01	9.9E+01	5.4E+05	2.2E+01	3.9E+03	1.6E+04	1.6E+04	3.2E+03	
				7.00E-02	I						1.36E+09		1	0.1	Tebuthiuron	34014-18-1					5.5E+03	2.3E+04	2.3E+04	4.4E+03	
				2.00E-02	H						1.36E+09		1	0.1	Temephos	3383-96-8					1.6E+03	6.6E+03	6.6E+03	1.3E+03	
				1.30E-02	I						1.36E+09		1	0.1	Terbacil	5902-51-2					1.0E+03	4.3E+03	4.3E+03	8.2E+02	
				2.50E-05	H				V	3.09E+01	1.36E+09	2.64E+05	1		Terbufos	13071-79-9					2.0E+00			2.0E+00	
				1.00E-03	I						1.36E+09		1	0.1	Terbutyn	886-50-0					7.8E+01	3.3E+02	3.3E+02	6.3E+01	
5.00E-03	C	1.30E-06	C	1.00E-04	I						1.36E+09	3.99E+03	1	0.1	Tert-Butyl Acetate	540-88-5	1.4E+02		8.6E+00	8.1E+00	7.8E+01	3.3E+01	3.3E+01	6.3E+00	
				3.00E-05	P						1.36E+09		1		Tetrabromodiphenyl ether, 2,2',4,4'-(BDE-47)	5436-43-1					2.3E+00			2.3E+00	
											1.36E+09	5.07E+04	1		Tetrachlorobenzene, 1,2,4,5-	95-94-3					7.8E+00	3.3E+01	3.3E+01	6.3E+00	
2.60E-02	I	7.40E-06	I	3.00E-02	I					6.80E+02	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	
2.00E-01	I	5.80E-05	C	2.00E-02	I					1.90E+03	1.36E+09	1.51E+04	1		Tetrachloroethane, 1,1,1,2,2-	79-34-5	3.5E+00		7.3E-01	6.0E-01	1.6E+03			1.6E+03	
2.10E-03	I	2.60E-07	I	6.00E-03	I	4.00E-02	V			1.66E+02	1.36E+09	2.35E+03	1		Tetrachloroethylene	127-18-4	3.5E+02		2.5E+01	2.4E+01	4.7E+02		9.8E+01	9.8E+01	
				3.00E-02	I						1.36E+09		1	0.1	Tetrachlorophenol, 2,3,4,6-	58-90-2					2.3E+03	9.9E+03	9.9E+03	1.9E+03	
1.60E+01	X			6.00E-05	X						1.36E+09	1.05E+05	1		Tetrachlorotoluene, p-alpha, alpha, alpha-	5216-25-1	4.3E-02				4.7E+00			4.7E+00	
				5.00E-04	I						1.36E+09		1	0.1	Tetraethyl Dithiopyrophosphate	3689-24-5					3.9E+01	1.6E+02	1.6E+02	3.2E+01	
						8.00E+01	V			2.05E+03	1.36E+09	1.22E+03	1		Tetrafluoroethane, 1,1,1,2-	811-97-2					7.8E+00	3.3E+01	3.3E+01	1.0E+05	
				1.00E-04	X						1.36E+09		1	0.1	Tetramethylphosphoramide, -N,N,N',N'-(TMPA)	16853-36-4					1.6E+00			6.3E+00	
				2.00E-03	P						1.36E+09		1	0.00065	Tetryl (Trinitrophenylmethylnitramine)	479-45-8					1.6E+02	1.0E+05	1.0E+05	1.6E+02	
				2.00E-05	G						1.36E+09		1		Thallic Oxide	1314-32-5					1.6E+00			1.6E+00	
				1.00E-05	X						1.36E+09		1		Thallium (I) Nitrate	10102-45-1					7.8E-01			7.8E-01	
				1.00E-05	X						1.36E+09		1		Thallium (Soluble Salts)	7440-28-0					7.8E-01			7.8E-01	
				1.00E-05	X						1.36E+09	1.40E+05	1		Thallium Acetate	563-68-8					7.8E-01			7.8E-01	
				2.00E-05	X						1.36E+09		1	0.1	Thallium Carbonate	6533-73-9					1.6E+00	6.6E+00	6.6E+00	1.3E+00	
				1.00E-05	X						1.36E+09		1		Thallium Chloride	7781-12-0					7.8E-01			7.8E-01	
				1.00E-05	G						1.36E+09		1		Thallium Selenite	12039-52-0					7.8E-01			7.8E-01	
				2.00E-05	X						1.36E+09		1		Thallium Sulfate	7446-18-6					1.6E+00			1.6E+00	
				4.30E-02	O						1.36E+09		1	0.1	Thifensulfuron-methyl	79277-27-3					3.4E+03	1.4E+04	1.4E+04	2.7E+03	
				1.00E-02	I						1.36E+09		1	0.1	Thiobencarb	28249-77-6					7.8E+02	3.3E+03	3.3E+03	6.3E+02	
				2.00E-04	P						1.36E+09		1		Thiocyanates	E1790665					1.6E+01			1.6E+01	
				2.00E-04	X						1.36E+09		1		Thiocyanic Acid	463-56-9					1.6E+01			1.6E+01	
				3.00E-02	H						1.36E+09		1	0.1	Thiocyanic acid, (2-benzothiazolythio)methyl ester (TCMTB)	21564-17-0					2.3E+03	9.9E+03	9.9E+03	1.9E+03	
				7.00E-02	X						1.36E+09		1	0.0075	Thiodiethylol	111-48-8					5.5E+03	3.1E+05	3.1E+05	5.4E+03	
				3.00E-04	H						1.36E+09		1	0.1	Thiofanox	39196-18-4					2.3E+01	9.9E+01	9.9E+01	1.9E+01	
1.16E-02	O			1.60E-01	O						1.36E+09		1	0.1	Thiophanate, Methyl	23564-05-8	6.0E+01	2.1E+02		4.7E+01	1.3E+04	5.3E+04	5.3E+04	1.0E+04	
				1.50E-02	O						1.36E+09		1	0.1	Thiram	137-26-8					1.2E+03	4.9E+03	4.9E+03	9.5E+02	
				6.00E-01	H						1.36E+09														

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; T = ATSDR DRAFT; C = Cal EPA; X = PPRTV Screening Level; H = HEAST; D = OW; R = ORD; N = WI; W = TEF applied; E = RPF applied; G = see user's guide; 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	RD ₅₀ (mg/kg-day)	k e y	RIC ₁ (mg/m ³)	k e y	v o l	mutagen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS ₀	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)		
9.00E-03	P			2.00E-04	O						1.36E+09		1	0.1	Tribufos	78-48-8					1.6E+01	6.6E+01			1.3E+01	
				1.00E-02	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	
				3.00E-04	P						1.36E+09		1	0.1	Tributyltin Compounds	E1790679					2.3E+01	9.9E+01			1.9E+01	
				3.00E-04	I						1.36E+09		1	0.1	Tributyltin Oxide	56-35-9					2.3E+01	9.9E+01			1.9E+01	
				3.00E+01	I	5.00E+00	P	V		9.10E+02	1.36E+09	1.29E+03	1	0.1	Trichloramine	10025-85-1					2.3E+06	6.6E+03	6.7E+03		6.7E+03	
7.00E-02	I			2.00E-02	I						1.36E+09		1	0.1	Trichloro-1,1,2-trifluoroethane, 1,1,2-	76-13-1					1.6E+03	6.6E+03			1.3E+03	
2.90E-02	H										1.36E+09		1	0.1	Trichloroacetic Acid	76-03-9	9.9E+00	3.5E+01		7.8E+00	7.8E+02	3.3E+03			6.3E+02	
7.00E-03	X			3.00E-05	X						1.36E+09		1	0.1	Trichloroaniline HCl, 2,4,6-	33663-50-2	2.4E+01	8.5E+01		1.9E+01	2.3E+01	9.9E+01			1.9E+00	
				8.00E-04	X						1.36E+09	3.22E+04	1	0.1	Trichloroaniline, 2,4,6-	634-93-5	9.9E+01	3.5E+02		7.8E+01	2.3E+00	9.9E+00			1.9E+00	
											1.36E+09		1	0.1	Trichlorobenzene, 1,2,3-	87-61-6					6.3E+01	9.9E+00			6.3E+01	
2.90E-02	P			1.00E-02	I	2.00E-03	P	V		4.04E+02	1.36E+09	2.99E+04	1	0.1	Trichlorobenzene, 1,2,4-	120-82-1	2.4E+01			2.4E+01	7.8E+02	3.3E+03	6.2E+01		5.8E+01	
				2.00E+00	I	5.00E+00	I	V		6.40E+02	1.36E+09	1.65E+03	1	0.1	Trichloroethane, 1,1,1-	71-55-6					1.6E+05	8.6E+03			8.1E+03	
5.70E-02	I	1.60E-05	I	4.00E-03	I	2.00E-04	X	V		2.16E+03	1.36E+09	7.22E+03	1	0.1	Trichloroethane, 1,1,2-	79-00-5	1.2E+01		1.3E+00	1.1E+00	3.1E+02	1.5E+00			1.5E+00	
4.60E-02	I	4.10E-06	I	5.00E-04	I	2.00E-03	I	V	M	6.92E+02	1.36E+09	2.21E+03	1	0.1	Trichloroethylene	79-01-6	8.8E+00		1.1E+00	9.4E-01	3.9E+01		4.6E+00	4.1E+00		
				3.00E-01	I					1.23E+03	1.36E+09	1.04E+03	1	0.1	Trichlorofluoromethane	75-69-4					2.3E+04			2.3E+04		
				1.00E-01	I						1.36E+09		1	0.1	Trichlorophenol, 2,4,5-	95-95-4					7.8E+03	3.3E+04			6.3E+03	
1.10E-02	I	3.10E-06	I	1.00E-03	P						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	
				1.00E-02	I						1.36E+09		1	0.1	Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5					7.8E+02	3.3E+03			6.3E+02	
				8.00E-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	
3.00E+01	I			5.00E-03	I		V			1.28E+03	1.36E+09	1.50E+04	1	0.1	Trichloropropane, 1,1,2-	598-77-6					3.9E+02				3.9E+02	
				4.00E-03	I	3.00E-04	I	V	M	1.40E+03	1.36E+09	1.57E+04	1	0.1	Trichloropropane, 1,2,3-	96-18-4	5.1E-03			5.1E-03	3.1E+02		4.9E+00		4.8E+00	
				3.00E-03	X	3.00E-04	P	V		3.11E+02	1.36E+09	2.34E+03	1	0.1	Trichloropropane, 1,2,3-	96-19-5					2.3E+02		7.3E-01		7.3E-01	
				2.00E-02	A						1.36E+09		1	0.1	Tricresyl Phosphate (TCP)	1330-78-5					1.6E+03	6.6E+03			1.3E+03	
				3.00E-03	I						1.36E+09		1	0.1	Tridiphane	58138-08-2					2.3E+02	9.9E+02			1.9E+02	
						7.00E-03	I	V		2.79E+04	1.36E+09	1.58E+04	1	0.1	Triethylamine	121-44-8							1.2E+02		1.2E+02	
				2.00E+00	P						1.36E+09		1	0.1	Triethylene Glycol	112-27-6					1.6E+05	6.6E+05			1.3E+05	
7.70E-03	I			7.50E-03	I	2.00E+01	P	V		4.81E+03	1.36E+09	7.12E+02	1	0.1	Trifluoroethane, 1,1,1-	420-46-2				9.0E+01		9.0E+01	1.5E+04		1.5E+04	
2.00E-02	P			1.00E-02	P						1.36E+09		1	0.1	Trifuralin	1582-09-8	9.0E+01				5.9E+02				5.9E+02	
				1.00E-02	I	6.00E-02	I	V		2.93E+02	1.36E+09	9.44E+03	1	0.1	Trimethyl Phosphate	512-56-1	3.5E+01	1.2E+02		2.7E+01	7.8E+02	3.3E+03			6.3E+02	
				1.00E-02	I	6.00E-02	I	V		2.19E+02	1.36E+09	7.91E+03	1	0.1	Trimethylbenzene, 1,2,3-	526-73-8					7.8E+02		5.9E+02		3.4E+02	
				1.00E-02	I	6.00E-02	I	V		2.19E+02	1.36E+09	7.91E+03	1	0.1	Trimethylbenzene, 1,2,4-	95-63-6					7.8E+02		5.0E+02		3.0E+02	
				1.00E-02	I	6.00E-02	I	V		1.82E+02	1.36E+09	6.61E+03	1	0.1	Trimethylbenzene, 1,3,5-	108-67-8					7.8E+02		4.1E+02		2.7E+02	
				1.00E-02	X					2.96E+01	1.36E+09	1.00E+03	1	0.1	Trimethylpentene, 2,4,4-	25167-70-8					7.8E+02				7.8E+02	
				3.00E-02	I						1.36E+09		1	0.019	Trinitrobenzene, 1,3,5-	99-35-4					2.3E+03	5.2E+04			2.2E+03	
3.00E-02	I			5.00E-04	I						1.36E+09		1	0.032	Trinitrotoluene, 2,4,6-	118-96-7	2.3E+01	2.6E+02		2.1E+01	3.9E+01	5.2E+02			3.6E+01	
				2.00E-02	P						1.36E+09		1	0.1	Triphenylphosphine Oxide	791-28-6					1.6E+03	6.6E+03			1.3E+03	
				2.00E-02	A						1.36E+09		1	0.1	Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8					1.6E+03	6.6E+03			1.3E+03	
				1.00E-02	X						1.36E+09		1	0.1	Tris(1-chloro-2-propyl)phosphate	13674-84-5					7.8E+02	3.3E+03			6.3E+02	
2.30E+00	C	6.60E-04	C							4.67E+02	1.36E+09	9.03E+05	1	0.1	Tris(2,3-dibromopropyl)phosphate	126-72-7	3.0E-01		3.8E+00	2.8E-01	5.5E+02	2.3E+03			4.4E+02	
2.00E-02	P			7.00E-03	P						1.36E+09		1	0.1	Tris(2-chloroethyl)phosphate	115-96-8	3.5E+01	1.2E+02		2.7E+01	5.5E+02	2.3E+03			6.3E+03	
3.20E-03	P			1.00E-01	P						1.36E+09		1	0.1	Tris(2-ethylhexyl)phosphate	78-42-2	2.2E+02	7.7E+02		1.7E+02	7.8E+03	3.3E+04			6.3E+03	
				8.00E-04	P						1.36E+09		1	0.1	Tunasten	7440-33-7					6.3E+01				6.3E+01	
				2.00E-04	A	4.00E-05	A				1.36E+09		1	0.1	Uranium	7440-61-1					1.6E+01		5.7E+04			1.6E+01
1.00E+00	C	2.90E-04	C								1.36E+09		1	0.1	Urethane	51-79-6	1.5E-01	6.0E-01	4.8E+03	1.2E-01	7.0E+02				6.6E+02	
				8.30E-03	P	9.00E-03	I	7.00E-06	P		1.36E+09		0.026	0.026	Vanadium Pentoxide	1314-62-1					3.9E+02				3.9E+02	
				5.04E-03	G	1.00E-04	A				1.36E+09		1	0.1	Vanadium and Compounds	7440-62-2					3.9E+02				3.9E+02	
				1.00E-03	I						1.36E+09	1.23E+05	1	0.1	Vernolate	1929-77-7					7.8E+01				7.8E+01	
				1.20E-03	O						1.36E+09		1	0.1	Vinclozolin	50471-44-8					9.4E+01	4.0E+02			7.6E+01	
				1.00E+00	H	2.00E-01	I	V		2.75E+03	1.36E+09	4.40E+03	1	0.1	Vinyl Acetate	108-05-4					7.8E+04		9.2E+02		9.1E+02	
7.20E-01	I	1.50E-05	P			3.00E-03	I	V		2.47E+03	1.36E+09	1.37E+03	1	0.1	Vinyl Bromide	593-60-2	9.4E-02		2.6E-01	2.6E-01	2.3E+02				4.3E+00	
				4.40E-06	I	5.11E-02	A	V	M	3.92E+03	1.36E+09	9.56E+02	1	0.1	Vinyl Chloride	75-01-4			1.6E-01	5.9E-02	2.3E+02		5.1E+01			