

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) = 0.1							
SFO (mg/kg-day) <sup>1</sup>	k <sub>e</sub> y	IUR (ug/m <sup>3</sup> ) <sup>1</sup>	k <sub>e</sub> y	RD <sub>50</sub> (mg/kg-day)	k <sub>e</sub> y	RIC <sub>1</sub> (mg/m <sup>3</sup> )	k <sub>e</sub> y	v <sub>o</sub> l	mutagen	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	GIABS	ABS <sub>0</sub>	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=0.1 (mg/kg)	Dermal SL THQ=0.1 (mg/kg)	Inhalation SL THQ=0.1 (mg/kg)	Noncarcinogenic SL THQ=0.1 (mg/kg)	
2.20E-06	I	3.00E-04	O	9.00E-03	I	V	1.07E+05	1.36E+09	1	0.1	1.36E+09	8.72E+03	1	0.1	Acetophthaldehyde	30560-19-1			1.1E+01	1.1E+01	2.3E+00	9.9E+00		1.9E+00	
2.00E-02	I	9.00E-01	I	2.00E-03	X	V	1.14E+05	1.36E+09	1	0.1	1.36E+09	1.37E+04	1	0.1	Acetochlor	75-07-0					1.6E+02	6.6E+02		8.2E+00	
9.00E-01	I	2.00E-02	I	6.00E-03	I	V	1.28E+05	1.36E+09	1	0.1	1.36E+09	1.30E+04	1	0.1	Acetone	67-64-1					7.0E+03			7.0E+03	
3.80E+00	C	1.30E-03	C	1.00E-01	I	V	2.52E+03	1.36E+09	1	0.1	1.36E+09	5.97E+04	1	0.1	Acetone Cyanohydrin	75-86-5								2.8E+05	
5.00E-01	I	1.00E-04	I	2.00E-03	I	V	2.27E+04	1.36E+09	1	0.1	1.36E+09	6.91E+03	1	0.1	Acetonitrile	75-05-8								8.1E+01	
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	0.1	Acetophenone	98-86-2	1.8E-01	6.5E-01	2.9E+03	1.4E-01	7.8E+02				7.8E+02	
5.60E-02	C	1.00E-02	I	1.00E-03	I	V	1.11E+05	1.36E+09	1	0.1	1.36E+09	3.42E+04	1	0.1	Acetylaminofluorene, 2-Acrolein	53-96-3	1.8E-01	6.5E-01	2.9E+03	1.4E-01	3.9E+00			1.4E-02	1.4E-02
1.70E+01	I	4.90E-03	I	3.00E-05	I	V	1.42E+03	1.36E+09	1	0.1	1.36E+09	1.58E+03	1	0.1	Acrylamide	79-06-1	3.1E-01	1.2E+00	1.4E+04	2.4E-01	1.6E+01	6.6E+01	8.5E+05	1.3E+01	
2.10E-02	C	6.00E-06	C	1.00E+00	P	5.00E-03	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Acrylic Acid	79-10-7	3.1E-01	1.2E+00	1.4E+04	2.4E-01	3.9E+03			2.0E+00	
2.10E+01	C	6.00E-03	C	4.00E-04	I	9.00E-03	I	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Acrylonitrile	107-13-1	1.3E+00		3.2E-01	2.5E-01	7.0E-01			1.6E+00	
4.00E-02	P	1.60E-06	C	2.00E-03	X	V	1.37E+04	1.36E+09	1	0.1	1.36E+09	2.62E+04	1	0.1	Adiponitrile	111-69-3	1.2E+01	4.4E+01		9.7E+00	7.8E+01	3.3E+02			6.3E+01
4.00E-02	P	2.00E-03	X	4.00E-04	I	3.00E-04	A	1.36E+09	0.15	0.1	1.36E+09	0.15	0.1	Alachlor	15972-60-8	1.2E+01	4.4E+01		9.7E+00	7.8E+01	3.3E+02			6.3E+00	
1.50E+00	I	4.30E-03	I	3.00E-04	I	1.50E-05	C	1.36E+09	1	0.03	1.36E+09	1.36E+09	1	0.1	Aldicarb	116-06-3	3.3E+01		7.4E-01	7.2E-01	7.8E+03			7.1E+05	
8.80E-01	C	2.50E-04	C	4.00E-04	I	3.00E-03	X	V	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Aldicarb Sulfone	1646-88-4	3.3E+01		7.4E-01	7.2E-01	7.8E+03			7.7E+03
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Aldicarb Sulfoxide	1646-87-3	3.3E+01		7.4E-01	7.2E-01	7.8E+03			3.1E+00	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Aldrin	309-00-2	4.1E-02		9.8E-01	3.9E-02	2.3E-01			2.3E-01	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	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+01			3.6E-01	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Allyl Chloride	107-05-1	3.3E+01		7.4E-01	7.2E-01	3.1E+01			3.5E-01	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Aluminum	7429-90-5	3.3E+01		7.4E-01	7.2E-01	3.1E+01			1.7E-01	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Aluminum Phosphide	20859-73-8	3.3E+01		7.4E-01	7.2E-01	3.1E+01			7.1E+05	
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Ameltryn	834-12-8	3.3E-02	1.2E-01	6.4E+02	2.6E-02	7.0E+01	3.0E+02			3.1E+00
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Aminobiphenyl, 4-	92-67-1	3.3E-02	1.2E-01	6.4E+02	2.6E-02	7.0E+01	3.0E+02			5.7E+01
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Aminophenol, m-	591-27-5	3.3E-02	1.2E-01	6.4E+02	2.6E-02	7.0E+01	3.0E+02			5.1E+02
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Aminophenol, o-	95-55-6	3.3E-02	1.2E-01	6.4E+02	2.6E-02	7.0E+01	3.0E+02			2.5E+01
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Aminophenol, p-	123-30-8	3.3E-02	1.2E-01	6.4E+02	2.6E-02	7.0E+01	3.0E+02			1.3E+02
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Amitraz	33089-61-1	3.3E-02	1.2E-01	6.4E+02	2.6E-02	7.0E+01	3.0E+02			1.6E+01
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Ammonia	7664-41-7	3.3E-02	1.2E-01	6.4E+02	2.6E-02	7.0E+01	3.0E+02			1.6E+01
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Ammonium Picrate	131-74-8	3.3E-02	1.2E-01	6.4E+02	2.6E-02	7.0E+01	3.0E+02			1.3E+01
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Ammonium Sulfamate	7773-06-0	3.3E-02	1.2E-01	6.4E+02	2.6E-02	7.0E+01	3.0E+02			1.6E+03
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	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+01	2.3E+02	8.2E+00	1.4E+05	1.6E+03
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Aniline	62-53-3	1.2E+02	4.3E+02	2.4E+06	9.5E+01	5.5E+01	2.3E+02	8.2E+00	1.4E+05	1.6E+03
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Anthraquinone, 9,10-	84-65-1	1.7E+01	6.2E+01	2.4E+06	9.5E+01	1.6E+01	6.6E+01	4.3E+04	1.4E+06	1.3E+01
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Antimony (metallic)	7440-36-0	1.7E+01	6.2E+01	2.4E+06	9.5E+01	1.6E+01	6.6E+01	4.3E+04	1.4E+06	1.3E+01
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Antimony Pentoxide	1314-60-9	1.7E+01	6.2E+01	2.4E+06	9.5E+01	1.6E+01	6.6E+01	4.3E+04	1.4E+06	1.3E+01
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Antimony Trioxide	1332-81-6	1.7E+01	6.2E+01	2.4E+06	9.5E+01	1.6E+01	6.6E+01	4.3E+04	1.4E+06	1.3E+01
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Antimony Trioxide	1309-64-4	1.7E+01	6.2E+01	2.4E+06	9.5E+01	1.6E+01	6.6E+01	4.3E+04	1.4E+06	1.3E+01
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	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+00	3.3E+01	2.1E+03	3.5E+00	2.8E+04
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Arsine	7784-42-1	7.7E-01	5.5E+00	8.9E+02	6.8E-01	3.9E+00	3.3E+01	2.1E+03	3.5E+00	2.8E+04
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Asbestos (units in fibers)	1332-21-4	7.7E-01	5.5E+00	8.9E+02	6.8E-01	3.9E+00	3.3E+01	2.1E+03	3.5E+00	2.8E+04
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Asulam	3337-71-1	7.7E-01	5.5E+00	8.9E+02	6.8E-01	3.9E+00	3.3E+01	2.1E+03	3.5E+00	2.8E+04
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.36E+09	1	0.1	1.36E+09	1.36E+09	1	0.1	Atrazine	1912-24-9	3.0E+00	1.1E+01	1.5E+04	2.4E+00	2.3E+01	9.9E+01	1.4E+06	1.9E+01	2.3E+03
1.10E-01	I	3.10E-05	I	1.00E+00	P	7.00E-06	P	1.3																	

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) = 0.1								
SFO (mg/kg-day) <sup>1</sup>	key	IUR (ug/m <sup>3</sup> ) <sup>1</sup>	key	RD <sub>50</sub> (mg/kg-day)	key	RIC <sub>1</sub> (mg/m <sup>3</sup> )	key	vo	mutagen	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	GIABS	ABS <sub>0</sub>	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=0.1 (mg/kg)	Dermal SL THQ=0.1 (mg/kg)	Inhalation SL THQ=0.1 (mg/kg)	Noncarcinogenic SL THQ=0.1 (mg/kg)		
7.90E-03	I	1.10E-06	I	2.00E-02	I	5.00E-03	I	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+02			1.6E+02		
				1.40E-03	I	5.00E-03	I	V		3.59E+03	1.36E+09	1.40E+03	1		Bromomethane	74-83-9					1.1E+01		7.3E-01	6.8E-01		
				5.00E-03	H			V			1.36E+09	1.24E+05	1		Bromophos	2104-96-3					3.9E+01			3.9E+01		
		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+01	2.2E+01		
1.03E-01	O			1.50E-02	O						1.36E+09		1	0.1	Bromoxynil	1689-84-5	6.7E+00	2.4E+01		5.3E+00	1.2E+02	4.9E+02			9.5E+01	
1.03E-01	O			1.50E-02	O						1.36E+09	4.74E+05	1		Bromoxynil Octanoate	1689-99-2	6.7E+00			6.7E+00	1.2E+02				1.2E+02	
6.00E-01	C	3.00E-05	I			2.00E-03	I	V		6.67E+02	1.36E+09	8.66E+02	1		Butadiene, 1,3-	106-99-0	1.2E+00		8.1E-02	7.6E-02				1.8E-01	7.8E-02	
				1.00E-01	I					7.64E+03	1.36E+09	3.00E+04	1		Butanol, N-	71-36-3										
5.00E-04	I			4.00E-01	I	5.00E+00	I	V			1.36E+09	2.87E+04	1		Butyl Alcohol, 1-	75-65-0	1.4E+03			1.4E+03	7.8E+02		3.1E+03	1.5E+04	2.6E+03	
				2.00E+00	P	3.00E+01	P	V		2.13E+04	1.36E+09	2.92E+04	1		Butyl alcohol, sec-	78-92-2					1.6E+04		9.1E+04	1.3E+04		
				5.00E-02	I						1.36E+09	8.63E+04	1		Butylate	2008-41-5					3.9E+02			3.9E+02		
2.00E-04	C	5.70E-08	C								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+03	
3.60E-03	P			3.00E-01	P						1.36E+09		1	0.1	Butylated hydroxytoluene	128-37-0	1.9E+02	6.9E+02		1.5E+02	2.3E+03	9.9E+03			1.9E+03	
				5.00E-02	P			V		1.08E+02	1.36E+09	8.14E+03	1		Butylbenzene, n-	104-51-8					3.9E+02				3.9E+02	
				1.00E-01	X			V		1.45E+02	1.36E+09	7.35E+03	1		Butylbenzene, sec-	135-98-8					7.8E+02				7.8E+02	
				1.00E-01	X			V		1.83E+02	1.36E+09	7.36E+03	1		Butylbenzene, tert-	98-06-6					7.8E+02				7.8E+02	
				2.00E-02	A						1.36E+09		1	0.1	Cacodylic Acid	75-60-5					1.6E+02	6.6E+02			1.3E+02	
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-01	8.2E+00	1.4E+03		7.1E-01	
1.80E-03	I	1.00E-04	A	1.00E-05	A						1.36E+09		0.05	0.001	Cadmium (Water)	7440-43-9										
				5.00E-01	I	2.20E-03	C				1.36E+09		1	0.1	Caprolactam	105-60-2					3.9E+03	1.6E+04	3.1E+05			3.1E+03
1.50E-01	C	4.30E-05	C			2.00E-03	I				1.36E+09		1	0.1	Captafol	2425-06-1	4.6E+00	1.6E+01	8.9E+04	3.6E+00	1.6E+01	6.6E+01			1.3E+01	
2.30E-03	C	6.60E-07	C			1.30E-01	I				1.36E+09		1	0.1	Captan	133-06-2	3.0E+02	1.1E+03	5.8E+06	2.4E+02	1.0E+03	4.3E+03			8.2E+02	
				1.00E-01	I						1.36E+09		1	0.1	Carbaryl	63-25-2					7.8E+02	3.3E+03			6.3E+02	
				5.00E-03	I						1.36E+09		1	0.1	Carbaryl	1563-66-2					3.9E+01	1.6E+02			3.2E+01	
				1.00E-01	I	7.00E-01	I	V		7.38E+02	1.36E+09	1.17E+03	1		Carbon Disulfide	75-15-0					7.8E+02		8.5E+01		7.7E+01	
7.00E-02	I	6.00E-06	I	4.00E-03	I	1.00E-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+01		1.6E+01		1.0E+01	
				1.00E-02	I					5.89E+03	1.36E+09	6.46E+02	1		Carbonyl Sulfide	463-58-1					7.8E+01		6.7E+00		6.7E+00	
				1.00E-01	I						1.36E+09		1	0.1	Carbosulfan	55285-14-8					7.8E+01	3.3E+02			6.3E+01	
				1.00E-01	I						1.36E+09		1	0.1	Carboxin	5234-68-4					7.8E+02	3.3E+03			6.3E+02	
				1.00E-01	I	9.00E-04	I				1.36E+09		1		Ceric oxide	1306-38-3							1.3E+05		1.3E+05	
				1.50E-02	I						1.36E+09	1.45E+05	1		Chloral Hydrate	302-17-0					7.8E+02				7.8E+02	
				5.00E-01	I						1.36E+09		1	0.1	Chloramben	138-90-4					1.2E+02	4.9E+02			9.5E+01	
4.03E-01	H			5.00E-04	G						1.36E+09		1	0.1	Chloramines, Organic	E701825										
				5.00E-04	G						1.36E+09		1	0.1	Chloranil	118-75-2	1.7E+00	6.1E+00		1.3E+00						
3.50E-01	I	1.00E-04	I	5.00E-04	G						1.36E+09	1.49E+06	1	0.04	Chlordane (alpha)	5103-71-9					3.9E+00	4.1E+01			3.6E+00	
				5.00E-04	G						1.36E+09	1.49E+06	1	0.04	Chlordane (gamma)	5103-74-2					3.9E+00	4.1E+01			3.6E+00	
				5.00E-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+00	4.1E+01	1.1E+02		3.5E+00	
1.00E+01	I	4.60E-03	C	3.00E-04	I						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+00	9.9E+00			1.9E+00	
				7.00E-04	A						1.36E+09		1	0.1	Chlorfenvinphos	470-90-6					5.5E+00	2.3E+01			4.4E+00	
				9.00E-02	O						1.36E+09		1	0.1	Chlorimuron, Ethyl-	90982-32-4					7.0E+02	3.0E+03			5.7E+02	
				1.00E-01	I	1.45E-04	A	V		2.78E+03	1.36E+09	1.22E+03	1		Chlorine	7782-50-5					7.8E+02		1.8E-02		1.8E-02	
				3.00E-02	I	2.00E-04	I	V			1.36E+09		1		Chlorine Dioxide	10049-04-4					2.3E+02		2.8E+04		2.3E+02	
				3.00E-02	I						1.36E+09		1		Chlorite (Sodium Salt)	7758-19-2					2.3E+02				2.3E+02	
4.60E-01	H	3.00E-04	I	2.00E-02	H	5.00E+01	I	V	M	1.15E+03	1.36E+09	1.03E+03	1		Chloro-1,1-difluoroethane, 1-	75-68-3					1.6E+02		5.4E+03		5.4E+03	
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+02		2.2E+00	2.2E+00		
2.70E-01	X			3.50E-03	C						1.36E+09		1	0.1	Chloro-2-methylaniline HCl, 4-	3165-93-3					1.2E+00				1.2E+00	
				3.50E-03	C					1.18E+04	1.36E+09	1.62E+04	1		Chloro-2-methylaniline, 4-	95-69-2	7.0E+00	2.5E+01	5.0E+04	5.4E+00	2.3E+01	9.9E+01			1.9E+01	
				3.50E-03	C						1.36E+09		1	0.1	Chloroacetaldehyde, 2-	107-20-0	2.6E+00			2.6E+00						2.6E+00
				3.00E-05	I						1.36E+09		1	0.1	Chloroacetic Acid	79-11-8					2.7E+01	1.2E+02			2.2E+01	
2.00E-01	P			5.00E-04	P						1.36E+09		1	0.1	Chloroacetophenone, 2-	532-27-4	3.5E+00	1.2E+01		2.7E+00	3.9E+00	1.6E+01		4.3E+03	4.3E+03	
				2.00E-02	I	5.00E-02	P	V		7.61E+02	1.36E+09	6.45E+03	1		Chloroaniline, p-	106-47-8					1.6E+02				2.8E+01	
				1.00E-01	X						1.36E+09		1	0.1	Chlorobenzene	108-90-7					3.9E+00	1.6E+01			3.2E+00	
1.10E-01	C	3.10E-05	C	2.00E-02	I						1.36E+09		1													

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) = 0.1						
SFO (mg/kg-day) <sup>1</sup>	key <sup>1</sup>	IUR (ug/m <sup>3</sup> ) <sup>1</sup>	key <sup>1</sup>	RD <sub>50</sub> (mg/kg-day)	key <sup>1</sup>	RIC <sub>1</sub> (mg/m <sup>3</sup> )	key <sup>1</sup>	vo <sup>1</sup>	mutagen	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	GIABS	ABS <sub>0</sub>	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=0.1 (mg/kg)	Dermal SL THQ=0.1 (mg/kg)	Inhalation SL THQ=0.1 (mg/kg)	Noncarcinogenic SL THQ=0.1 (mg/kg)
				4.00E-02	H						1.36E+09				Copper	7440-50-8					3.1E+02			3.1E+02
				5.00E-02	I	6.00E-01	C				1.36E+09		0.1	Cresol, m-	108-39-4						3.9E+02	1.6E+03	8.5E+07	3.2E+02
				5.00E-02	I	6.00E-01	C				1.36E+09		0.1	Cresol, o-	95-48-7						3.9E+02	1.6E+03	8.5E+07	3.2E+02
				2.00E-02	P	6.00E-01	C				1.36E+09		0.1	Cresol, p-	106-44-5						1.6E+02	6.6E+02	8.5E+07	1.3E+02
				1.00E-01	A						1.36E+09		0.1	Cresol, p-chloro-m-	59-50-7						7.8E+02	3.3E+03		6.3E+02
1.90E+00	H			1.00E-01	A	6.00E-01	C				1.36E+09		0.1	Cresols	1319-77-3						7.8E+02	3.3E+03	8.5E+07	6.3E+02
				1.00E-03	P					1.66E+04	1.36E+09	1.89E+04			Crotonaldehyde, trans-	123-73-9	3.7E-01			3.7E-01	7.8E+00			7.8E+00
				1.00E-01	I	4.00E-01	I	V		2.68E+02	1.36E+09	6.21E+03			Cumene	98-82-8					7.8E+02		2.6E+02	1.9E+02
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	2.5E+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+01	6.6E+01		1.3E+01
				1.00E-03	I	9.00E-03	C				1.36E+09		1	-Calcium Cyanide	592-01-8						7.8E+00		1.3E+06	7.8E+00
				5.00E-03	I						1.36E+09		1	-Copper Cyanide	544-92-3						3.9E+01			3.9E+01
				6.00E-04	I	8.00E-04	G	V		9.54E+05	1.36E+09	5.89E+04			-Cyanide (CN-)	57-12-5					4.7E+00		4.9E+00	2.4E+00
				1.00E-03	I						1.36E+09		1	-Cyanogen	460-19-5						7.8E+00			7.8E+00
				9.00E-02	I						1.36E+09		1	-Cyanogen Bromide	506-68-3						7.0E+02			7.0E+02
				5.00E-02	I						1.36E+09		1	-Cyanogen Chloride	506-77-4						3.9E+02			3.9E+02
				6.00E-04	I	8.00E-04	I	V		1.00E+07	1.36E+09	5.22E+04			-Hydrocyan Cyanide	74-90-8					4.7E+00		4.4E+00	2.3E+00
				2.00E-03	I	9.00E-03	C				1.36E+09		0.04	-Potassium Cyanide	151-50-8						1.6E+01		1.3E+06	1.6E+01
				5.00E-03	I						1.36E+09		0.04	-Potassium Silver Cyanide	506-61-6						3.9E+01			3.9E+01
				1.00E-01	I						1.36E+09		0.04	-Silver Cyanide	506-64-9						7.8E+02			7.8E+02
				1.00E-03	I	9.00E-03	C				1.36E+09		1	-Sodium Cyanide	143-33-9						7.8E+00		1.3E+06	7.8E+00
				5.00E-02	I						1.36E+09		1	-Zinc Cyanide	557-21-1						3.9E+02			3.9E+02
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								
				5.00E+00	I	7.00E-01	P	V		5.11E+03	1.36E+09	4.17E+04			Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	3.5E+01	1.2E+02		2.7E+01	1.6E+02	6.6E+02	6.5E+02	1.3E+02
				2.00E-01	I						1.36E+09		1	Cyclohexanone	108-94-1						3.9E+04		3.0E+03	2.8E+03
				5.00E-03	P	1.00E+00	X	V		2.83E+02	1.36E+09	1.46E+03			Cyclohexene	110-83-8					3.9E+01		1.5E+02	3.1E+01
				2.00E-01	I					2.93E+05	1.36E+09	7.46E+04			Cyclohexylamine	108-91-8					1.6E+03			1.6E+03
				2.50E-02	I						1.36E+09		0.1	Cyfluthrin	68359-37-5						2.0E+02	8.2E+02		1.6E+02
				5.00E-01	O						1.36E+09		0.1	Cyromazine	66215-27-8						3.9E+03	1.6E+04		3.2E+03
				3.00E-02	I						1.36E+09		0.1	Dalapon	75-99-0						2.3E+02	9.9E+02		1.9E+02
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+03	4.9E+03			9.5E+02
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+01	2.3E+02	6.5E+01	2.3E+02	4.4E+01
				4.00E-05	I						1.36E+09		0.1	Demeton	8065-48-3						3.1E-01	1.3E+00		2.5E-01
1.20E-03	I			6.00E-01	I						1.36E+09		0.1	Di(2-ethylhexyl)adipate	103-23-1				4.5E+02	4.7E+03	2.0E+04			3.8E+03
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	2.00E-04	P	2.00E-04	I	V	9.79E+02	1.36E+09	3.20E+04			Diazinon	333-41-5				5.5E+00	2.3E+01		4.4E+00
8.00E-01	P	6.00E-03	P								1.36E+09		0.1	Dibromo-3-chloropropane, 1,2-	96-12-8	1.9E-01		5.4E-03	5.3E-03	1.6E+00		6.7E-01		4.7E-01
2.50E-01	C			4.00E-04	X					1.59E+02	1.36E+09	1.93E+04			Dibromoaecetic acid	631-64-1	2.8E+00	9.9E+00		2.2E+00	2.3E+00	9.9E+00		1.9E+00
				1.00E-02	I						1.36E+09		1	Dibromobenzene, 1,3-	108-36-1						3.1E+00			3.1E+00
				1.00E-02	I						1.36E+09		1	Dibromobenzene, 1,4-	106-37-6						7.8E+01			7.8E+01
				2.00E-02	I					8.02E+02	1.36E+09	7.95E+03			Dibromochloromethane	124-48-1	8.3E+00			8.3E+00	1.6E+02			1.6E+02
8.40E-02	I			2.00E-02	I					1.34E+03	1.36E+09	8.64E+03			Dibromomethane, 1,2-	106-93-4	3.5E-01		4.0E-02	3.6E-02	7.0E+01		8.1E+00	7.3E+00
2.00E+00	I	6.00E-04	I			4.00E-03	X	V		2.82E+03	1.36E+09	5.64E+03			Dibromomethane (Methylene Bromide)	74-95-3					7.0E+01		2.4E+00	2.4E+00
				3.00E-04	P						1.36E+09		0.1	Dibutyltin Compounds	E1790661						2.3E+00	9.9E+00		1.9E+00
				3.00E-02	I						1.36E+09		0.1	Dicamba	1918-00-9						2.3E+02	9.9E+02		1.9E+02
				4.20E-03	P					5.54E+02	1.36E+09	3.21E+03			Dichloramine	3400-09-7					2.3E+02	9.9E+02		1.9E+02
				4.20E-03	P					5.19E+02	1.36E+09	1.11E+04			Dichloro-2-butene, 1,4-	764-41-0			2.1E-03	2.1E-03				
				4.20E-03	P					7.60E+02	1.36E+09	1.11E+04			Dichloro-2-butene, cis-1,4-	1476-11-5			7.4E-03	7.4E-03				
				5.00E-02	I						1.36E+09		0.1	Dichloro-2-butene, trans-1,4-	110-57-6			7.4E-03	7.4E-03					
				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+01	1.3E+02		2.5E+01
				9.00E-02	I	2.00E-01	H	V		3.76E+02	1.36E+09	1.17E+04			Dichlorobenzene, 1,2-	95-50-1					7.0E+02		2.4E+02	1.8E+02
5.40E-03	C	1.10E-05	C			7.00E-02	A	8.00E-01	I	V	1.36E+09	1.04E+04			Dichlorobenzene, 1,4-	106-46-7	1.3E+02		2.7E+00	2.6E+00	5.5E+02		8.7E+02	3.4E+02
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		7.0E+01	3.0E+02		5.7E+01
				9.00E-03	X						1.36E+09		0.1	Dichlorobenzophenone, 4,4'-	90-98-2						1.6E+03	3.0E+02		8.8E+00
				2.00E-01	I	1.00E-01	X	V		8.45E+02	1.36E+09	8.41E+02			Dichlorodifluoromethane	75-71-8					7.0E+01			7.7E+00
2.40E-01	I	6.90E-05	C			5.00E-04	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) = 0.1					
SFO (mg/kg-day) <sup>1</sup>	k e y	IUR (ug/m <sup>3</sup> ) <sup>1</sup>	k e y	RD <sub>50</sub> (mg/kg-day)	k e y	RIC <sub>1</sub> (mg/m <sup>3</sup> )	k e y	v o l	mutagen	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	GIABS	ABS <sub>0</sub>	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=0.1 (mg/kg)	Dermal SL THQ=0.1 (mg/kg)	Inhalation SL THQ=0.1 (mg/kg)	Noncarcinogenic SL THQ=0.1 (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		420-45-1	1.6E+01		2.7E+01	9.9E+00				2.4E+03	2.4E+03		
				8.00E-02	I	7.00E-01	P	V		2.26E+03	1.36E+09	3.06E+03	1		1445-75-6					6.3E+02	7.2E+02		6.3E+02			
				2.18E-02	O					5.30E+02	1.36E+09	3.81E+04	1		55290-64-7					1.7E+02	7.2E+01		1.4E+02			
				2.20E-03	O						1.36E+09		0.1		60-51-5					1.7E+01	7.3E+01		1.4E+01			
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	119-90-4	9.6E-02	3.7E-01		7.6E-02			6.2E-02	6.2E-02			
1.70E-03	P			6.00E-02	P						1.36E+09		0.1	75-18-3	4.1E+02	1.5E+03	3.2E+02		4.7E+02	2.0E+03		3.8E+02				
4.60E+00	C	1.30E-03	C								1.36E+09		0.1	60-11-7	1.5E-01	5.4E-01	2.9E+03	1.2E-01								
5.80E-01	H										1.36E+09		0.1	21436-96-4	1.2E+00	4.3E+00		9.4E-01								
2.00E-01	P			2.00E-03	X						1.36E+09		0.1	95-68-1	3.5E+00	1.2E+01		2.7E+00								
2.70E-02	P			2.00E-03	I			V	M	8.30E+02	1.36E+09	3.13E+04	1	0.1	121-69-7	2.6E+01			2.6E+01			1.6E+01	6.6E+01	1.3E+01		
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	119-93-7	1.4E-02	5.4E-02		1.1E-02			7.8E+02	4.0E+02	2.6E+02		
				1.00E-04	X	2.00E-06	X	V		1.72E+05	1.36E+09	2.77E+04	1		68-12-2					7.8E-01		5.8E-03	5.7E-03			
5.50E+02	C	1.60E-01	C							1.89E+05	1.36E+09	1.68E+05	1	0.1	540-73-8	1.3E-03		2.9E-03	8.8E-04			1.6E+02	6.6E+02	1.3E+02		
				2.00E-02	I						1.36E+09		0.1	105-67-9							4.7E+00	2.0E+01		3.8E+00		
				6.00E-04	I						1.36E+09		0.1	576-26-1							7.8E+00	3.3E+01		6.3E+00		
4.50E-02	C	1.30E-05	C					V		4.73E+02	1.36E+09	5.48E+03	1	0.1	95-05-9	1.5E+01		1.2E+00	1.1E+00							
				1.00E-03	I						1.36E+09		0.1	513-37-1												
				2.00E-03	X						1.36E+09		0.1	534-52-1							6.3E-01	2.6E+00		5.1E-01		
				2.00E-03	X						1.36E+09		0.1	131-89-5							1.6E+01	6.6E+01		1.3E+01		
				4.00E-04	X	2.00E-03	X				1.36E+09		0.1	618-87-1							3.1E+00	1.3E+01	2.8E+05	2.5E+00		
				1.00E-04	P						1.36E+09		0.1	528-29-0							7.8E-01	3.3E+00		6.3E-01		
				1.00E-04	I						1.36E+09		0.1	99-65-0							7.8E-01	3.3E+00		6.3E-01		
				1.00E-04	P						1.36E+09		0.1	100-25-4							7.8E-01	3.3E+00		6.3E-01		
				2.00E-03	I						1.36E+09		0.1	51-28-5							1.6E+01	6.6E+01		1.3E+01		
6.80E-01	I										1.36E+09		0.1	E1615210	1.0E+00	3.6E+00		8.0E-01		2.2E+00	7.8E+00	4.3E+04	1.7E+00	1.6E+01	6.5E+01	1.3E+01
3.10E-01	C	8.90E-05	C								1.36E+09		0.102	121-14-2	4.6E-01	1.7E+00		3.6E-01		2.3E+00	1.0E+01		1.9E+00			
1.50E+00	P			3.00E-04	X						1.36E+09		0.099	606-20-2						2.3E+00	1.0E+01		1.9E+00			
				1.00E-04	X						1.36E+09		0.006	Dinitrotoluene, 2-Amino-4,6-						7.8E-01	5.5E+01		7.7E-01			
				1.00E-04	X						1.36E+09		0.009	Dinitrotoluene, 4-Amino-2,6-						7.8E-01	3.7E+01		7.7E-01			
4.50E-01	X			9.00E-04	X						1.36E+09		0.1	Dinitrotoluene, Technical grade						1.5E+00	5.5E+00		1.2E+00			
				1.00E-03	I						1.36E+09		0.1	88-85-7						7.0E+00	3.0E+01		5.7E+00			
1.00E-01	I	5.00E-06	I			3.00E-02	I	V		1.16E+05	1.36E+09	3.96E+04	1		Dinoseb					7.8E+00	3.3E+01		6.3E+00			
				3.00E-02	I						1.36E+09		0.1	123-91-1			7.0E+00		2.2E+01	5.3E+00	2.3E+02		1.2E+02	8.1E+01		
											1.36E+09		0.03	Dioxins												
6.20E+03	I	1.30E+00	I								1.36E+09		0.03	-Hexachlorodibenzo-p-dioxin, Mixture	34465-46-8	1.1E-04	1.3E-03	2.9E+00	1.0E-04	5.5E-06	7.7E-05	8.2E-03	5.1E-06			
1.30E+05	C	3.80E+01	C			3.00E-02	I				1.36E+09		0.03	-TCDD, 2,3,7,8-	1746-01-6	5.3E-06	6.3E-05	1.4E-04	4.8E-06	2.3E+02	9.9E+02		1.9E+02			
				8.00E-04	X						1.36E+09		0.1	Diphenamid	957-51-7											
				1.00E-01	O						1.36E+09		0.1	Diphenyl Ether	101-84-8								3.4E+00	3.4E+00		
				2.20E-03	I						1.36E+09		0.1	Diphenyl Sulfone	127-63-9						6.3E+00	2.6E+01		5.1E+00		
8.00E-01	I	2.20E-04	I								1.36E+09		0.1	Diphenylamine	122-39-4	8.7E-01	3.1E+00	1.7E+04	6.8E-01	7.8E+02	3.3E+03		6.3E+02			
				2.20E-03	I						1.36E+09		0.1	Diphenylhydrazine, 1,2-	2764-72-9											
7.40E+00	C	2.10E-03	C								1.36E+09		0.1	Diquat	1937-37-7	9.4E-02	3.3E-01	1.8E+03	7.3E-02	1.7E+01	7.3E+01		1.4E+01			
7.40E+00	C	2.10E-03	C								1.36E+09		0.1	Direct Black 38	2602-46-2	9.4E-02	3.3E-01	1.8E+03	7.3E-02							
6.70E+00	C	1.90E-03	C								1.36E+09		0.1	Direct Blue 6	16071-86-6	1.0E-01	3.7E-01	2.0E+03	8.1E-02							
				4.00E-05	I						1.36E+09		0.1	Disulfoton	298-04-4						3.1E-01	1.3E+00		2.5E-01		
				1.00E-02	I			V			1.36E+09	4.54E+04	1	0.1	Dithiane, 1,4-	505-29-3					7.8E+01			7.8E+01		
				2.00E-03	I						1.36E+09		0.1	Diuron	330-54-1						1.6E+01	6.6E+01		1.3E+01		
				2.00E-02	O						1.36E+09		0.1	Dodine	2439-10-3						1.6E+02	6.6E+02		1.3E+02		
				5.00E-02	O			V			1.36E+09	1.17E+05	1	0.1	EPTC	759-94-4					3.9E+02			3.9E+02		
				6.00E-03	I			V			1.36E+09	4.10E+05	1	0.1	Endosulfan	115-29-7					4.7E+01			4.7E+01		
				6.00E-03	P						1.36E+09		0.1	Endosulfan Sulfate	1031-07-8						4.7E+01	2.0E+02		3.8E+01		
				2.00E-02	I						1.36E+09		0.1	Endothall	145-73-3						1.6E+02	6.6E+02		1.3E+02		
				3.00E-04	I						1.36E+09		0.1	Endrin	72-20-8						2.3E+00	9.9E+00		1.9E+00		
9.90E-03	I	1.20E-06	I			6.00E-03	P	1.00E-03	I	V	1.05E+04	1.36E+09	1.89E+04	1		Epichlorohydrin	106-89-8	7.0E+01		4.4E+01	2.7E+01		2.0E+00	1.9E+00		
				4.00E-02	P						1.36E+09		0.1	Epoxycyclohexane, 1,2-	106-88-7						4.7E+01			1.6E+01		
				5.00E-03	I						1.36E+09		0.1	Ethanol, 2-(2-methoxyethoxy)-	111-77-3						3.1E+02	1.3E+03		2.5E+02		

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) = 0.1				
SFO (mg/kg-day) <sup>1</sup>	k <sub>e</sub> y	IUR (ug/m <sup>3</sup> ) <sup>1</sup>	k <sub>e</sub> y	RD <sub>50</sub> (mg/kg-day)	k <sub>e</sub> y	RIC <sub>1</sub> (mg/m <sup>3</sup> )	k <sub>e</sub> y	v <sub>o</sub> I	mutagen	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	GIABS	ABS <sub>0</sub>	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=0.1 (mg/kg)	Dermal SL THQ=0.1 (mg/kg)	Inhalation SL THQ=0.1 (mg/kg)	Noncarcinogenic SL THQ=0.1 (mg/kg)	
				1.30E-02							1.36E+09		1	0.1	Fluometuron	2164-17-2					1.0E+02	4.3E+02			8.2E+01
				4.00E-02	C	1.30E-02	C				1.36E+09		1		Fluoride	16984-48-8					3.1E+02		1.8E+06		3.1E+02
				6.00E-02	I	1.30E-02	C				1.36E+09		1		Fluorine (Soluble Fluoride)	7782-41-4					4.7E+02		1.8E+06		4.7E+02
				8.00E-02	I						1.36E+09		1	0.1	Fluridone	59756-60-4					6.3E+02	2.6E+03			5.1E+02
				4.00E-02	O						1.36E+09		1	0.1	Flurprimidol	56425-91-3					3.1E+02	1.3E+03			2.5E+02
				2.00E-03	O						1.36E+09		1	0.1	Flusilazole	85509-19-9					1.6E+01	6.6E+01			1.3E+01
				5.00E-01	O						1.36E+09		1	0.1	Flutolanil	96332-98-5					3.9E+03	1.6E+04			3.2E+03
				1.00E-02	I						1.36E+09		1	0.1	Fluvalinate	69403-94-5					7.8E+01	3.3E+02			6.3E+01
				9.00E-02	O						1.36E+09		1	0.1	Folpet	133-07-3					7.0E+02	3.0E+03			5.7E+02
				1.00E-02	O						1.36E+09		1	0.1	Fomesafen	72178-02-0					7.8E+01	3.3E+02			6.3E+01
2.10E-02	C	7.40E-06	I	2.00E-03	I	7.00E-03	I	V	M	4.24E+04	1.36E+09	7.77E+04	1	0.1	Fonofos	944-22-9				4.3E+00	1.6E+01	6.6E+01		1.3E+01	
				2.00E-01	P	3.00E-04	X	V		1.06E+05	1.36E+09	9.30E+04	1		Formaldehyde	50-00-0	7.3E+00		1.1E+01		1.6E+03		5.7E+01		5.5E+01
				9.00E-01	P						1.36E+09		1		Formic Acid	64-18-6					7.0E+03		2.9E+00		2.9E+00
				2.50E+00	O						1.36E+09		1	0.1	Fosetyl-AL	39148-24-8					2.0E+04	8.2E+04			1.6E+04
				1.00E-03	X			V			1.36E+09	1.56E+05	1		Furans										
				1.00E-03	I			V		6.22E+03	1.36E+09	2.62E+03	1		-Dibenzofuran	132-64-9					7.8E+00				7.8E+00
				9.00E-01	I	2.00E+00	I	V		1.65E+05	1.36E+09	1.20E+04	1		-Furan	110-00-9					7.8E+00				7.8E+00
3.80E+00	H			3.00E-03	I	5.00E-02	H	V			1.36E+09		1	0.1	-Tetrahydrofuran	109-99-9	1.8E-01	6.5E-01		1.4E-01	7.0E+03		2.5E+03		1.8E+03
				1.00E-01	I			V			1.36E+09		1	0.1	Furazolidone	67-45-8									
1.50E+00	C	4.30E-04	C	3.00E-03	I	5.00E-02	H	V		1.01E+04	1.36E+09	4.86E+04	1	0.1	Furfural	98-01-1					2.3E+01		2.5E+02		2.1E+01
3.00E-02	I	8.60E-06	C	6.00E-03	O						1.36E+09		1	0.1	Furium	531-82-8	4.6E-01	1.6E+00	8.9E+03	3.6E-01					
				1.00E-01	A	8.00E-05	C				1.36E+09		1	0.1	Furmecycloz	60568-05-0	2.3E-01	8.2E+01	4.4E+05	1.8E+01					
				4.00E-04	I	1.00E-03	X	V		1.06E+05	1.36E+09	8.43E+04	1		Glufosinate, Ammonium	77182-82-2					4.7E+01	2.0E+02			3.8E+01
				1.00E-01	I						1.36E+09		1	0.1	Glutaraldehyde	111-30-8					7.8E+02	3.3E+03	1.1E+04		6.0E+02
				1.00E-02	X			V			1.36E+09	1.45E+05	1		Glycidaldehyde	765-34-4					3.1E+00		8.8E+00		2.3E+00
				2.00E-02	P						1.36E+09		1	0.1	Glyphosate	1071-83-6					7.8E+02	3.3E+03			6.3E+02
				3.00E-02	X						1.36E+09		1	0.1	Guanidine	113-00-8					7.8E+01				7.8E+01
				5.00E-05	I						1.36E+09		1	0.1	Guanidine Chloride	50-01-1					1.6E+02	6.6E+02			1.3E+02
4.50E+00	I	1.30E-03	I	1.00E-04	A			V			1.36E+09	4.79E+05	1		Guanidine Nitrate	506-93-4					2.3E+02	9.9E+02			1.9E+02
9.10E+00	I	2.60E-03	I	1.30E-05	I			V		3.00E-03	1.36E+09	8.43E+05	1		Haloxypol, Methyl	69806-40-2	1.5E-01		1.0E+00	1.3E-01	3.9E-01	1.6E+00			3.2E-01
				3.00E-04	X	4.00E-01	P	V		5.79E+01	1.36E+09	8.95E+02	1		Heptachlor	76-44-8	7.6E-02		9.1E-01	7.0E-02	7.8E-01				7.8E-01
				2.00E-03	I			V			1.36E+09	3.80E+05	1		Heptachlor Epoxide	1024-57-3					1.0E-01				1.0E-01
1.60E+00	I	4.60E-04	I	1.00E-05	P			V		2.09E+02	1.36E+09	7.80E+03	1		Heptanal, n-	111-71-7							2.4E+00		
7.80E-02	I	2.20E-05	I	1.00E-03	P			V		1.68E+01	1.36E+09	1.08E+04	1		Heptane, n-	142-92-5					2.3E+00		3.7E+01		2.2E+00
6.30E+00	I	1.80E-03	I	9.00E-04	A						1.36E+09		1	0.1	Hexabromobenzene	87-82-1	4.3E-01		4.1E-01	2.1E-01	1.6E+01				1.6E+01
1.80E+00	I	5.30E-04	I	6.00E-08	X			A			1.36E+09		1	0.1	Hexachlorobutadiene	86631-49-2					1.6E+00	6.6E+00			1.3E+00
				2.00E-04	I						1.36E+09		1	0.1	Hexachlorodiphenyl ether, 2,2',4,4',5,5'-(BDE-153)	118-74-1					7.8E-02				7.8E-02
6.30E+00	I	1.80E-03	I	9.00E-04	A					1.68E+01	1.36E+09	1.08E+04	1		Hexachlorobenzene	87-82-1	8.9E+00		1.4E+00	1.2E+00	7.8E+00		3.0E+01		7.8E+00
1.80E+00	I	5.30E-04	I	6.00E-08	X			A			1.36E+09		1	0.1	Hexachlorocyclohexane, Alpha-	319-84-6	1.1E-01	3.9E-01	2.1E+03	8.6E-02	7.0E+00				5.7E+00
				8.00E-07	A						1.36E+09		1	0.1	Hexachlorocyclohexane, Beta-	319-85-7	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		1	0.04	Hexachlorocyclohexane, Delta-	319-86-8	6.3E-01	5.6E+00	1.2E+04	5.7E-01	4.7E-04	2.0E-03			3.8E-04
1.80E+00	I	5.10E-04	I	6.00E-03	I	2.00E-04	I	V		1.57E+01	1.36E+09	8.51E+03	1		Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	3.9E-01	1.4E+00	7.5E+03	3.0E-01	6.3E-03	6.6E-02			5.7E-03
4.00E-02	I	1.10E-05	C	6.00E-03	I	3.00E-02	I	V			1.36E+09	8.01E+03	1		Hexachlorocyclohexane, Technical	608-73-1									
8.00E-02	I	4.00E-03	I	3.00E-04	I						1.36E+09		1	0.1	Hexachlorocyclopentadiene	77-47-4	1.7E+01		2.0E+00	1.8E+00	4.7E+01		1.8E-01		1.8E-01
				1.00E-05	I	V				3.39E+03	1.36E+09	3.00E+05	1		Hexachloroethane	67-72-1					5.5E+00		2.5E+01		4.5E+00
				4.00E-04	C						1.36E+09		1	0.1	Hexachloroethene	70-30-4					2.3E+00	9.9E+00			1.9E+00
				1.00E-05	I	V				3.39E+03	1.36E+09	3.00E+05	1		Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	8.7E+00	2.1E+02		8.3E+00	3.1E+01		8.8E+02		3.1E-01
				4.00E-04	C						1.36E+09		1	0.1	Hexamethylene Diisocyanate, 1,6-	822-06-0								3.1E-01	3.1E-01
				4.00E-04	P						1.36E+09		1	0.1	Hexamethylene diisocyanate biuret	4035-89-6							5.7E+04		5.7E+04
				6.00E-01	P	V				1.41E+02	1.36E+09	8.29E+02	1		Hexamethylene diisocyanate isocyanurate	3779-63-3					3.1E+00	1.3E+01			5.7E+04
2.00E-07	X			7.00E-01	I	V				1.41E+02	1.36E+09	8.29E+02	1		Hexamethylphosphoramide	680-31-9							1.3E+01		2.5E+00
				2.00E+00	P						1.36E+09		1	0.1	Hexane, Commercial	E5241997			1.2E+01	1.2E+01	3.1E+00		5.2E+01		5.2E+01
9.50E-03	P			7.00E-01	I	V				1.41E+02	1.36E+09	8.29E+02	1		Hexane, n-	110-54-3									

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) = 0.1							
SFO (mg/kg-day) <sup>1</sup>	k <sub>e</sub> y <sup>-1</sup>	IUR (ug/m <sup>3</sup> ) <sup>1</sup>	k <sub>e</sub> y <sup>-1</sup>	RD <sub>50</sub> (mg/kg-day)	k <sub>e</sub> y <sup>-1</sup>	RIC <sub>1</sub> (mg/m <sup>3</sup> )	k <sub>e</sub> y <sup>-1</sup>	v <sub>o</sub> l	mutagen	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	GIABS	ABS <sub>0</sub>	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=0.1 (mg/kg)	Dermal SL THQ=0.1 (mg/kg)	Inhalation SL THQ=0.1 (mg/kg)	Noncarcinogenic SL THQ=0.1 (mg/kg)	
				1.87E-05	P						1.36E+09		1	Lanthanum Chloride Heptahydrate	10025-84-0						1.5E-01			1.5E-01	
				2.83E-05	P						1.36E+09		1	Lanthanum Chloride, Anhydrous	10099-58-8						2.2E-01			2.2E-01	
				1.60E-05	P						1.36E+09		1	Lanthanum Nitrate Hexahydrate	10277-43-7						1.3E-01			1.3E-01	
8.50E-03	C	1.20E-05	C								1.36E+09		1	Lead Compounds		8.2E+01		3.2E+05	8.2E+01						
2.10E-01	C	8.00E-05	C								1.36E+09		0.1	-Lead Phosphate	7446-27-7	3.3E+00	1.2E+01	4.8E+04	2.6E+00					2.0E+02	
											1.36E+09		1	-Lead acetate	301-04-2									1.0E+02	
											1.36E+09		1	-Lead and Compounds	7439-92-1										
											1.36E+09		1	-Lead and Compounds (with other sources of lead present, see Guidance)	7439-92-1										
3.80E-02	C	1.10E-05	C								1.36E+09		0.1	-Lead subacetate	1335-32-6	1.8E+01	6.5E+01	3.5E+05	1.4E+01						
				1.00E-07	I		V			2.43E+00	1.36E+09	1.91E+03	1	-Tetraethyl Lead	78-00-2						7.8E-04			7.8E-04	
				5.00E-06	P		V			3.83E+02	1.36E+09	2.55E+04	1	Lewisite	541-25-3						3.9E-02			3.9E-02	
				7.70E-03	O						1.36E+09		0.1	Linuron	330-55-2						6.0E+01	2.5E+02		4.9E+01	
				2.00E-03	P						1.36E+09		1	Lithium	7439-93-2						1.6E+01			1.6E+01	
				5.00E-04	I						1.36E+09		0.1	MCPA	94-74-6						3.9E+00	1.6E+01		3.2E+00	
				4.40E-02	O						1.36E+09		0.1	MCPB	94-81-5						3.4E+02	1.5E+03		2.8E+02	
				1.00E-03	I						1.36E+09		0.1	MCPP	93-65-2						7.8E+00	3.3E+01		6.3E+00	
				2.00E-02	I						1.36E+09		0.1	Malathion	121-75-5						1.6E+02	6.6E+02		1.3E+02	
				1.00E-01	I	7.00E-04	C				1.36E+09		0.1	Maleic Anhydride	108-31-6						7.8E+02	3.3E+03	9.9E+04	6.3E+02	
				5.00E-01	I						1.36E+09		0.1	Maleic Hydrazide	123-33-1						3.9E+03	1.6E+04		3.2E+03	
				1.00E-04	P						1.36E+09		0.1	Malononitrile	109-27-3						7.8E-01	3.3E+00		6.3E-01	
				3.00E-02	H						1.36E+09		0.1	Mancozeb	8018-01-7						2.3E+02	9.9E+02		1.9E+02	
				5.00E-03	I						1.36E+09		0.1	Maneb	12427-38-2						3.9E+01	1.6E+02		3.2E+01	
				1.40E-01	I	5.00E-05	I				1.36E+09		1	Manganese (Diet)	7439-96-5										
				2.40E-02	G	5.00E-05	I		0.04		1.36E+09		1	Manganese (Non-diet)	7439-96-5						1.9E+02		7.1E+03	1.8E+02	
				9.00E-05	H						1.36E+09		0.1	Meposfolan	950-10-7						7.0E-01	3.0E+00		5.7E-01	
				3.00E-02	I						1.36E+09		0.1	Mepiquat Chloride	24307-26-4						2.3E+02	9.9E+02		1.9E+02	
1.10E-02	P			4.00E-03	P						1.36E+09		0.1	Mercaptobenzothiazole, 2-Mercury Compounds	149-30-4	6.3E+01	2.2E+02		4.9E+01		3.1E+01	1.3E+02		2.5E+01	
				3.00E-04	I	3.00E-04	G		0.07		1.36E+09		1	-Mercuric Chloride (and other Mercury salts)	7487-94-7						2.3E+00		4.3E+04	2.3E+00	
											1.36E+09	2.27E+04	1	-Mercury (elemental)	7439-97-6							7.1E-01		7.1E-01	
				1.00E-04	I						1.36E+09		1	-Methyl Mercury	22967-92-6						7.8E-01			7.8E-01	
				8.00E-05	I						1.36E+09		0.1	-Phenylmercuric Acetate	62-38-4						6.3E-01	2.6E+00		5.1E-01	
				3.00E-05	I	V				1.94E+06	1.36E+09		1	Mephos	159-50-5						7.8E-01			6.3E-01	
				6.00E-02	I						1.36E+09		0.1	Metaxyl	57837-19-1						4.7E+02	2.0E+03		3.8E+02	
				1.00E-04	I	3.00E-02	P V			4.58E+03	1.36E+09	6.79E+03	1	Methacrylonitrile	126-98-7						7.8E-01		2.1E+01	7.5E-01	
				5.00E-05	I						1.36E+09		0.1	Methamidophos	10265-92-6						3.9E-01	1.6E+00		3.2E-01	
				2.00E+00	I	2.00E+01	I V			1.06E+05	1.36E+09	2.90E+04	1	Methanol	67-56-1						1.6E+04		6.1E+04	1.2E+04	
				1.50E-03	O						1.36E+09		0.1	Methidathion	950-37-8						1.2E+01	4.9E+01		9.5E+00	
				2.50E-02	I						1.36E+09		0.1	Methomyl	16752-77-5						2.0E+02	8.2E+02		1.6E+02	
4.90E-02	C			5.00E-03	I						1.36E+09		0.1	Methoxy-5-nitroaniline, 2-Methoxychlor	99-59-2	1.4E+01	5.0E+01		1.1E+01		3.9E+01	1.6E+02		3.2E+01	
											1.36E+09		0.1	Methoxyethanol Acetate, 2-Methoxyethanol, 2-Methyl Acetate	110-49-6						6.3E+01		1.3E+01	1.1E+01	
				5.00E-03	P	1.00E-03	P V			1.15E+05	1.36E+09	1.24E+05	1	Methyl Acrylate	109-86-4						3.9E+01		7.4E+01	2.6E+01	
				1.00E+00	X					2.90E+04	1.36E+09	8.12E+03	1	Methyl Acrylate	79-20-9						7.8E+03			7.8E+03	
						2.00E-02	P V			6.75E+03	1.36E+09	6.97E+03	1	Methyl Ethyl Ketone (2-Butanone)	96-33-3									1.5E+01	
				6.00E-01	I	5.00E+00	I V			2.84E+04	1.36E+09	1.22E+04	1	Methyl Hydrazine	78-93-3						4.7E+03		6.4E+03	2.7E+03	
1.00E-03	X			1.00E-03	P	2.00E-05	X V			1.80E+05	1.36E+09	5.04E+04	1	Methyl Isobutyl Ketone (4-methyl-2-pentanone)	60-34-4			1.4E-01	1.4E-01		7.8E+00		1.1E-01	1.0E-01	
						3.00E+00	I V			3.36E+03	1.36E+09	1.06E+04	1	Methyl Isocyanate	108-10-1						6.3E+01		3.3E+03	3.3E+03	
						1.00E-03	C V			1.01E+04	1.36E+09	4.42E+03	1	Methyl Methacrylate	624-83-9						1.1E+04		4.6E+01	4.6E-01	
				1.40E+00	I	7.00E-01	I V			2.36E+03	1.36E+09	6.33E+03	1	Methyl Parathion	80-62-6						4.7E+02	2.0E+03		4.4E+02	
				2.50E-04	I						1.36E+09		0.1	Methyl Phosphonic Acid	298-00-0						2.0E+00	8.2E+00		1.6E+00	
				6.00E-02	X						1.36E+09		0.1	Methyl Styrene (Mixed Isomers)	993-13-5						4.7E+02	2.0E+03		3.8E+02	
				6.00E-03	H	4.00E-02	H V			3.93E+02	1.36E+09	2.43E+04	1	Methyl methanesulfonate	25013-15-4						4.7E+01		1.0E+02	3.2E+01	
9.90E-02	C	2.80E-05	C								1.36E+09		0.1	Methyl tert-Butyl Ether (MTBE)	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	4.90E+03	1	Methyl-1,4-benzenediamine dihydrochloride, 2-Methyl-2-Pentanol, 4-Methyl-5-Nitroaniline, 2-Methyl-N-nitro-N-nitrosoguanidine, N-Methylamine Hydrochloride, 2-Methylarsinic acid	1634-04-4	3.9E+02		5.3E+01	4.7E+01						
				3.00E-04	X						1.36E+09		0.1	Methylbenzene, 1,4-diamine monohydrochloride, 2-Methylbenzene-1,4-diamine monohydrochloride, 2-Methylchlorohexane	615-45-2						2.3E+00	9.9E+00		1.5E+03	
											1.36E+09	1.72E+04	1	Methylbenzene, 1,4-diamine monohydrochloride, 2-Methylbenzene-1,4-diamine monohydrochloride, 2-Methylchlorohexane	108-11-2									5.4E+03	
9.00E-03	P			2.00E-02	X						1.36E+09		0.1	Methylchlorohexane	99-55-8	7.7E+01	2.7E+02		6.0E+01		1.6E+02	6.6E+02		5.4E+03	
8.30E+00	C	2.40E-03	C								1.36E+09		0.1	Methylchlorohexane	70-25-7	8.4E-02	3.0E-01	1.6E+03	6.5E-02						

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) = 0.1						
SFO (mg/kg-day) <sup>1</sup>	k <sub>e</sub> y <sup>-1</sup>	IUR (ug/m <sup>3</sup> ) <sup>1</sup>	k <sub>e</sub> y <sup>-1</sup>	RD <sub>50</sub> (mg/kg-day)	k <sub>e</sub> y <sup>-1</sup>	RIC <sub>1</sub> (mg/m <sup>3</sup> )	k <sub>e</sub> y <sup>-1</sup>	v <sub>o</sub> l/m <sup>3</sup>	mutagen	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	GIABS	ABS <sub>0</sub>	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=0.1 (mg/kg)	Dermal SL THQ=0.1 (mg/kg)	Inhalation SL THQ=0.1 (mg/kg)	Noncarcinogenic SL THQ=0.1 (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	300-76-5					1.6E+01				1.6E+01
				1.20E-01	O	1.40E-05	C				1.36E+09	1	0.1	Naphtha, High Flash Aromatic (HFAN)	64742-95-6	3.9E-01	1.4E+00		3.0E-01	1.6E+01			1.4E+07	2.3E+02
				2.60E-04	C	1.10E-02	C	1.40E-05	C		1.36E+09	1	0.1	Naphthylamine, 2-	91-59-8					1.6E+01				1.6E+01
				2.60E-04	C	1.10E-02	C	1.40E-05	C		1.36E+09	1	0.1	Napropamide	15299-99-7			1.5E+04	1.5E+04	9.4E+02	4.0E+03	3.6E+02	2.0E+03	7.6E+02
				2.60E-04	C	1.10E-02	C	1.40E-05	C		1.36E+09	1	0.1	Nickel Acetate	373-02-4			1.5E+04	1.5E+04	8.6E+01	3.6E+02	3.6E+02	2.0E+03	6.7E+01
				2.60E-04	C	1.10E-02	C	1.40E-05	C	V	1.36E+09	1		Nickel Carbonate	3333-67-3			1.5E+04	1.5E+04	8.6E+01	3.6E+02	3.6E+02	2.0E+03	6.7E+01
				2.60E-04	C	1.10E-02	C	1.40E-05	C		1.36E+09	0.04		Nickel Carbonyl	13463-39-3			1.5E+04	1.5E+04	8.6E+01	3.6E+02	3.6E+02	2.0E+03	8.2E+01
				2.60E-04	C	1.10E-02	C	2.00E-05	C		1.36E+09	0.04		Nickel Hydroxide	12054-49-7			1.5E+04	1.5E+04	8.6E+01	3.6E+02	3.6E+02	2.0E+03	8.2E+01
				2.40E-04	I	1.10E-02	C	1.40E-05	C		1.36E+09	0.04		Nickel Oxide	1313-99-1			1.5E+04	1.5E+04	8.6E+01	3.6E+02	3.6E+02	2.0E+03	8.4E+01
				2.60E-04	C	2.00E-02	I	1.00E-05	T		1.36E+09	0.04		Nickel Refinery Dust	E715532			1.6E+04	1.6E+04	8.6E+01	3.6E+02	3.6E+02	2.0E+03	8.2E+01
1.70E+00	C	4.80E-04	I	1.10E-02	C	1.40E-05	C				1.36E+09	0.04		Nickel Soluble Salts	7440-02-0	4.1E-01	2.7E+00	8.0E+03	4.1E-01	1.6E+01	1.6E+01	2.0E+03	1.4E+03	1.4E+02
9.10E-01	C	2.60E-04	C	1.10E-02	C	1.40E-05	C				1.36E+09	0.04	0.1	Nickel Sub sulfide	12035-72-2	7.6E-01	2.7E+00	8.0E+03	4.1E-01	8.6E+01	3.6E+02	3.6E+02	2.0E+03	8.2E+01
				1.60E+00	I						1.36E+09	1	0.1	Nickelocene	1271-28-9			1.5E+04	6.0E-01	8.6E+01	3.6E+02	3.6E+02	2.0E+03	6.7E+01
				1.00E-01	I						1.36E+09	1		Nitrate (measured as nitrogen)	14797-55-8					1.3E+04				1.3E+04
				1.00E-02	X	5.00E-05	X				1.36E+09	1	0.1	Nitrate + Nitrite (measured as nitrogen)	E701177					7.8E+02				7.8E+02
2.00E-02	P			4.00E-03	P	6.00E-03	P				1.36E+09	1	0.1	Nitro (measured as nitrogen)	14797-65-0					7.8E+02	3.3E+02	7.1E+03	8.5E+05	2.5E+01
				1.00E-01	I						1.36E+09	1	0.1	Nitroaniline, 2-	88-74-4	3.5E+01	1.2E+02		2.7E+01	7.8E+02	3.3E+02	7.1E+03	8.5E+05	2.5E+01
				4.00E-05	I	2.00E-03	I	9.00E-03	I	V	3.05E+03	7.32E+04	1	0.1	Nitroaniline, 4-	100-01-6			5.1E+00	5.1E+00	3.1E+01	1.3E+02	8.5E+05	2.5E+01
				3.00E+03	P						1.36E+09	1	0.1	Nitrobenzene	98-95-3			5.1E+00	5.1E+00	1.6E+01	9.9E+07	6.9E+01	1.9E+07	1.3E+01
				7.00E-02	H						1.36E+09	1	0.1	Nitrocellulose	9004-70-0					2.3E+07	5.5E+02	2.3E+03	4.4E+02	1.9E+07
1.30E+00	C	3.70E-04	C								1.36E+09	1	0.1	Nitrofurantoin	67-20-9			1.0E+04	4.2E-01	5.3E-01	1.9E+00	1.0E+04	4.2E-01	4.2E-01
1.70E-02	P			1.00E-04	P						1.36E+09	1	0.1	Nitrofurazone	59-87-0	4.1E+01	1.5E+02	1.0E+04	3.2E+01	7.8E-01	3.3E+00	3.3E+00	3.3E+00	6.3E-01
				1.00E-01	I						1.36E+09	1	0.1	Nitroglycerin	55-63-0					7.8E-01	3.3E+00	3.3E+00	3.3E+00	6.3E-01
				1.00E-01	I						1.36E+09	1	0.1	Nitroquinidine	556-88-7					7.8E+02	3.3E+03	3.3E+03	3.3E+03	6.3E+02
				8.80E-06	P	5.00E-03	P	V		1.80E+04	1.36E+09	1.69E+04	1	0.1	Nitromethane	75-52-5			5.4E+00	5.4E+00	6.4E-02	6.4E-02	6.4E-02	8.8E+00
2.70E+01	C	7.70E-03	C	5.80E-04	X	2.00E-02	I	V		4.86E+03	1.36E+09	1.31E+04	1	0.1	Nitropropane, 2-	79-46-9			6.4E-02	6.4E-02	6.4E-02	6.4E-02	6.4E-02	2.7E+01
				3.40E-02	C				M		1.36E+09	1	0.1	Nitroso-N-ethylurea, N-	759-73-9	5.7E-03	2.2E-02	1.8E+02	4.5E-03					
5.40E+00	I	1.60E-03	I								1.36E+09	2.43E+05	1	0.1	Nitroso-N-methylurea, N-	684-93-5	1.3E-03	5.0E-03	4.1E+01	1.0E-03	1.3E-01	4.3E-01	9.9E-02	
2.80E+00	I	8.00E-04	C								1.36E+09	1	0.1	Nitrosodibutylamine, N-	924-16-3	1.3E-01	4.3E-01	9.9E-02	9.9E-02	2.5E-01	8.8E-01	4.8E+03	1.9E-01	
				1.50E+02	I	4.30E-02	I				1.36E+09	1	0.1	Nitrosodiethanolamine, N-	1116-54-7	2.5E-01	8.8E-01	4.8E+03	1.9E-01					
5.10E+01	I	1.40E-02	I	8.00E-06	P	4.00E-05	X	V	M	2.37E+05	1.36E+09	8.23E+04	1	0.1	Nitrosodimethylamine, N-	55-18-5	1.0E-03	4.0E-03	3.2E+01	8.1E-04				6.3E-02
4.90E-03	I	2.60E-06	C								1.36E+09	1	0.1	Nitrosodiphenylamine, N-	62-75-9	3.0E-03	6.0E-03	2.0E-03	2.0E-03	1.4E+02	5.0E+02	1.5E+06	1.1E+02	
7.00E+00	I	2.00E-03	C								1.36E+09	1	0.1	Nitrosodipropylamine, N-	95-30-6	1.4E+02	5.0E+02	1.5E+06	1.1E+02					
2.20E+01	I	6.30E-03	C								1.36E+09	1.21E+05	1	0.1	Nitrosodipropylamine, N-	621-64-7	9.9E-02	3.5E-01	1.9E+03	7.8E-02				
6.70E+00	C	1.90E-03	C								1.36E+09	1	0.1	Nitrosomethylthylamine, N-	10595-95-6	3.2E-02	5.4E-02	2.0E-02	2.0E-02					
9.40E+00	C	2.70E-03	C								1.36E+09	1	0.1	Nitrosomorpholine [N-]	59-89-2	1.0E-01	3.7E-01	2.0E+03	8.1E-02					
2.10E+00	I	6.10E-04	I								1.36E+09	1	0.1	Nitrosopiperidine [N-]	100-75-4	7.4E-02	2.6E-01	1.4E+03	5.8E-02					
				1.00E-04	X						1.36E+09	1	0.1	Nitrosopyrrolidine, N-	930-55-2	3.3E-01	1.2E+00	6.3E+03	2.6E-01					
2.20E-01	P			9.00E-04	P			V		1.51E+03	1.36E+09	1.37E+05	1	0.1	Nitrotoluene, o-	88-72-2	3.2E+00			3.2E+00	7.0E+00			7.0E+00
1.60E-02	P			4.00E-03	P						1.36E+09	1	0.1	Nitrotoluene, p-	99-99-0	4.3E+01	1.5E+02		3.4E+01	3.1E+01	1.3E+02	1.3E+02	2.5E+01	2.5E+01
				3.00E-04	X	2.00E-02	P	V		6.86E+00	1.36E+09	1.04E+03	1	0.1	Nonane, n-	111-84-2					2.3E+00		2.2E+00	1.1E+00
				1.50E-03	O						1.36E+09	1	0.1	Norflurazon	27314-13-2					1.2E+01	4.9E+01	4.9E+01	4.9E+01	9.5E+00
				3.00E-03	I						1.36E+09	1	0.1	Octabromodiphenyl Ether	32536-52-0					2.3E+01	9.9E+01	9.9E+01	9.9E+01	1.9E+01
				5.00E-02	I						1.36E+09	1	0.006	Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0					3.9E+02	2.7E+04	2.7E+04	2.7E+04	3.9E+02
7.79E-03	O			2.00E-03	H						1.36E+09	1	0.1	Octamethylphosphoramide	152-18-9					1.6E+01	6.6E+01	6.6E+01	6.6E+01	1.3E+01
				1.90E-01	O						1.36E+09	1	0.1	Oryzalin	19044-88-3	8.9E+01	3.2E+02		7.0E+01	1.5E+03	6.3E+03	6.3E+03	6.3E+03	1.2E+03
				5.00E-03	I						1.36E+09	1	0.1	Oxadiazon	19666-30-9					3.9E+01	1.6E+02	1.6E+02	1.6E+02	3.2E+01
				2.50E-02	I						1.36E+09	1	0.1	Oxamyl	23135-22-0					2.0E+02	8.2E+02	8.2E+02	8.2E+02	1.6E+02
				4.00E-02	O						1.36E+09	1	0.1	Oxylufen	42874-03-3	9.5E+00	3.4E+01		7.4E+00	3.1E+02	1.3E+03	1.3E+03	2.5E+02	2.5E+02
				1.30E-02	I						1.36E+09	1	0.1	Paclitaxel	76738-62-0					1.0E+02				

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) = 0.1					
SFO (mg/kg-day) <sup>1</sup>	key	IUR (ug/m <sup>3</sup> ) <sup>1</sup>	key	RD <sub>50</sub> (mg/kg-day)	key	RIC <sub>1</sub> (mg/m <sup>3</sup> )	key	vo	mutagen	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	GIABS	ABS <sub>0</sub>	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 TH=0.1 (mg/kg)	Dermal SL TH=0.1 (mg/kg)	Inhalation SL TH=0.1 (mg/kg)	Noncarcinogenic SL TH=0.1 (mg/kg)
		5.00E-04	I								1.36E+09		1	0.1	-Perfluorohexanoate	92612-52-7					3.9E+00	1.6E+01		3.2E+00
		5.00E-04	I								1.36E+09		1	0.1	-Perfluorohexanoic acid (PFHxA)	307-24-4					3.9E+00	1.6E+01		3.2E+00
		3.00E-06	A								1.36E+09		1	0.1	-Perfluorononanoate	72007-68-2					2.3E-02	9.9E-02		1.9E-02
		3.00E-06	A								1.36E+09		1	0.1	-Perfluorononanoic acid (PFNA)	375-95-1					2.3E-02	9.9E-02		1.9E-02
3.95E+01	D	4.00E-02	N								1.36E+09		1	0.1	-Perfluorooctadecanoic acid (PFODA)	16517-11-6					3.1E+02	3.3E+03		2.5E+02
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-04	1.3E+03		6.3E-04
2.93E+04	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-04	3.3E+03		6.3E-04
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-04	9.9E-04		1.9E-04
		3.00E-08	D								1.36E+09		1	0.1	-Perfluorooctanoic acid (PFOA)	335-87-1	2.4E-05	8.4E-05		1.9E-05	2.3E-04	9.9E-04		1.9E-04
		5.00E-04	R					V		1.36E+04	1.36E+09	2.65E+04	1		-Perfluoropropanoic acid (PFPrA)	422-64-0					3.9E+00			3.9E+00
		1.00E-03	N								1.36E+09		1	0.1	-Perfluorotetradecanoic acid (PFTeTDA)	376-06-7					7.8E+00	3.3E+01		6.3E+00
		3.00E-04	N								1.36E+09		1	0.1	-Perfluoroundecanoic acid (PFUDA)	2058-94-8					2.3E+00	9.9E+00		1.9E+00
		3.00E-04	P								1.36E+09		1	0.1	-Potassium perfluorobutanesulfonate	29420-49-3					2.3E+00	9.9E+00		1.9E+00
3.95E+01	D	2.00E-03	I					V		9.61E+04	1.36E+09	6.13E+04	1		-Potassium perfluorobutanoate	2966-54-3				1.4E-02			1.6E+01	
		2.15E-09	I								1.36E+09		1	0.1	-Potassium perfluorodecanoate	51604-85-4					1.7E-05	7.1E-05		1.4E-05
		1.00E-07	D								1.36E+09		1	0.1	-Potassium perfluorooctanesulfonate	2795-39-3					7.8E-04	3.3E+03		6.3E-04
		1.00E-03	I					V		8.99E+04	1.36E+09	6.02E+04	1		-Sodium perfluorobutanoate	2218-54-4					7.8E+00			7.8E+00
		2.09E-09	I								1.36E+09		1	0.1	-Sodium perfluorodecanoate	3830-45-3					1.6E-05	6.9E-05		1.3E-05
		5.00E-04	I								1.36E+09		1	0.1	-Sodium perfluorohexanoate	2923-26-4					3.9E+00	1.6E+01		3.2E+00
		7.00E-04	I								1.36E+09		1		Perchlorates						5.5E+00			5.5E+00
		7.00E-04	I								1.36E+09		1		-Ammonium Perchlorate	7790-98-9					5.5E+00			5.5E+00
		7.00E-04	I								1.36E+09		1		-Lithium Perchlorate	7791-03-9					5.5E+00			5.5E+00
		7.00E-04	I								1.36E+09		1		-Perchlorate and Perchlorate Salts	14797-73-0					5.5E+00			5.5E+00
		7.00E-04	I								1.36E+09		1		-Potassium Perchlorate	7778-74-7					5.5E+00			5.5E+00
		7.00E-04	I								1.36E+09		1		-Sodium Perchlorate	7601-89-0					5.5E+00			5.5E+00
2.20E-03	C	6.30E-07	C								1.36E+09		1	0.1	Permethrin	52645-53-1	3.2E+02	1.1E+03	6.1E+06	2.5E+02	3.9E+02	1.6E+03		3.2E+02
		2.40E-01	O								1.36E+09		1	0.1	Phenacetin	62-44-2					1.9E+03	7.9E+03		1.5E+03
		3.00E-01	I	2.00E-01	C						1.36E+09		1	0.1	Phenmedipham	13684-63-4					2.3E+03	9.9E+03	2.8E+07	1.9E+03
		4.00E-03	I								1.36E+09		1	0.1	Phenol	108-95-2					3.1E+01	1.3E+02		2.5E+01
		5.00E-04	X								1.36E+09		1	0.1	Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1					3.9E+00	1.6E+01		3.2E+00
		2.00E-04	X					V		1.29E+02	1.36E+09	7.06E+03	1		Phenothiazine	92-84-2					1.6E+00			1.6E+00
1.20E-01	P	6.00E-03	I								1.36E+09		1	0.1	Phenyl isothiocyanate	103-72-0					4.7E+01	2.0E+02		3.8E+01
		4.00E-03	P					M			1.36E+09		1	0.1	Phenylethylenediamine, m-	103-45-2	1.3E+00	5.0E+00		1.0E+00	3.1E+01	1.3E+02		2.5E+01
1.94E-03	H	1.00E-03	X								1.36E+09		1	0.1	Phenylethylenediamine, o-	95-54-5					3.1E+01	1.3E+02		2.5E+01
		2.00E-04	H								1.36E+09		1	0.1	Phenylphenol, p-	106-50-3					7.8E+00	3.3E+01		6.3E+00
		3.00E-04	I	V						1.61E+03	1.36E+09	9.81E+02	1		Phenylphenol, 2-	90-43-7	3.6E+02	1.3E+03		2.8E+02	1.6E+00	6.6E+00		1.3E+00
		2.00E-02	I								1.36E+09		1	0.1	Phorate	298-02-2					1.6E+00	6.6E+02		3.1E-02
		2.93E+00	X								1.36E+09		1		Phosgene	75-44-5					1.6E+02	6.6E+02		1.3E+02
		3.00E-01	X								1.36E+09		1		Phosmet	732-11-6					1.6E+02	6.6E+02		1.3E+02
		1.00E+00	P								1.36E+09		1		Phosphates, inorganic						2.3E+04			2.3E+04
		1.00E+00	P								1.36E+09		1		-Aluminum metaphosphate	13776-88-0					2.3E+03			2.3E+03
		1.00E+00	P								1.36E+09		1		-Aluminum salts of inorganic phosphates	E524680405					7.8E+03			7.8E+03
		1.00E+00	P								1.36E+09		1		-Dipotassium phosphate	7758-11-4					7.8E+03			7.8E+03
		1.00E+00	P								1.36E+09		1		-Disodium phosphate	7558-79-4					7.8E+03			7.8E+03
		3.54E+00	X								1.36E+09		1	0.1	-Monosodium phosphate	13530-50-2					2.8E+04			2.8E+04
		1.00E+00	P								1.36E+09		1		-Monopotassium phosphate	7778-77-0					7.8E+03			7.8E+03
		1.00E+00	P								1.36E+09		1		-Monosodium phosphate	7558-80-7					7.8E+03			7.8E+03
		1.00E+00	P	1.00E-02	I						1.36E+09		1		-Phosphoric Acid	7664-38-2					7.8E+03		1.4E+06	7.8E+03
		1.36E+00	X								1.36E+09		1	0.1	-Phosphoric acid, aluminum salt (1:1) [aluminum phosphate]	7784-30-7					1.1E+04	4.5E+04		8.6E+03
		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+04			3.3E+04
		1.00E+00	P								1.36E+09		1		-Polyphosphoric acid	8017-16-1					7.8E+03			7.8E+03
		1.00E+00	P								1.36E+09		1		-Potassium salts of inorganic phosphates	E524680403					7.8E+03			7.8E+03
		1.00E+00	P								1.36E+09		1		-Potassium tripolyphosphate	13845-36-8					7.8E+03			7.8E+03
		4.99E+00	X								1.36E+09		1		-Sodium aluminum phosphate (anhydrous)	10279-59-1					3.9E+04			3.9E+04
		3.52E+00	X								1.36E+09		1		-Sodium aluminum phosphate (tetrahydrate)	10305-76-7					2.8E+04			2.8E+04
		1.00E+00	P								1.36E+09		1		-Sodium hexametaphosphate	10124-56-8					7.8E+03			7.8E+03
		1.00E+00	P								1.36E+09		1		-Sodium polyphosphate	68915-31-1					7.8E+03			7.8E+03
		1.00E+00	P								1.36E+09		1		-Sodium pyrophosphate	7758-16-9					7.8E+03			7.8E+03
		1.00E+00	P																					



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) = 0.1							
SFO (mg/kg-day) <sup>1</sup>	key	IUR (ug/m <sup>3</sup> ) <sup>1</sup>	key	RD <sub>50</sub> (mg/kg-day)	key	RIC <sub>1</sub> (mg/m <sup>3</sup> )	key	vo	mutagen	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	GIABS	ABS <sub>0</sub>	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=0.1 (mg/kg)	Dermal SL THQ=0.1 (mg/kg)	Inhalation SL THQ=0.1 (mg/kg)	Noncarcinogenic SL THQ=0.1 (mg/kg)	
1.20E-01	H			5.00E-03	I	3.00E-03	C				1.36E+09		1	Silica (crystalline, respirable)	7631-86-9									4.3E+05	4.3E+05
				5.00E-03	I						1.36E+09		0.04	Silver	7440-22-4						3.9E+01	1.6E+02	4.3E+02		3.9E+01
				1.30E-02	I						1.36E+09		0.1	Simazine	122-34-9	5.8E+00	2.1E+01		4.5E+00		3.9E+01	1.6E+02	4.3E+02		3.2E+01
2.70E-01	H			4.00E-03	I						1.36E+09		1	Sodium Acifluorfen	62476-59-9						1.0E+02	4.3E+02			8.2E+01
				3.00E-02	I						1.36E+09		0.1	Sodium Azide	26628-22-8	2.6E+00	9.2E+00		2.0E+00		3.1E+01	9.9E+02			3.1E+01
				5.00E-02	A	1.40E-02	C				1.36E+09		1	Sodium Diethyldithiocarbamate	148-18-5						2.3E+02	9.9E+02			1.9E+02
				2.00E-05	I						1.36E+09		1	Sodium Fluoracetate	7681-49-4						3.9E+02	6.6E-01	2.0E+06		3.9E+02
				1.00E-03	H						1.36E+09		1	Sodium Melavanadate	62-74-8						1.6E-01				1.3E-01
				8.00E-04	P						1.36E+09		1	Sodium Tungstate	13472-45-2						6.3E+00				6.3E+00
2.40E-02	H			8.00E-04	P						1.36E+09		1	Sodium Tungstate Dihydrate	10213-10-2						6.3E+00				6.3E+00
				3.00E-02	I						1.36E+09		0.1	Stirofos (Tetrachlorovinphos)	961-11-5	2.9E+01	1.0E+02		2.3E+01		2.3E+02	9.9E+02			1.9E+02
				6.00E-01	I						1.36E+09		1	Strontium, Stable	7440-24-6						4.7E+03				4.7E+03
				3.00E-04	I						1.36E+09		0.1	Strychnine	57-24-9						2.3E+00	9.9E+00			1.9E+00
				2.00E-01	I	1.00E+00	V			8.67E+02	1.36E+09	9.35E+03	1	Styrene	100-42-5						1.6E+03		9.7E+02		6.0E+02
				3.00E-03	P						1.36E+09		0.1	Styrene-Acrylonitrile (SAN) Trimer (THNA isomer)	57964-39-3						2.3E+01	9.9E+01			1.9E+01
				3.00E-03	P						1.36E+09		0.1	Styrene-Acrylonitrile (SAN) Trimer (THNP isomer)	57964-40-6						2.3E+01	9.9E+01			1.9E+01
				1.00E-03	P	2.00E-03	X				1.36E+09		0.1	Sulfolane	126-33-0						7.8E+00	3.3E+01	2.8E+05		6.3E+00
				8.00E-04	P						1.36E+09		1	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9						6.3E+00	2.6E+01			5.1E+00
						1.00E-03	C	V			1.36E+09		1	Sulfur Trioxide	7446-11-9										1.4E+05
						1.00E-03	C				1.36E+09		1	Sulfuric Acid	7664-93-9										1.4E+05
2.50E-02	I	7.10E-06	I	5.00E-02	H						1.36E+09		0.1	Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethyl)phenoxy]-1-methylethyl ester	140-57-8	2.8E+01	9.9E+01	5.4E+05	2.2E+01	3.9E+02	1.6E+03				3.2E+02
				7.00E-02	I						1.36E+09		0.1	Tebuthiuron	34014-18-1						5.5E+02	2.3E+03			4.4E+02
				2.00E-02	H						1.36E+09		0.1	Temephos	3383-96-8						1.6E+02	6.6E+02			1.3E+02
				1.30E-02	I						1.36E+09		0.1	Terbacil	5902-51-2						1.0E+02	4.3E+02			8.2E+01
				2.50E-05	H				V	3.09E+01	1.36E+09	2.64E+05	1	Terbufos	13071-79-9						2.0E-01				2.0E-01
				1.00E-03	I						1.36E+09		0.1	Terbutyn	886-50-0						7.8E+00	3.3E+01			6.3E+00
5.00E-03	C	1.30E-06	C	1.00E-04	I						1.36E+09	3.99E+03	1	Tert-Butyl Acetate	540-88-5	1.4E+02		8.6E+00	8.1E+00		7.8E-01	3.3E+00			6.3E-01
				3.00E-05	P						1.36E+09		0.1	Tetrabromodiphenyl ether, 2,2',4,4'-(BDE-47)	5436-43-1						2.3E-01				2.3E-01
											1.36E+09	5.07E+04	1	Tetrachlorobenzene, 1,2,4,5-	95-94-3						7.8E-01	3.3E+00			6.3E-01
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	2.0E+00	2.3E+02				2.3E+02
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	6.0E-01	1.6E+02				1.6E+02
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	3.3E+02	2.5E+01	2.4E+01	4.7E+01	9.9E+00		9.8E+00	8.1E+00	
				3.00E-02	I						1.36E+09		0.1	Tetrachlorophenol, 2,3,4,6-	58-90-2						2.3E+02	9.9E+02			1.9E+02
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			4.3E-02	4.7E-01					4.7E-01
				5.00E-04	I						1.36E+09		0.1	Tetraethyl Dithiopyrophosphate	3689-24-5						3.9E+00	1.6E+01			3.2E+00
						8.00E+01	V			2.05E+03	1.36E+09	1.22E+03	1	Tetrafluoroethane, 1,1,1,2-	811-97-2						7.8E-01	3.3E+00		1.0E+04	1.0E+04
				1.00E-04	X						1.36E+09		0.1	Tetramethylphosphoramide, -N,N,N',N'-(TMPA)	16853-36-4						1.6E-01	1.0E+04			6.3E-01
				2.00E-03	P						1.36E+09		0.00065	Tetryl (Trinitrophenylmethylnitramine)	479-45-8						1.6E+01	1.0E+04			1.6E+01
				2.00E-05	G						1.36E+09		1	Thallic Oxide	1314-32-5						1.6E-01				1.6E-01
				1.00E-05	X						1.36E+09		1	Thallium (I) Nitrate	10102-45-1						7.8E-02				7.8E-02
				1.00E-05	X						1.36E+09		1	Thallium (Soluble Salts)	7440-28-0						7.8E-02				7.8E-02
				1.00E-05	X						1.36E+09	1.40E+05	1	Thallium Acetate	563-68-8						7.8E-02				7.8E-02
				2.00E-05	X						1.36E+09		0.1	Thallium Carbonate	6533-73-9						1.6E-01	6.6E-01			1.3E-01
				1.00E-05	X						1.36E+09		1	Thallium Chloride	7731-12-0						7.8E-02				7.8E-02
				1.00E-05	G						1.36E+09		1	Thallium Selenite	12039-52-0						4.7E-02				7.8E-02
				2.00E-05	X						1.36E+09		1	Thallium Sulfate	7446-18-6						1.6E-01				1.6E-01
				4.30E-02	O						1.36E+09		0.1	Thifensulfuron-methyl	79277-27-3						3.4E+02	1.4E+03			2.7E+02
				1.00E-02	I						1.36E+09		0.1	Thiobencarb	28249-77-6						7.8E+01	3.3E+02			6.3E+01
				2.00E-04	P						1.36E+09		1	Thiocyanates	E1790665						1.6E+00				1.6E+00
				2.00E-04	X						1.36E+09		1	Thiocyanic Acid	463-56-9						1.6E+00				1.6E+00
				3.00E-02	H						1.36E+09		0.1	Thiocyanic acid, (2-benzothiazolythio)methyl ester (TCMTB)	21564-17-0						2.3E+02	9.9E+02			1.9E+02
				7.00E-02	X						1.36E+09		0.0075	Thiodiethylol	111-48-8						5.5E+02	3.1E+04			5.4E+02
				3.00E-04	H						1.36E+09		0.1	Thiofanox	39196-18-4						2.3E+00	9.9E+00			1.9E+00
1.16E-02	O			1.60E-01	O						1.36E+09		0.1	Thiophanate, Methyl	23564-05-8	6.0E+01	2.1E+02		4.7E+01	1.3E+03	5.3E+03				1.0E+03
				1.50E-02	O						1.36E+09		0.1	Thiram	137-26-8						1.2E+02	4.9E+02			9.5E+01
				6.00E-01	H																				

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) = 0.1						
SFO (mg/kg-day) <sup>1</sup>	k <sub>e</sub> y	IUR (ug/m <sup>3</sup> ) <sup>1</sup>	k <sub>e</sub> y	RD <sub>50</sub> (mg/kg-day)	k <sub>e</sub> y	RIC <sub>1</sub> (mg/m <sup>3</sup> ) <sup>1</sup>	k <sub>e</sub> y	v <sub>o</sub> l	mutagen	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	GIABS	ABS <sub>0</sub>	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=0.1 (mg/kg)	Dermal SL THQ=0.1 (mg/kg)	Inhalation SL THQ=0.1 (mg/kg)	Noncarcinogenic SL Child THQ=0.1 (mg/kg)	
9.00E-03	P			2.00E-04	O						1.36E+09		1	0.1	Tribufos	78-48-8					1.6E+00	6.6E+00		1.3E+00	
				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+01	3.3E+02		6.3E+01	
				3.00E-04	P						1.36E+09		1	0.1	Tributyltin Compounds	E1790679					2.3E+00	9.9E+00		1.9E+00	
				3.00E-04	I						1.36E+09		1	0.1	Tributyltin Oxide	56-35-9					2.3E+00	9.9E+00		1.9E+00	
				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+05	6.6E+02	6.7E+02	6.7E+02	
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+02	6.6E+02		1.3E+02	
2.90E-02	H										1.36E+09		1	0.1	Trichloroacetic Acid	76-03-9	9.9E+00	3.5E+01		7.8E+00					
7.00E-03	X			3.00E-05	X						1.36E+09		1	0.1	Trichloroaniline, 2,4,6-	33663-50-2	2.4E+01	8.5E+01		1.9E+01					
				8.00E-04	X			V			1.36E+09	3.22E+04	1		Trichloroaniline, 2,4,6-	634-93-5	9.9E+01	3.5E+02		7.8E+01					
											1.36E+09		1	0.1	Trichlorobenzene, 1,2,3-	87-61-6					6.3E+00	9.9E-01		1.9E-01	
2.90E-02	P			1.00E-02	I	2.00E-03	P	V		4.04E+02	1.36E+09	2.99E+04	1		Trichlorobenzene, 1,2,4-	120-82-1	2.4E+01			2.4E+01					
				2.00E+00	I	5.00E+00	I	V		6.40E+02	1.36E+09	1.65E+03	1		Trichloroethane, 1,1,1-	71-55-6					1.6E+04	8.6E+02		8.1E+02	
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		Trichloroethane, 1,1,2-	79-00-5	1.2E+01		1.3E+00	1.1E+00	3.1E+01	1.5E-01		1.5E-01	
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		Trichloroethylene	79-01-6	8.8E+00		1.1E+00	9.4E-01	3.9E+00		4.6E-01	4.1E-01	
				3.00E-01	I			V		1.23E+03	1.36E+09	1.04E+03	1		Trichlorofluoromethane	75-69-4					2.3E+03			2.3E+03	
				1.00E-01	I						1.36E+09		1	0.1	Trichlorophenol, 2,4,5-	95-95-4					7.8E+02	3.3E+03		6.3E+02	
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+00	3.3E+01		5.8E+00	
				1.00E-02	I						1.36E+09		1	0.1	Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5					7.8E+01	3.3E+02		6.3E+01	
				8.00E-03	I						1.36E+09		1	0.1	Trichloroxypropionic acid, -2,4,5	93-72-1					6.3E+01	2.6E+02		5.1E+01	
3.00E+01	I			5.00E-03	I		V			1.28E+03	1.36E+09	1.50E+04	1		Trichloropropane, 1,1,2-	598-77-6					3.9E+01			3.9E+01	
				4.00E-03	I	3.00E-04	I	V	M	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+01		4.9E-01		4.8E-01
				3.00E-03	X	3.00E-04	P	V		3.11E+02	1.36E+09	2.34E+03	1		Trichloropropane, 1,2,3-	96-19-5					2.3E+01		7.3E-02		7.3E-02
				2.00E-02	A						1.36E+09		1	0.1	Tricresyl Phosphate (TCP)	1330-78-5					1.6E+02	6.6E+02		1.3E+02	
				3.00E-03	I						1.36E+09		1	0.1	Tridiphane	58138-08-2					2.3E+01	9.9E+01		1.9E+01	
						7.00E-03	I	V		2.79E+04	1.36E+09	1.58E+04	1		Triethylamine	121-44-8							1.2E+01		1.2E+01
				2.00E+00	P						1.36E+09		1	0.1	Triethylene Glycol	112-27-6					1.6E+04	6.6E+04		1.3E+04	
7.70E-03	I			7.50E-03	I		P	V		4.81E+03	1.36E+09	7.12E+02	1		Trifluoroethane, 1,1,1-	420-46-2							1.5E+03	1.5E+03	
2.00E-02	P			1.00E-02	P						1.36E+09		1	0.1	Trifuralin	1582-09-8	9.0E+01			9.0E+01	5.9E+01				5.9E+01
				1.00E-02	P						1.36E+09		1	0.1	Trimethyl Phosphate	512-56-1	3.5E+01	1.2E+02		2.7E+01	7.8E+01	3.3E+02			6.3E+01
				1.00E-02	I	6.00E-02	I	V		2.93E+02	1.36E+09	9.44E+03	1		Trimethylbenzene, 1,2,3-	526-73-8					7.8E+01		5.9E+01		3.4E+01
				1.00E-02	I	6.00E-02	I	V		2.19E+02	1.36E+09	7.91E+03	1		Trimethylbenzene, 1,2,4-	95-63-6					7.8E+01		5.0E+01		3.0E+01
				1.00E-02	I	6.00E-02	I	V		1.82E+02	1.36E+09	6.61E+03	1		Trimethylbenzene, 1,3,5-	108-67-8					7.8E+01		4.1E+01		2.7E+01
				1.00E-02	X			V		2.96E+01	1.36E+09	1.00E+03	1		Trimethylpentene, 2,4,4-	25167-70-8					7.8E+01				7.8E+01
				3.00E-02	I						1.36E+09		1	0.019	Trinitrobenzene, 1,3,5-	99-35-4					2.3E+02	5.2E+03			2.2E+02
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+00	5.2E+01			3.6E+00
				2.00E-02	P						1.36E+09		1	0.1	Triphenylphosphine Oxide	791-28-6					1.6E+02	6.6E+02			1.3E+02
				2.00E-02	A						1.36E+09		1	0.1	Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8					1.6E+02	6.6E+02			1.3E+02
				1.00E-02	X						1.36E+09		1	0.1	Tris(1-chloro-2-propyl)phosphate	13674-84-5					7.8E+01	3.3E+02			6.3E+01
2.30E+00	C	6.60E-04	C					V		4.67E+02	1.36E+09	9.03E+05	1		Tris(2,3-dibromopropyl)phosphate	126-72-7	3.0E-01		3.8E+00	2.8E-01					
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+01	2.3E+02			4.4E+01
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+02	3.3E+03			6.3E+02
				8.00E-04	P						1.36E+09		1		Tunasten	7440-33-7					6.3E+00				6.3E+00
				2.00E-04	A	4.00E-05	A				1.36E+09		1		Uranium	7440-61-1					1.6E+00		5.7E+03		1.6E+00
1.00E+00	C	2.90E-04	C						M		1.36E+09		1	0.1	Urethane	51-79-6	1.5E-01	6.0E-01	4.8E+03	1.2E-01					
		8.30E-03	P	9.00E-03	I	7.00E-06	P				1.36E+09		0.026		Vanadium Pentoxide	1314-62-1			4.6E+02	4.6E+02	7.0E+01		9.9E+02		6.6E+01
				5.04E-03	G	1.00E-04	A				1.36E+09		0.026		Vanadium and Compounds	7440-62-2					3.9E+01		1.4E+04		3.9E+01
				1.00E-03	I			V			1.36E+09	1.23E+05	1		Vernolate	1929-77-7					7.8E+00				7.8E+00
				1.20E-03	O						1.36E+09		1	0.1	Vinclozolin	50471-44-8					9.4E+00	4.0E+01			7.6E+00
				1.00E+00	H	2.00E-01	I	V		2.75E+03	1.36E+09	4.40E+03	1		Vinyl Acetate	108-05-4					7.8E+03		9.2E+01		9.1E+01
7.20E-01	I	1.50E-05	P	3.00E-03	I	5.11E-02	A	V	M	3.92E+03	1.36E+09	9.56E+02	1		Vinyl Bromide	593-60-2			2.6E-01	2.6E-01					4.3E-01
				3.00E-04	I						1.36E+09		1	0.1	Vinyl Chloride	75-01-4	9.4E-02		1.6E-01	5.9E-02	2.3E+01		5.1E+00		4.2E+00
											1.36E+09		1		Warfarin	81-81-2					2.3E+00	9.9E+00			1.9E+00
				2.00E-01	G	1.00E-01	G	V		3.88E+02	1.36E+09	5.47E+03	1		Xylene, m-	108-38-3					1.6E+03		5.7E+01		5.5E+01