

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (see FO #31); H = HEAST; F = See FAQ; W = see user guide Section 2.3.5; E = see user guide Section 2.3.6; L = see user's guide Section 5.2; M = mutagen; S = see user's guide Section 5; V = volatile; R = RBA applied (see user's guide Section 5.10); c = cancer; n = noncancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; SSL values are based on DAF=1; n = Concentration may exceed ceiling limit (see user's guide Section 5.13); s = concentration may exceed Csat (see user's guide Section 5.12)															Toxicity and Chemical-specific Information		Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k _e (y ⁻¹)	IUR (ug/m ³ -y ⁻¹)	k _e (y ⁻¹)	RFD _o (mg/kg-day)	k _e (y ⁻¹)	RfC _o (mg/m ³)	k _e (y ⁻¹)	muta gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	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 TH=1 (mg/kg)			
2.2E-06	I			1.2E-03	O	9.0E-03	I	V	1.07E+05	1.36E+09	8.72E+03	1	0.1	Acetophenone	30560-19-1				4.9E+01	4.9E+01	1.4E+03	3.3E+03	3.4E+02	9.8E+02		
				2.0E-02	I				1.36E+09			1	0.1	Acetaldehyde	75-07-0								3.4E+02	1.6E+04		
				9.0E-01	I	3.1E+01	A	V	1.14E+05	1.36E+09	1.37E+04	1		Acetone	67-64-1					2.3E+04	5.5E+04		1.8E+06	6.7E+05		
						2.0E-03	X		1.36E+09			1	0.1	Acetone Cyanohydrin	75-86-5							1.2E+07	1.2E+07			
				6.0E-02	I			V	1.28E+05	1.36E+09	1.30E+04	1		Acetonitrile	75-05-8							3.4E+03	3.4E+03			
3.8E+00	C	1.3E-03	C	1.0E-01	I			V	2.52E+03	1.36E+09	5.97E+04	1		Acetophenone	98-86-2	8.6E-01	2.0E+00	1.3E+04	6.0E-01	1.2E+05			1.2E+05	1.2E+05		
				5.0E-04	I	2.0E-05	I	V	2.27E+04	1.36E+09	6.91E+03	1		Acetylaminofluorene, 2-	53-96-3					5.8E+02		6.1E-01	6.0E-01			
5.0E-01	I	1.0E-04	I	2.0E-03	I	6.0E-03	I	M	1.36E+09			1	0.1	Acrolein	107-02-8	6.5E+00	1.5E+01	1.7E+05	4.6E+00	2.3E+03	5.5E+03	3.6E+07	1.6E+03			
				5.0E-01	I	1.0E-03	I	V	1.09E+05	1.36E+09	9.53E+04	1		Acrylic Acid	79-10-7					5.8E+05		4.2E+02	4.2E+02			
5.4E-01	I	6.8E-05	I	4.0E-02	A	2.0E-03	I	V	1.13E+04	1.36E+09	7.69E+03	1		Acrylonitrile	107-13-1	6.1E+00		1.4E+00	1.1E+00	4.7E+04		6.7E+01	6.7E+01			
						6.0E-03	P		1.36E+09			1	0.1	Adiponitrile	111-69-3							3.6E+07	3.6E+07			
5.6E-02	C			1.0E-02	I				1.36E+09			1	0.1	Alachlor	15972-60-8	5.8E+01	1.4E+02		4.1E+01	1.2E+04	2.8E+04		8.2E+03	8.2E+03		
				1.0E-03	I				1.36E+09			1	0.1	Aldicarb	116-06-3					1.2E+03	2.8E+03		8.2E+02	8.2E+02		
				1.0E-03	I				1.36E+09			1	0.1	Aldicarb Sulfone	1646-88-4					1.2E+03	2.8E+03		8.2E+02	8.2E+02		
1.7E+01	I	4.9E-03	I	3.0E-05	I			V	1.36E+09	1.72E+06		1		Aldrin	309-00-2	1.9E-01		4.3E+00	1.8E-01	3.5E+01			3.5E+01	3.5E+01		
				5.0E-03	I	1.0E-04	X	V	1.11E+05	1.36E+09	3.42E+04	1		Allyl Alcohol	107-18-6					5.8E+03		1.5E+01	1.5E+01			
2.1E-02	C	6.0E-06	C	1.0E+00	P	5.0E-03	P		1.42E+03	1.36E+09	1.58E+03	1		Allyl Chloride	107-05-1	1.6E+02		3.2E+00	3.2E+00			6.9E+00	6.9E+00			
				4.0E-04	I				1.36E+09			1		Aluminum	7429-90-5					1.2E+06		3.0E+07	1.1E+06	1.1E+06		
				9.0E-03	I				1.36E+09			1	0.1	Aluminum Phosphide	20859-73-8					4.7E+02			4.7E+02	4.7E+02		
									1.36E+09			1	0.1	Ametryn	834-12-8					1.1E+04	2.5E+04		7.4E+03	7.4E+03		
				8.0E-02	P				1.36E+09			1	0.1	Aminobiphenyl, 4-	92-67-1	1.6E-01	3.7E-01	2.8E+03	1.1E-01					7.4E+03		
				4.0E-03	X				1.36E+09			1	0.1	Aminophenol, m-	591-27-5					9.3E+04	2.2E+05		6.6E+04	6.6E+04		
				2.0E-02	P				1.36E+09			1	0.1	Aminophenol, o-	95-55-6					4.7E+03	1.1E+04		3.3E+03	3.3E+03		
				2.5E-03	I				1.36E+09			1	0.1	Aminophenol, p-	123-30-8					2.3E+04	5.5E+04		1.6E+04	1.6E+04		
						5.0E-01	I	V	1.36E+09			1		Amirtraz	33089-81-1					2.9E+03	6.9E+03		2.1E+03	2.1E+03		
				2.0E-01	I				1.36E+09			1		Ammonia	7664-41-7											
						3.0E-03	X	V	1.37E+04	1.36E+09	2.62E+04	1		Ammonium Sulfamate	7773-06-0					2.3E+05			2.3E+05	2.3E+05		
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I		1.36E+09			1	0.1	Amyl Alcohol, tert-	75-85-4					8.2E+03	1.9E+04	3.4E+02	3.4E+02	3.4E+02		
4.0E-02	P			2.0E-03	X				1.36E+09			1	0.1	Aniline	62-53-3	5.7E+02	1.4E+03	1.0E+07	4.0E+02	2.3E+03	5.5E+03	6.0E+06	5.7E+03	5.7E+03		
				4.0E-04	I				1.36E+09		0.15			Anthraquinone, 9,10-	84-65-1	8.2E+01	1.9E+02		5.7E+01				1.6E+03	1.6E+03		
				5.0E-04	H				1.36E+09		0.15			Antimony (metallic)	7440-36-0					4.7E+02			4.7E+02	4.7E+02		
				4.0E-04	H				1.36E+09		0.15			Antimony Pentoxide	1314-60-9					5.8E+02			5.8E+02	5.8E+02		
				4.0E-04	H				1.36E+09		0.15			Antimony Tetroxide	1332-81-6					4.7E+02			4.7E+02	4.7E+02		
1.5E+00	I	4.3E-03	I	2.0E-04	I	1.5E-05	C		1.36E+09			0.03		Antimony Trioxide	1309-64-4							1.2E+06	1.2E+06	1.2E+06		
				3.5E-06	C	5.0E-05	I		1.36E+09			1		Arsenic, Inorganic	7440-38-2	3.6E+00	1.7E+01	3.9E+03	3.0E+00	5.8E+02	2.8E+03	8.9E+04	4.8E+02	4.8E+02		
				3.6E-02	O				1.36E+09			1	0.1	Arsine	7784-42-1					4.1E+00		3.0E+05	4.1E+00	4.1E+00		
2.3E-01	C			3.5E-02	I				1.36E+09			1	0.1	Asulam	3337-71-1	1.4E+01	3.4E+01		1.0E+01	4.2E+04	9.9E+04		3.0E+04	3.0E+04		
8.8E-01	C	2.5E-04	C	1.0E-01	I				1.36E+09			1	0.1	Atrazine	1912-24-9	3.7E+00	8.8E+00	6.7E+04	2.6E+00	4.1E+04	9.7E+04		2.9E+04	2.9E+04		
				4.0E-04	I				1.36E+09			1	0.1	Auramine	492-80-8											
				3.0E-03	A	1.0E-02	A		1.36E+09			1	0.1	Avermectin B1	65195-53-3					4.7E+02	1.1E+03		3.3E+02	3.3E+02		
1.1E-01	I	3.1E-05	I	2.0E-01	I	5.0E-04	H		1.36E+09			0.07		Azinphos-methyl	86-50-0	3.0E+01		2.1E+02	2.6E+01	3.5E+03	8.3E+03	6.0E+07	2.5E+03	2.5E+03		
				1.0E+00	P	7.0E-06	P		1.36E+09			1	0.1	Azobenzene	103-33-3											
				2.0E-01	I	5.0E-04	H		1.36E+09			1		Azodicarbonamide	123-77-3					1.2E+06	2.8E+06	4.2E+04	4.0E+04	4.0E+04		
				5.0E-03	O			V	1.36E+09			1		Barium	7440-39-3					2.3E+05		3.0E+06	2.2E+05	2.2E+05		
									1.36E+09			1	0.1	Benfluralin	1861-40-1					5.8E+03			5.8E+03	5.8E+03		
				5.0E-02	I				1.36E+09			1	0.1	Benmethyl	17804-35-2					5.8E+04	1.4E+05		4.1E+04	4.1E+04		
				2.0E-01	I				1.36E+09			1	0.1	Bensulfuron-methyl	83055-99-6					2.3E+05	5.5E+05		1.6E+05	1.6E+05		
				3.0E-02	I				1.36E+09			1	0.1	Bentazon	25057-89-0					3.5E+04	8.3E+04		2.5E+04	2.5E+04		
4.0E-03	P			1.0E-01	I			V	1.16E+03	1.36E+09	2.25E+04	1		Benzaldehyde	100-52-7	8.2E+02			8.2E+02	1.2E+05			1.2E+05	1.2E+05		
5.5E-02	I	7.8E-06	I	4.0E-03	I	3.0E-02	I	V	1.82E+03	1.36E+09	3.54E+03	1		Benzene	71-43-2	5.9E+01		5.6E+00	5.1E+00	4.7E+03		4.6E+02	4.2E+02	4.2E+02		
1.0E-01	X			3.0E-04	X				1.36E+09			1	0.1	Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1	3.3E+01	7.7E+01		2.3E+01	3.5E+02	8.3E+02		2.5E+02	2.5E+02		
2.3E+02	I	6.7E-02	I	1.0E-03	P			V	1.26E+03	1.36E+09	1.94E+04	1		Benzenethiol	108-98-5					1.2E+03			1.2E+03	1.2E+03		
				3.0E-03	I			M	1.36E+09			1	0.1	Benzidine	92-87-5	1.4E-02	3.4E-02	2.5E+02	1.0E-02	3.5E+03	8.3E+03		2.5E+03	2.5E+03		
				4.0E+00	I				1.36E+09																	

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Toxicity and Chemical-specific Information													Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -y)	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³)	k _e (y)	muta gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	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 THI=1 (mg/kg)
				3.0E-04	X		V		8.96E+02	1.36E+09	1.12E+04	1		Bromo-3-fluorobenzene, 1-	1073-06-9					3.5E+02			3.5E+02
				3.0E-04	X		V		3.23E+02	1.36E+09	1.14E+04	1		Bromo-4-fluorobenzene, 1-	460-00-4					3.5E+02			3.5E+02
				8.0E-03	I	6.0E-02	I	V	6.79E+02	1.36E+09	8.37E+03	1		Bromobenzene	108-86-1					9.3E+03		2.2E+03	1.8E+03
						4.0E-02	X	V	4.04E+03	1.36E+09	3.58E+03	1		Bromochloromethane	74-97-5							6.3E+02	6.3E+02
6.2E-02	I	3.7E-05	C	2.0E-02	I		V		9.32E+02	1.36E+09	3.97E+03	1		Bromodichloromethane	75-27-4	5.3E+01		1.3E+00	1.3E+00	2.3E+04			2.3E+04
7.9E-03	I	1.1E-06	I	2.0E-02	I		V		9.15E+02	1.36E+09	9.70E+03	1		Bromofom	75-25-2	4.1E+02		1.1E+02	8.6E+01	2.3E+04			2.3E+04
				1.4E-03	I	5.0E-03	I	V	3.59E+03	1.36E+09	1.40E+03	1		Bromomethane	74-83-9					1.6E+03		3.1E+01	3.0E+01
				5.0E-03	H		V		1.36E+09	1.24E+05		1		Bromophos	2104-96-3					5.8E+03			5.8E+03
1.0E-01	O			1.5E-02	O	1.0E-01	A	V	9.66E+02	1.36E+09	2.14E+03	1	0.1	Bromopropane, 1-	106-94-5				2.2E+01	1.8E+04	4.1E+04	9.4E+02	9.4E+02
1.0E-01	O			1.5E-02	O		V		1.36E+09	4.74E+05		1		Bromoxynil	1689-84-5	3.2E+01	7.5E+01			1.8E+04			1.2E+04
3.4E+00	C	3.0E-05	I			2.0E-03	I	V	6.67E+02	1.36E+09	8.66E+02	1		Bromoxynil Octanoate	1689-99-2	3.2E+01			3.2E+01	1.8E+04			1.8E+04
				3.0E-02	O		V		1.36E+09			1	0.1	Butadiene, 1,3-	106-99-0	9.6E-01		3.5E-01	2.6E-01			7.6E+00	7.6E+00
				1.0E-01	I		V		7.64E+03	1.36E+09	3.00E+04	1		Butanoic acid, 4-(2,4-dichlorophenoxy)-	94-82-6					3.5E+04	8.3E+04		2.5E+04
				2.0E+00	P	3.0E+01	P	V	2.13E+04	1.36E+09	2.92E+04	1		Butanol, N-	71-36-3					1.2E+05			1.2E+05
				5.0E-02	I		V		1.36E+09	8.63E+04		1		Butyl alcohol, sec-	78-92-2					2.3E+06		3.8E+06	1.5E+06
														Butylate	2008-41-5					5.8E+04			5.8E+04
2.0E-04	C	5.7E-08	C							1.36E+09		1	0.1	Butylated hydroxyanisole	25013-16-5	1.6E+04	3.9E+04	2.9E+08	1.1E+04	3.5E+05	8.3E+05		2.5E+05
3.6E-03	P			3.0E-01	P		V		1.08E+02	1.36E+09	8.14E+03	1		Butylated hydroxytoluene	128-37-0	9.1E+02	2.1E+03		6.4E+02	5.8E+04			5.8E+04
				5.0E-02	P		V		1.08E+02	1.36E+09	8.14E+03	1		Butylbenzene, n-	104-51-8					1.2E+05			1.2E+05
1.0E-01	X			1.0E-01	X		V		1.45E+02	1.36E+09	7.35E+03	1		Butylbenzene, sec-	135-98-8					1.2E+05			1.2E+05
1.0E-01	X			1.0E-01	X		V		1.83E+02	1.36E+09	7.36E+03	1		Butylbenzene, tert-	98-06-6					1.2E+05			1.2E+05
				2.0E-02	A		V		1.36E+09			1	0.1	Cacodylic Acid	75-60-5					2.3E+04	5.5E+04		1.6E+04
		1.8E-03	I	1.0E-03	I	1.0E-05	A			1.36E+09		0.025	0.001	Cadmium (Diet)	7440-43-9			9.3E+03	9.3E+03	1.2E+03	6.9E+03	6.0E+04	9.8E+02
		1.8E-03	I	5.0E-04	I	1.0E-05	A			1.36E+09		0.05	0.001	Cadmium (Water)	7440-43-9					1.2E+03	6.9E+03	6.0E+04	9.8E+02
				5.0E-01	I	2.2E-03	C			1.36E+09		1	0.1	Caprolactam	105-60-2					5.8E+05	1.4E+06	1.3E+07	4.0E+05
1.5E-01	C	4.3E-05	C	2.0E-03	I		V			1.36E+09		1	0.1	Captafol	2425-06-1	2.2E+01	5.2E+01	3.9E+05	1.5E+01	2.3E+03	5.5E+03		1.6E+03
2.3E-03	C	6.6E-07	C	1.3E-01	I		V			1.36E+09		1	0.1	Captan	133-06-2	1.4E+03	3.4E+03	2.5E+07	1.0E+03	1.5E+05	3.6E+05		1.1E+05
				1.0E-01	I		V			1.36E+09		1	0.1	Carbaryl	63-25-2					1.2E+05	2.8E+05		8.2E+04
				5.0E-03	I		V			1.36E+09		1	0.1	Carbaryl	1563-66-2					5.8E+03	1.4E+04		4.1E+03
				1.0E-01	I	7.0E-01	I	V	7.38E+02	1.36E+09	1.17E+03	1		Carbon Disulfide	75-15-0					1.2E+05			3.6E+03
7.0E-02	I	6.0E-06	I	4.0E-03	I	1.0E-01	I	V	4.58E+02	1.36E+09	1.49E+03	1		Carbon Tetrachloride	56-23-5	4.7E+01		3.1E+00	2.9E+00	4.7E+05		6.5E+02	5.7E+02
						1.0E-01	P	V	5.89E+03	1.36E+09	6.46E+02	1		Carbonyl Sulfide	463-58-1					1.2E+04	2.8E+04		2.8E+02
				1.0E-01	I		V			1.36E+09		1	0.1	Carbosulfan	55285-14-8					1.2E+04	2.8E+04		8.2E+03
				1.0E-01	I		V			1.36E+09		1	0.1	Carboxin	5234-68-4					1.2E+05	2.8E+05		8.2E+04
						9.0E-04	I			1.36E+09		1		Ceric oxide	1306-38-3					1.2E+05		5.4E+06	5.4E+06
				1.0E-01	I		V			1.36E+09	1.45E+05	1		Chloral Hydrate	302-17-0					1.2E+05			1.2E+05
				1.5E-02	I		V			1.36E+09		1	0.1	Chloramben	133-90-4					1.8E+04	4.1E+04		1.2E+04
4.0E-01	H						V			1.36E+09		1	0.1	Chloranil	118-75-2	8.1E+00	1.9E+01		5.7E+00	5.8E+02	3.4E+03	4.7E+03	4.5E+02
3.5E-01	I	1.0E-04	I	5.0E-04	I	7.0E-04	I	V		1.36E+09	1.53E+06	1	0.04	Chlorodane	12789-03-6	9.3E+00	5.5E+01	1.9E+02	7.7E+00	3.5E+02	8.3E+02		2.5E+02
1.0E+01	I	4.6E-03	C	3.0E-04	I		V			1.36E+09		1	0.1	Chlordecone (Kepone)	143-50-0	3.3E-01	7.7E-01	3.6E+03	2.3E-01	3.5E+02	2.8E+05		8.2E+02
				7.0E-04	A		V			1.36E+09		1	0.1	Chlorfenvinphos	470-90-6					8.2E+02	1.9E+03		5.7E+02
				9.0E-02	O		V			1.36E+09		1	0.1	Chlorimuron, Ethyl-	90982-32-4					1.1E+05	2.5E+05		7.4E+04
				1.0E-01	I	1.5E-04	A	V	2.78E+03	1.36E+09	1.22E+03	1		Chlorine	7782-50-5					1.2E+05		7.8E-01	7.8E-01
				3.0E-02	I	2.0E-04	I	V		1.36E+09		1		Chlorine Dioxide	10049-04-4					3.5E+04		1.2E+06	3.4E+04
				3.0E-02	I		V			1.36E+09		1		Chlorite (Sodium Salt)	7758-19-2					3.5E+04			3.5E+04
						5.0E+01	I	V	1.15E+03	1.36E+09	1.03E+03	1		Chloro-1,1-difluoroethane, 1-	75-68-3							2.3E+05	2.3E+05
		3.0E-04	I	2.0E-02	H	2.0E-02	I	V	7.86E+02	1.36E+09	1.08E+03	1		Chloro-1,3-butadiene, 2-	126-99-8			4.4E-02	4.4E-02	2.3E+04		9.4E+01	9.4E+01
4.6E-01	H						V			1.36E+09		1	0.1	Chloro-2-methylaniline HCl, 4-	3165-93-3	7.1E+00	1.7E+01		5.0E+00	3.5E+03	8.3E+03		2.5E+03
1.0E-01	P	7.7E-05	C	3.0E-03	X		V			1.36E+09		1	0.1	Chloro-2-methylaniline, 4-	95-69-2	3.3E+01	7.7E+01	2.2E+05	2.3E+01	3.5E+03	8.3E+03		2.5E+03
2.7E-01	X						V		1.18E+04	1.36E+09	1.62E+04	1		Chloroacetaldehyde, 2-	107-20-0								
						3.0E-05	I			1.36E+09		1	0.1	Chloroacetic Acid	79-11-8								
							V			1.36E+09		1	0.1	Chloroacetophenone, 2-	532-27-4							1.8E+05	1.8E+05
2.0E-01	P			4.0E-03	I		V			1.36E+09		1	0.1	Chloroaniline, p-	106-47-8	1.6E+01	3.9E+01		1.1E+01	4.7E+03	1.1E+04		3.3E+03
				2.0E-02	I	5.0E-02	P	V	7.61E+02	1.36E+09	6.45E+03	1		Chlorobenzene	108-90-7					2.3E+04		1.4E+03	1.3E+03
				1.0E-01	X		V			1.36E+09		1	0.1	Chlorobenzene sulfonic acid, p									

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (see FO #31); H = HEAST; F = See FAQ; W = see user guide Section 2.3.5; E = see user guide Section 2.3.6; L = see user's guide Section 5.2; M = mutagen; S = see user's guide Section 5; V = volatile; R = RBA applied (see user's guide Section 5.10); c = cancer; n = noncancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; SSL values are based on DAF=1; m = Concentration may exceed ceiling limit (see user's guide Section 5.13); s = concentration may exceed Csat (see user's guide Section 5.12)														Contaminant										
Toxicity and Chemical-specific Information														Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k _e (y ⁻¹)	IUR (ug/m ³ -y)	k _e (y ⁻¹)	RFD _o (mg/kg-day)	k _e (y ⁻¹)	RfC _o (mg/m ³)	k _e (y ⁻¹)	muta	gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	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 THI=1 (mg/kg)
				1.0E-02	H						1.36E+09		1	0.1	Chlorpyrifos Methyl	5598-13-0					1.2E+04	2.8E+04		8.2E+03
				5.0E-02	O						1.36E+09		1	0.1	Chlorsulfuron	64902-72-3					5.8E+04	1.4E+05		4.1E+04
				1.0E-02	I						1.36E+09		1	0.1	Chlorthal-dimethyl	1861-32-1					1.2E+04	2.8E+04		8.2E+03
				8.0E-04	H						1.36E+09		1	0.1	Chlorthiophos	60238-56-4					9.3E+02	2.2E+03		6.6E+02
				1.5E+00	I						1.36E+09		0.013		Chromium(III), Insoluble Salts	16065-83-1					1.8E+06			1.8E+06
5.0E-01	C	8.4E-02	S	3.0E-03	I	1.0E-04	I		M		1.36E+09	0.025			Chromium(VI)	18540-29-9	6.5E+00		2.0E+02	6.3E+00	3.5E+03		6.0E+05	3.5E+03
				1.3E-02	I						1.36E+09		0.013		Chromium, Total	7440-47-3					1.5E+04	3.6E+04		1.1E+04
				1.3E-02	I						1.36E+09		1	0.1	Clofentazine	74115-24-5					1.5E+04	3.6E+04		1.1E+04
9.0E-03	P	6.2E-04	I	3.0E-04	P	6.0E-06	P		V	M	1.36E+09		1		Cobalt	7440-48-4		1.9E+03	1.9E+03		3.5E+02		3.6E+04	3.5E+02
				4.0E-02	H						1.36E+09		1		Coke Oven Emissions	8007-45-2							3.6E+04	3.5E+02
				4.0E-02	H						1.36E+09		1		Copper	7440-50-8					4.7E+04			4.7E+04
				5.0E-02	I	6.0E-01	C				1.36E+09		1	0.1	Cresol, m-	108-39-4					5.8E+04	1.4E+05	3.6E+09	4.1E+04
				5.0E-02	I	6.0E-01	C				1.36E+09		1	0.1	Cresol, o-	95-48-7					5.8E+04	1.4E+05	3.6E+09	4.1E+04
				1.0E-01	A	6.0E-01	C				1.36E+09		1	0.1	Cresol, p-	106-44-5					1.2E+05	2.8E+05	3.6E+09	8.2E+04
				1.0E-01	A	6.0E-01	C				1.36E+09		1	0.1	Cresol, p-chloro-m-	59-50-7					1.2E+05	2.8E+05	3.6E+09	8.2E+04
				1.0E-01	A	6.0E-01	C				1.36E+09		1	0.1	Cresols	1319-77-3					1.2E+05	2.8E+05	3.6E+09	8.2E+04
1.9E+00	H			1.0E-03	P		V			1.66E+04	1.36E+09	1.89E+04			Crotonaldehyde, trans-	123-73-9	1.7E+00		1.7E+00		1.2E+03			1.2E+03
				1.0E-01	I	4.0E-01	I	V		2.68E+02	1.36E+09	6.21E+03	1		Cumene	98-82-8					1.2E+05		1.1E+04	9.9E+03
2.2E-01	C	6.3E-05	C	2.0E-03	H						1.36E+09		1	0.1	Cupferron	135-20-6	1.5E+01	3.5E+01	2.6E+05	1.0E+01	2.3E+03	5.5E+03		1.6E+03
8.4E-01	H			2.0E-03	H						1.36E+09		1	0.1	Cyanazine	21725-46-2	3.9E+00	9.2E+00	2.7E+00		2.3E+03	5.5E+03		1.6E+03
				1.0E-03	I						1.36E+09		1		Cyanides						1.2E+03			1.2E+03
				5.0E-03	I						1.36E+09		1		~Calcium Cyanide	592-01-8					5.8E+03			5.8E+03
				6.0E-04	I	8.0E-04	S	V		9.54E+05	1.36E+09	5.33E+04	1		~Copper Cyanide	544-92-3					7.0E+02		1.9E+02	1.5E+02
				1.0E-03	I		V				1.36E+09		1		~Cyanide (CN-)	57-12-5					1.2E+03			1.2E+03
				9.0E-02	I		V				1.36E+09		1		~Cyanogen	460-19-5					1.1E+05			1.1E+05
				5.0E-02	I		V				1.36E+09		1		~Cyanogen Bromide	506-68-3					5.8E+04			5.8E+04
				5.0E-02	I		V				1.36E+09		1		~Cyanogen Chloride	506-77-4					7.0E+02		1.8E+02	1.5E+02
				6.0E-04	I	8.0E-04	I	V		1.00E+07	1.36E+09	5.22E+04	1		~Hydrogen Cyanide	74-90-8					2.3E+03			2.3E+03
				2.0E-03	I						1.36E+09		1		~Potassium Cyanide	151-50-8					5.8E+03			5.8E+03
				5.0E-03	I						1.36E+09		0.04		~Potassium Silver Cyanide	506-61-6					1.2E+05			1.2E+05
				1.0E-01	I						1.36E+09		0.04		~Silver Cyanide	506-64-9					1.2E+03			1.2E+03
				1.0E-03	I						1.36E+09		1		~Sodium Cyanide	143-33-9					1.2E+03			1.2E+03
				2.0E-04	P						1.36E+09		1		~Thiocyanates	E1790664					2.3E+02			2.3E+02
				2.0E-04	X		V				1.36E+09		1		~Thiocyanic Acid	463-56-9					2.3E+02			2.3E+02
				5.0E-02	I						1.36E+09		1		~Zinc Cyanide	557-21-1					5.8E+04			5.8E+04
2.0E-02	X			2.0E-02	X	6.0E+00	I	V		1.17E+02	1.36E+09	1.04E+03	1		Cyclohexane	110-82-7	1.6E+02	3.9E+02	1.1E+02		2.3E+04	5.5E+04	2.7E+04	2.7E+04
				5.0E+00	I	7.0E-01	P	V		5.11E+03	1.36E+09	4.17E+04	1	0.1	Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3					5.8E+06		1.3E+05	1.3E+05
				5.0E-03	P	1.0E+00	X	V		2.83E+02	1.36E+09	1.46E+03	1		Cyclohexanone	108-94-1					5.8E+03		6.4E+03	3.1E+03
				2.0E-01	I		V			2.93E+05	1.36E+09	7.46E+04	1		Cyclohexene	110-83-8					2.3E+05			2.3E+05
				2.5E-02	I						1.36E+09		1	0.1	Cyclohexylamine	108-91-8					2.9E+04	6.9E+04		2.1E+04
				1.0E-03	O						1.36E+09		1	0.1	Cyfluthrin	68359-37-5					1.2E+03			1.2E+03
				5.0E-01	O						1.36E+09		1	0.1	Cyhalothrin	68085-85-8					5.8E+05	1.4E+06		4.1E+05
2.4E-01	I	6.9E-05	C	3.0E-05	X						1.36E+09		1	0.1	Cyromazine	66215-27-8					3.5E+01	8.3E+01		2.5E+01
3.4E-01	I	9.7E-05	C	3.0E-04	X		V				1.36E+09	2.10E+06	1		DDD, p,p'- (DDD)	72-54-8	1.4E+01	3.2E+01	2.4E+05	9.6E+00	3.5E+02			3.5E+02
3.4E-01	I	9.7E-05	I	5.0E-04	I						1.36E+09		1	0.03	DDE, p,p'-	72-55-9	9.6E+00	2.7E+02	2.7E+02	9.3E+00	5.8E+02			5.8E+02
				3.0E-02	I						1.36E+09		1	0.1	DDT	50-29-3	9.6E+00	7.6E+01	1.7E+05	8.5E+00	5.8E+02	4.6E+03		5.2E+02
				1.8E-02	C	5.1E-06	C				1.36E+09		1	0.1	Dalapon	75-99-0					3.5E+04	8.3E+04		2.5E+04
7.0E-04	I			7.0E-03	I						1.36E+09		1	0.1	Daminozide	1596-84-5	1.8E+02	4.3E+02	3.3E+06	1.3E+02	1.8E+05	4.1E+05		1.2E+05
				4.0E-05	I						1.36E+09		1	0.1	Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5	4.7E+03	1.1E+04		3.3E+03	8.2E+03	1.9E+04		5.7E+03
				6.0E-01	I						1.36E+09		1	0.1	Dermeton	8065-48-3					4.7E+01	1.1E+02		3.3E+01
1.2E-03	I			6.0E-01	I						1.36E+09		1	0.1	Di(2-ethylhexyl)adipate	103-23-1	2.7E+03	6.4E+03		1.9E+03	7.0E+05	1.7E+06		4.9E+05
6.1E-02	H			7.0E-04	A						1.36E+09		1	0.1	Diallate	2303-16-4	5.4E+01	1.3E+02		3.8E+01	8.2E+02	1.9E+03		5.7E+02
				1.0E-02	X		V				1.36E+09	5.24E+05	1		Diazinon	333-41-5					1.2E+04			1.2E+04
8.0E-01	P	6.0E-03	P	2.0E-04	P	2.0E-04	I	V	M	9.79E+02	1.36E+09	3.20E+04	1		Dibenzothioephene	132-65-0	4.1E+00		6.5E-02	6.4E-02	2.3E+02		2.8E+01	1.2E+04
				4.0E-04	X		V			1.59E+02	1.36E+09	1.93E+04	1		Dibromo-3-chloropropane, 1,2-	96-12-8					4.7E+02			4.7E+02
				1.0E-02	I		V																	

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (see FO #31); H = HEAST; F = See FAQ; W = see user guide Section 2.3.5; E = see user guide Section 2.3.6; L = see user's guide Section 5.2; M = mutagen; S = see user's guide Section 5; V = volatile; R = RBA applied (see user's guide Section 5.10); c = cancer; n = noncancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; SSL values are based on DAF=1; n = Concentration may exceed ceiling limit (see user's guide Section 5.13); s = concentration may exceed Csat (see user's guide Section 5.12)														Toxicity and Chemical-specific Information														Contaminant														Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1			
SFO	k _e	IUR	k _e	RFD _c	k _e	RfC _c	k _e	Vo	muta	C _{sat}	PEF	VF	GIABS	ABS	Analyte	CAS No.	Ingestion SL	Dermal SL	Inhalation SL	Carcinogenic SL	Ingestion SL	Dermal SL	Inhalation SL	Noncarcinogenic SL	THQ=1	THQ=1	THQ=1	THI=1																					
(mg/kg-day) ⁻¹	y	(ug/m ³ -y)	y	(mg/kg-day)	y	(mg/m ³)	y	y	gen	(mg/kg)	(m ³ /kg)	(m ³ /kg)					TR=1E-06	TR=1E-06	TR=1E-06	TR=1E-06	THQ=1	THQ=1	THQ=1	THQ=1	THQ=1	THQ=1	THQ=1	THQ=1																					
				2.0E-03	I			V		2.37E+03	1.36E+09	2.50E+03	1		Dichloroethylene, 1,2-cis-	156-59-2						2.3E+03						2.3E+03																					
				2.0E-02	I			V		1.85E+03	1.36E+09	1.75E+03	1		Dichloroethylene, 1,2-trans-	156-60-5						2.3E+04					2.3E+04																						
				3.0E-03	I						1.36E+09		1	0.1	Dichlorophenol, 2,4-	120-83-2						3.5E+03	8.3E+03				2.5E+03																						
3.7E-02	P	3.7E-06	P	1.0E-02	I						1.36E+09		1	0.05	Dichlorophenoxy Acetic Acid, 2,4-	94-75-7						1.2E+04	5.5E+04				9.6E+03																						
				4.0E-02	P	4.0E-03	I	V		1.36E+03	1.36E+09	3.79E+03	1		Dichloropropane, 1,2-	78-87-5	8.8E+01		1.3E+01	1.1E+01		4.7E+04		6.6E+01		6.6E+01																							
				2.0E-02	P			V		1.49E+03	1.36E+09	6.76E+03	1		Dichloropropane, 1,3-	142-28-9						2.3E+04					2.3E+04																						
1.0E-01	I	4.0E-06	I	3.0E-03	I						1.36E+09		1	0.1	Dichloropropanol, 2,3-	616-23-9	3.3E+01		1.1E+01	8.2E+00		3.5E+03	8.3E+03		3.1E+02	2.5E+03																							
2.9E-01	I	8.3E-05	C	5.0E-04	I	5.0E-04	I				1.36E+09		1	0.1	Dichloropropene, 1,3-	542-75-6	1.1E+01	2.7E+01	2.0E+05	7.9E+00		3.5E+04	1.4E+03	3.0E+06	3.1E+02	4.1E+02																							
				3.0E-05	O						1.36E+09		1	0.1	Dichlorvos	62-73-7						5.8E+02	1.4E+03			4.1E+02																							
1.6E+01	I	4.6E-03	I	8.0E-02	P	3.0E-04	X	V		2.56E+02	1.36E+09	4.11E+03	1		Dicrotophos	141-66-2	2.0E-01	4.8E-01	3.6E+03	1.4E-01		3.5E+01	8.3E+01			2.5E+01																							
				5.0E-05	I						1.36E+09		1	0.1	Dicyclopentadiene	77-73-6						9.3E+04		5.4E+00			5.4E+00																						
				3.0E-04	C						1.36E+09		1	0.1	Dieldrin	60-57-1						5.8E+01	1.4E+02				4.1E+01																						
				2.0E-03	P	2.0E-04	P				1.36E+09		1	0.1	Diesel Engine Exhaust	E17136615											1.6E+03																						
				3.0E-02	P	1.0E-04	P				1.36E+09		1	0.1	Diethanolamine	111-42-2											6.0E+05																						
				6.0E-02	P	3.0E-04	P				1.36E+09		1	0.1	Diethylene Glycol Monobutyl Ether	112-34-5						7.0E+04	1.7E+05	1.8E+06			2.4E+04																						
3.5E+02	C	1.0E-01	C	1.0E-03	P			V		1.12E+05	1.36E+09	1.39E+05	1		Diethylene Glycol Monoethyl Ether	111-90-0	9.3E-03	2.2E-02	1.7E+02	6.6E-03		1.2E+03				4.8E+04																							
				8.3E-02	O						1.36E+09		1	0.1	Diethylformamide	617-84-5											1.2E+03																						
				2.0E-02	I						1.36E+09		1	0.1	Diethylstilbestrol	56-53-1											1.2E+03																						
				4.0E+01	I	V				1.43E+03	1.36E+09	1.15E+03	1		Difluoropropane, 2,2-	43222-48-6						9.7E+04	2.3E+05				6.8E+04																						
4.4E-02	C	1.3E-05	C	3.0E+01	X	V				6.91E+02	1.36E+09	7.58E+02	1		Difluoropropane, 2,2-	420-45-1	7.4E+01		1.2E+02	4.5E+01		2.3E+04	5.5E+04		2.0E+05	1.0E+05	2.0E+05																						
				7.0E-01	P	V				2.26E+03	1.36E+09	3.06E+03	1		Dihydroisoflavanone	94-58-6											1.0E+05																						
				8.0E-02	I			V		5.30E+02	1.36E+09	3.81E+04	1		Diisopropyl Ether	108-20-3											9.4E+03																						
				2.2E-02	O						1.36E+09		1	0.1	Diisopropyl Methylphosphonate	1445-75-6						9.3E+04					9.3E+04																						
				2.2E-03	O						1.36E+09		1	0.1	Dimethipin	55290-64-7						2.5E+04	6.0E+04				1.8E+04																						
1.6E+00	P			6.0E-02	P						1.36E+09		1	0.1	Dimethoate	60-51-5	2.0E+00	4.8E+00		1.4E+00		2.6E+03	6.1E+03				1.8E+03																						
1.7E-03	P										1.36E+09		1	0.1	Dimethoxybenzidine, 3,3'-	119-90-4	1.9E+03	4.5E+03		1.4E+03							4.9E+04																						
4.6E+00	C	1.3E-03	C								1.36E+09		1	0.1	Dimethyl methylphosphonate	756-79-6	7.1E-01	1.7E+00	1.3E+04	5.0E-01		7.0E+04	1.7E+05																										
5.8E-01	H										1.36E+09		1	0.1	Dimethylamino azobenzene [p-]	60-11-7	5.6E+00	1.3E+01		4.0E+00							1.6E+03																						
2.0E-01	P			2.0E-03	X						1.36E+09		1	0.1	Dimethylaniline HCl, 2,4-	21436-96-4	1.6E+01	3.9E+01		1.1E+01		2.3E+03	5.5E+03				2.3E+03																						
2.7E-02	P			2.0E-03	I			V		8.30E+02	1.36E+09	3.13E+04	1		Dimethylaniline, N,N-	121-69-7	1.2E+02			1.2E+02																													
1.1E+01	P										1.36E+09		1	0.1	Dimethylaniline, N,N-	121-69-7	3.0E-01	7.0E-01		2.1E-01							1.6E+03																						
				1.0E-01	P	3.0E-02	I	V		1.06E+05	1.36E+09	1.28E+05	1		Dimethylbenzidine, 3,3'-	119-93-7						1.2E+05		1.7E+04			1.5E+04																						
				1.0E-04	X	2.0E-06	X	V		1.72E+05	1.36E+09	2.77E+04	1		Dimethylformamide	68-12-2											2.4E-01																						
5.5E+02	C	1.6E-01	C					V		1.89E+05	1.36E+09	1.68E+05	1		Dimethylhydrazine, 1,1-	57-14-7	5.9E-03		1.3E-02	4.1E-03		1.2E+02		2.4E-01																									
				2.0E-02	I						1.36E+09		1	0.1	Dimethylhydrazine, 1,2-	540-73-8											1.6E+04																						
				6.0E-04	I						1.36E+09		1	0.1	Dimethylphenol, 2,4-	105-67-9						7.0E+02	1.7E+03				4.9E+02																						
				1.0E-03	I						1.36E+09		1	0.1	Dimethylphenol, 2,6-	576-26-1											8.2E+02																						
4.5E-02	C	1.3E-05	C					V		4.73E+02	1.36E+09	5.48E+03	1		Dimethylphenol, 3,4-	95-65-8	7.3E+01		5.2E+00	4.8E+00		1.2E+03	2.8E+03			8.2E+02																							
				8.0E-05	X						1.36E+09		1	0.1	Dimethylvinylchloride	513-37-1											8.2E+02																						
				2.0E-03	I						1.36E+09		1	0.1	Dinitro-o-cresol, 4,6-	534-52-1						9.3E+01	2.2E+02				6.6E+01																						
				1.0E-04	P						1.36E+09		1	0.1	Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5						2.3E+03	5.5E+03				1.6E+03																						
				1.0E-04	P						1.36E+09		1	0.1	Dinitrobenzene, 1,2-	528-29-0						1.2E+02	2.8E+02				8.2E+01																						
				1.0E-04	P						1.36E+09		1	0.1	Dinitrobenzene, 1,3-	99-65-0						1.2E+02	2.8E+02				8.2E+01																						
				1.0E-04	P						1.36E+09		1	0.1	Dinitrobenzene, 1,4-	100-25-4						1.2E+02	2.8E+02				8.2E+01																						
6.8E-01	I			2.0E-03	I						1.36E+09		1	0.1	Dinitrophenol, 2,4-	51-28-5	4.8E+00	1.1E+01		3.4E+00		2.3E+03	5.5E+03				1.6E+03																						
3.1E-01	C	8.9E-05	C	2.0E-03	I						1.36E+09		1	0.102	Dinitrotoluene, 2,4,6-	E1615210	1.1E+01	2.4E+01	1.9E+05	7.4E+00		2.3E+03	5.4E+03				1.6E+03																						
1.5E+00	P			3.0E-04	X						1.36E+09		1	0.099	Dinitrotoluene, 2,6-	606-20-2	2.2E+00	5.2E+00		1.5E+00		3.5E+02	8.4E+02				2.5E+02																						
				2.0E-03	S						1.36E+09		1	0.006	Dinitrotoluene, 2-Amino-4,6-	35572-78-2						2.3E+03	9.2E+04				2.3E+03																						
4.5E-01	X			2.0E-03	S						1.36E+09		1	0.009	Dinitrotoluene, 4-Amino-2,6-	19406-51-0	7.3E+00	1.7E+01		5.1E+00		2.3E+03	6.1E+04				2.3E+03																						
				9.0E-04	X						1.36E+09		1	0.1	Dinitrotoluene, Technical grade	25321-14-6						1.1E+03	2.5E+03				7.4E+02																						

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (see FO #31); H = HEAST; F = See FAQ; W = see user guide Section 2.3.5; E = see user guide Section 2.3.6; L = see user's guide Section 5.2; M = mutagen; S = see user's guide Section 5; V = volatile; R = RBA applied (see user's guide Section 5.10); c = cancer; n = noncancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; SSL values are based on DAF=1; n = Concentration may exceed ceiling limit (see user's guide Section 5.13); s = concentration may exceed Csat (see user's guide Section 5.12)														Contaminant										Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1			
Toxicity and Chemical-specific Information														Analyte										Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -y) ⁻¹	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³)	k _e (y)	muta	gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	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 THI=1 (mg/kg)								
9.9E-03	I	1.2E-06	I	3.0E-04	I	6.0E-03	P	1.0E-03	I	V	1.05E+04	1.36E+09	1.89E+04	1	0.1	72-20-8	3.3E+02		1.9E+02	1.2E+02	3.5E+02	8.3E+02		2.5E+02							
				6.0E-03	I	2.0E-02	I	V			1.53E+04	1.36E+09	7.66E+03	1		106-89-8							8.3E+01	8.2E+01							
				4.0E-02	P						1.36E+09		1	0.1	106-88-7					4.7E+04	1.1E+05		6.7E+02	3.3E+04							
				5.0E-03	I						1.36E+09		1	0.1	111-77-3					5.8E+03	1.4E+04		4.1E+03								
				5.0E-04	I						1.36E+09		1	0.1	16672-87-0					5.8E+02	1.4E+03		4.1E+02								
				1.0E-01	P	6.0E-02	P	V		2.38E+04	1.36E+09	6.15E+04	1		563-12-2					1.2E+05		1.6E+04	1.4E+04								
				9.0E-02	P	2.0E-01	I	V		1.06E+05	1.36E+09	9.84E+04	1		111-15-9					1.1E+05		8.6E+04	4.7E+04								
				9.0E-01	I	7.0E-02	P	V		1.08E+04	1.36E+09	8.62E+03	1		141-78-6					1.1E+06		2.6E+03	2.6E+03								
				5.0E-03	P	8.0E-03	P	V		2.50E+03	1.36E+09	6.34E+03	1		140-88-5					5.8E+03		2.2E+02	2.1E+02								
				2.0E-01	I	1.0E+01	I	V		2.12E+03	1.36E+09	1.29E+03	1		75-00-3							5.7E+04	5.7E+04								
											1.01E+04	1.36E+09	3.12E+03	1		60-29-7						2.3E+05	2.3E+05								
											1.10E+03	1.36E+09	5.77E+03	1		97-63-2						7.6E+03	7.6E+03								
1.1E-02	C	2.5E-06	C	1.0E-05	I						1.36E+09		1	0.1	2104-64-5	3.0E+02		2.8E+01	2.5E+01	1.2E+01	2.8E+01		8.2E+00								
				1.0E-01	I	1.0E+00	I	V		4.80E+02	1.36E+09	5.67E+03	1		100-41-4					1.2E+05		2.5E+04	2.0E+04								
				7.0E-02	P						1.36E+09		1	0.1	109-78-4					8.2E+04	1.9E+05		5.7E+04								
				9.0E-02	P					1.89E+05	1.36E+09	1.80E+05	1		107-15-3					1.1E+05			1.1E+05								
				2.0E+00	I	4.0E-01	C				1.36E+09		1	0.1	107-21-1					2.3E+06	5.5E+06	2.4E+09	1.6E+06								
				1.0E-01	I	1.6E+00	I				1.36E+09		1	0.1	111-76-2					1.2E+05	2.8E+05	9.5E+09	8.2E+04								
3.1E-01	C	3.0E-03	I	3.0E-02	C	V	M			1.21E+05	1.36E+09	6.09E+03	1		75-21-8	1.1E+01		2.5E-02	2.5E-02				8.0E+02								
4.5E-02	C	1.3E-05	C	8.0E-05	I						1.36E+09		1	0.1	96-45-7	7.3E+01	1.7E+02	1.3E+06	5.1E+01	9.3E+01	2.2E+02		8.0E+02	6.6E+01							
6.5E+01	C	1.9E-02	C							1.54E+05	1.36E+09	2.39E+04	1		151-86-4	5.0E-02		1.5E-02	1.2E-02												
				3.0E+00	I						1.36E+09		1	0.1	84-72-0					3.5E+06	8.3E+06		2.5E+06								
				2.5E-04	I						1.36E+09		1	0.1	22224-92-6					2.9E+02	6.9E+02		2.1E+02								
				2.5E-02	I						1.36E+09		1	0.1	39515-41-8					2.9E+04	6.9E+04		2.1E+04								
				2.5E-02	I						1.36E+09		1	0.1	51630-58-1					2.9E+04	6.9E+04		2.1E+04								
				1.3E-02	I						1.36E+09		1	0.1	2164-17-2					1.5E+04	3.6E+04		1.1E+04								
				4.0E-02	C	1.3E-02	C				1.36E+09		1	0.1	16984-48-8					4.7E+04		7.7E+07	4.7E+04								
				6.0E-02	I	1.3E-02	C				1.36E+09		1		7782-41-4					7.0E+04		7.7E+07	7.0E+04								
				8.0E-02	I						1.36E+09		1	0.1	59756-80-4					9.3E+04	2.2E+05		6.6E+04								
				4.0E-02	O						1.36E+09		1	0.1	56425-91-3					4.7E+04	1.1E+05		3.3E+04								
				2.0E-03	O						1.36E+09		1	0.1	85509-19-9					2.3E+03	5.5E+03		1.6E+03								
				5.0E-01	O						1.36E+09		1	0.1	66332-96-5					5.8E+05	1.4E+06		4.1E+05								
				1.0E-02	I						1.36E+09		1	0.1	69409-94-5					1.2E+04	2.8E+04		8.2E+03								
				9.0E-02	O						1.36E+09		1	0.1	133-07-3					1.1E+05	2.5E+05		7.4E+04								
				2.5E-03	O						1.36E+09		1	0.1	72178-02-0					2.9E+03	6.9E+03		2.1E+03								
				2.0E-03	I						1.36E+09		1	0.1	944-22-9					2.3E+03	5.5E+03		1.6E+03								
1.3E-05	I			2.0E-01	I	9.8E-03	A	V		4.24E+04	1.36E+09	7.77E+04	1		50-00-0			7.3E+01	7.3E+01	2.3E+05		3.3E+03	3.3E+03								
				9.0E-01	P	3.0E-04	X	V		1.06E+05	1.36E+09	9.30E+04	1		64-18-6					1.1E+06		1.2E+02	1.2E+02								
				2.5E+00	O						1.36E+09		1	0.1	39148-24-8					2.9E+06	6.9E+06		2.1E+06								
				1.0E-03	X						1.36E+09	1.56E+05	1	0.03	132-64-9					1.2E+03	9.2E+03		1.0E+03								
				1.0E-03	I					6.22E+03	1.36E+09	2.62E+03	1	0.03	110-00-9					1.2E+03	9.2E+03		1.0E+03								
3.8E+00	H			9.0E-01	I	2.0E+00	I	V		1.65E+05	1.36E+09	1.20E+04	1	0.03	109-99-9	8.6E-01	2.0E+00		6.0E-01	1.1E+06	8.3E+06	1.0E+05	9.4E+04								
				3.0E-03	I	5.0E-02	H	V		1.01E+04	1.36E+09	4.86E+04	1		67-45-8					3.5E+03		1.1E+04	2.6E+03								
											1.36E+09		1	0.1	98-01-1																
1.5E+00	C	4.3E-04	C	6.0E-03	O						1.36E+09		1	0.1	531-82-8	2.2E+00	5.2E+00	3.9E+04	1.5E+00				4.9E+03								
3.0E-02	I	8.6E-06	C	1.0E-01	I						1.36E+09		1	0.1	60568-05-0	1.1E+02	2.6E+02	1.9E+06	7.7E+01												
				1.0E-01	I						1.36E+09		1	0.1	77182-82-2					7.0E+03	1.7E+04		4.9E+03								
				1.0E-01	A	8.0E-05	C				1.36E+09		1	0.1	111-30-8					1.2E+05	2.8E+05	4.8E+05	7.0E+04								
				4.0E-04	I	1.0E-03	H	V		1.06E+05	1.36E+09	8.43E+04	1		765-34-4					4.7E+02		3.7E+02	2.1E+02								
				1.0E-01	I						1.36E+09		1	0.1	1071-83-6					1.2E+05	2.8E+05		8.2E+04								
				1.0E-02	X						1.36E+09	1.45E+05	1		113-00-8					1.2E+04			1.2E+04								
				2.0E-02	P						1.36E+09		1	0.1	50-01-1					2.3E+04	5.5E+04		1.6E+04								
				3.0E-02	X						1.36E+09		1	0.1	506-93-4					3.5E+04	8.3E+04		2.5E+04								
4.5E+00	I	1.3E-03	I	5.0E-05	I						1.36E+09		1	0.1	69806-40-2	7.3E-01		4.5E+00	6.3E-01	5.8E+01	1.4E+02		4.1E+01								
9.1E+00	I	2.6E-03	I	1.3E-05	I						1.36E+09	8.43E+05	1		76-44-8	3.6E-01		4.0E+00	3.3E-01	5.8E+02			5.8E+02								
											1.36E+09		1	0.1	1024-57-3					1.5E+01			1.5E+01								
				3.0E-04	X	3.0E-03	X	V		2.09E+02	1.																				

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (see FO #31); H = HEAST; F = See FAQ; W = see user guide Section 2.3.5; E = see user guide Section 2.3.6; L = see user's guide Section 5.2; M = mutagen; S = see user's guide Section 5; V = volatile; R = RBA applied (see user's guide Section 5.10); c = cancer; n = noncancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; SSL values are based on DAF=1; n = Concentration may exceed ceiling limit (see user's guide Section 5.13); s = concentration may exceed Csat (see user's guide Section 5.12)

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1				
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -y)	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³)	k _e (y)	muta gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	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 THI=1 (mg/kg)
				2.0E+00	P	7.0E-01	I V		1.41E+02	1.36E+09	8.29E+02	1	0.1	Hexane, N-Hexanedioic Acid	110-54-3					2.3E+06	5.5E+06	2.5E+03	2.5E+03
				5.0E-03	I	3.0E-02	I V		3.28E+03	1.36E+09	1.33E+04	1		Hexanone, 2-Hexazinone	591-78-6					5.8E+03		1.7E+03	1.3E+03
				3.3E-02	I					1.36E+09		1	0.1	Hexythiazox	51235-04-2					3.9E+04	9.1E+04		2.7E+04
				2.5E-02	I					1.36E+09		1	0.1	Hydramethylnon	78587-05-0					2.9E+04	6.9E+04		2.1E+04
3.0E+00	I	4.9E-03	I	1.7E-02	O				1.12E+05	1.36E+09	6.51E+04	1		Hydrazine	67485-29-4					2.0E+04	4.7E+04		1.4E+04
3.0E+00	I	4.9E-03	I			3.0E-05	P V			1.36E+09		1		Hydrazine Sulfate	302-01-2	1.1E+00		1.6E-01	1.4E-01			8.6E+00	8.6E+00
				4.0E-02	C	2.0E-02	I V			1.36E+09		1		Hydrogen Chloride	7647-01-0					4.7E+04		1.2E+08	1.2E+08
						1.4E-02	C V			1.36E+09		1		Hydrogen Fluoride	7664-39-3							8.3E+07	4.7E+04
						2.0E-03	I V			1.36E+09		1		Hydrogen Sulfide	7783-06-4							1.2E+07	1.2E+07
6.0E-02	P			4.0E-02	P					1.36E+09		1	0.1	Hydroquinone	123-31-9	5.5E+01	1.3E+02		3.8E+01	4.7E+04	1.1E+05		3.3E+04
6.1E-02	O			2.5E-03	O					1.36E+09		1	0.1	Imazalil	35554-44-0	5.4E+01	1.3E+02		3.8E+01	2.9E+03	6.9E+03		2.1E+03
				2.5E-01	I					1.36E+09		1	0.1	Imazazaquin	81335-37-7					2.9E+05	6.9E+05		2.1E+05
				2.5E+00	O					1.36E+09		1	0.1	Imazethapyr	81335-77-5					2.9E+06	6.9E+06		2.1E+06
				1.0E-02	A					1.36E+09		1		Iodine	7553-56-2					1.2E+04			1.2E+04
				4.0E-02	I					1.36E+09		1	0.1	Iprodione	36734-19-7					4.7E+04	1.1E+05		3.3E+04
				7.0E-01	P					1.36E+09		1		Iron	7439-89-6					8.2E+05			8.2E+05
				3.0E-01	I		V		1.00E+04	1.36E+09	2.81E+04	1		Isobutyl Alcohol	78-83-1					3.5E+05			3.5E+05
9.5E-04	I			2.0E-01	I	2.0E+00	C			1.36E+09		1	0.1	Isophorone	78-59-1	3.4E+03	8.1E+03		2.4E+03	2.3E+05	5.5E+05	1.2E+10	1.6E+05
				1.5E-02	I		V			1.36E+09	4.20E+05	1		Isopropalin	33820-53-0					1.8E+04			1.8E+04
				2.0E+00	P	2.0E-01	P V		1.09E+05	1.36E+09	2.77E+04	1		Isopropanol	67-63-0					2.3E+06		2.4E+04	2.4E+04
				1.0E-01	I					1.36E+09		1	0.1	Isopropyl Methyl Phosphonic Acid	1832-54-8					1.2E+05	2.8E+05		8.2E+04
				5.0E-02	I					1.36E+09		1	0.1	Isosaberenol	82558-50-7					5.8E+04	1.4E+05		4.1E+04
						3.0E-01	A V			1.36E+09		1		JP-7	E 1737665							1.8E+09	1.8E+09
				8.0E-03	O					1.36E+09		1	0.1	Lactofen	77501-63-4					9.3E+03	2.2E+04		6.6E+03
				2.0E-04	X					1.36E+09		1		Lactonitrile	78-97-7					2.3E+02	5.5E+02		1.6E+02
				5.0E-05	P					1.36E+09		1		Lanthanum	7439-91-0					5.8E+01			5.8E+01
				2.1E-05	P					1.36E+09		1	0.1	Lanthanum Acetate Hydrate	100587-80-4					2.4E+01	5.7E+01		1.7E+01
				1.9E-05	P					1.36E+09		1		Lanthanum Chloride Heptahydrate	10025-84-0					2.2E+01			2.2E+01
				2.8E-05	P					1.36E+09		1		Lanthanum Chloride, Anhydrous	10099-58-8					3.3E+01			3.3E+01
				1.6E-05	P					1.36E+09		1		Lanthanum Nitrate Hexahydrate	10277-43-7					1.9E+01			1.9E+01
8.5E-03	C	1.2E-05	C							1.36E+09		1		Lead Compounds		3.8E+02		1.4E+06	3.8E+02				
8.5E-03	C	1.2E-05	C							1.36E+09		1	0.1	~Lead Phosphate	7446-27-7	3.8E+02		1.4E+06	3.8E+02				
										1.36E+09		1	0.1	~Lead acetate	301-04-2	3.8E+02	9.1E+02	1.4E+06	2.7E+02				8.0E+02
8.5E-03	C	1.2E-05	C							1.36E+09		1	0.1	~Lead and Compounds	7439-92-1	3.8E+02	9.1E+02	1.4E+06	2.7E+02				
				1.0E-07	I		V		2.43E+00	1.36E+09	1.91E+03	1		~Lead subacetate	1335-32-6					1.2E-01			1.2E-01
				5.0E-06	P		V		3.83E+02	1.36E+09	2.55E+04	1		~Tetraethyl Lead	78-00-2								
				7.7E-03	O					1.36E+09		1	0.1	Lewisite	541-25-3					9.0E+03	2.1E+04		6.3E+03
				2.0E-03	P					1.36E+09		1		Linuron	330-55-2					2.3E+03			2.3E+03
				5.0E-04	I					1.36E+09		1	0.1	Lithium	7439-93-2								
				4.4E-03	O					1.36E+09		1	0.1	MCPA	94-74-6					5.8E+02	1.4E+03		4.1E+02
				1.0E-03	I					1.36E+09		1	0.1	MCPB	94-81-5					5.1E+03	1.2E+04		3.6E+03
				1.0E-03	I					1.36E+09		1	0.1	MCPP	93-65-2					1.2E+03	2.8E+03		8.2E+02
				2.0E-02	I					1.36E+09		1	0.1	Malathion	121-75-5					2.3E+04	5.5E+04	4.2E+06	1.6E+04
				1.0E-01	I	7.0E-04	C			1.36E+09		1	0.1	Maleic Anhydride	108-31-6					1.2E+05	2.8E+05		8.0E+04
				5.0E-01	I					1.36E+09		1	0.1	Maleic Hydrazide	123-33-1					5.8E+05	1.4E+06		4.1E+05
				1.0E-04	P					1.36E+09		1	0.1	Malononitrile	109-77-3					1.2E+02	2.8E+02		8.2E+01
				3.0E-02	H					1.36E+09		1	0.1	Mancozeb	8018-01-7					3.5E+04	8.3E+04		2.5E+04
				5.0E-03	I					1.36E+09		1	0.1	Maneb	12427-38-2					5.8E+03	1.4E+04		4.1E+03
				1.4E-01	I	5.0E-05	I			1.36E+09		1		Manganese (Diet)	7439-96-5					2.8E+04		3.0E+05	2.6E+04
				2.4E-02	S	5.0E-05	I			1.36E+09	0.04	1		Manganese (Non-diet)	7439-96-5					1.1E+02	2.5E+02		7.4E+01
				9.0E-05	H					1.36E+09		1	0.1	Mephsfolan	950-10-7					3.5E+04	8.3E+04		2.5E+04
1.1E-02	P			3.0E-02	I					1.36E+09		1	0.1	Mepiquat Chloride	24307-26-4	3.0E+02		7.0E+02	2.1E+02	4.7E+03	1.1E+04		3.3E+03
				4.0E-03	P					1.36E+09		1	0.1	Mercaptobenzothiazole, 2-Mercuro Compounds	149-30-4								
				3.0E-04	I	3.0E-04	S			1.36E+09	0.07	1		~Mercuric Chloride (and other Mercury salts)	7487-94-7					3.5E+02		1.8E+06	3.5E+02
						3.0E-04	I V		3.13E+00	1.36E+09	3.47E+04	1		~Mercury (elemental)	7439-97-6						4.6E+01		4.6E+01
				1.0E-04	I					1.36E+09		1		~Methyl Mercury	22967-92-6					1.2E+02			1.2E+02
				8.0E-05	I					1.36E+09		1	0.1	~Phenylmercuric Acetate	62-38-4					9.3E+01	2.2E+02		6.6E+01
				3.0E-05	I		V			1.36E+09	1.94E+06	1		Merphos	150-50-5					3.5E+01			3.5E+01
				1.0E-04	O					1.36E+09		1	0.1	Merphos Oxide	78-48-8					1.2E+02	2.8E+02		8.2E+01
				6.0E-02	I					1.36E+09		1											

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (see FO #31); H = HEAST; F = See FAQ; W = see user guide Section 2.3.5; E = see user guide Section 2.3.6; L = see user's guide Section 5.2; M = mutagen; S = see user's guide Section 5; V = volatile; R = RBA applied (see user's guide Section 5.10); c = cancer; n = noncancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; SSL values are based on DAF=1; n = Concentration may exceed ceiling limit (see user's guide Section 5.13); s = concentration may exceed Csat (see user's guide Section 5.12)

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k _e IUR (ug/m ³ -y) ⁻¹	k _e RfD _o (mg/kg-day)	k _e RfC _o (mg/m ³)	k _e V _o I	muta gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	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 THI=1 (mg/kg)			
	1.0E-03 X	6.0E-01 I 1.0E-03 P	5.0E+00 I 2.0E-05 X 3.0E+00 I	V		2.84E+04 1.80E+05 3.36E+03	1.36E+09 1.36E+09 1.36E+09	1.22E+04 5.04E+04 1.06E+04	1		Methyl Ethyl Ketone (2-Butanone) Methyl Hydrazine Methyl Isobutyl Ketone (4-methyl-2-pentanone)	78-93-3 60-34-4 108-10-1			6.2E-01	6.2E-01	7.0E+05 1.2E+03		2.7E+05 4.4E+00 1.4E+05	1.9E+05 4.4E+00 1.4E+05			
		1.4E+00 I 2.5E-04 I	1.0E-03 C 7.0E-01 I	V		1.01E+04 2.36E+03	1.36E+09 1.36E+09	4.42E+03 6.33E+03	1	0.1	Methyl Isocyanate Methyl Methacrylate Methyl Parathion	624-83-9 80-62-6 298-00-0					1.6E+06 2.9E+02	6.9E+02	1.9E+01 1.9E+04	1.9E+01 1.9E+04			
9.9E-02 C 1.8E-03 C	2.8E-05 C 2.6E-07 C		3.0E+00 I 3.0E-04 X	V		8.87E+03 1.36E+09	4.90E+03 1.36E+09		1	0.1	Methyl tert-Butyl Ether (MTBE) Methyl-1,4-benzenediamine dihydrochloride, 2- Methyl-2-Pentanol, 4-	1634-04-4 615-45-2 108-11-2	3.3E+01 1.8E+03	7.8E+01 2.3E+02	6.0E+05 2.1E+02	2.3E+01		6.4E+04	6.4E+04	2.5E+02 2.3E+05			
9.0E-03 P 8.3E+00 C 1.3E-01 C	2.4E-03 C 3.7E-05 C	2.0E-02 X				1.36E+09 1.36E+09 1.36E+09			1	0.1	Methyl-5-Nitroaniline, 2- Methyl-N-nitro-N-nitrosoguanidine, N- Methylaniline Hydrochloride, 2-	99-55-8 70-25-7 636-21-5	3.6E+02 3.9E-01 2.5E+01	8.6E+02 9.3E-01 5.9E+01	6.9E+03 2.3E-01 4.5E+05	2.6E+02 2.6E-01 1.8E+01	2.3E+04 5.5E+04		1.6E+04				
1.0E-01 X		1.0E-02 A 2.0E-04 X 3.0E-04 X				1.36E+09 1.36E+09 1.36E+09			1	0.1	Methylarsonic acid Methylbenzene, 1,4-diamine monohydrochloride, 2- Methylbenzene-1,4-diamine sulfate, 2-	124-58-3 74612-12-7 615-50-9	3.3E+01 1.5E-01 1.6E+03	7.7E+01 3.5E-01 2.7E+03	2.6E+03 1.0E-01 1.0E+03	2.3E+01		2.3E+03 3.5E+02 8.3E+02	5.5E+02 8.3E+02	8.2E+03 1.6E+02 2.5E+02			
2.2E+01 C 2.0E-03 I 1.0E-01 P	6.3E-03 C 4.3E-04 C	6.0E-03 I 2.0E-03 P	6.0E-01 I 2.0E-03 P	V	M	3.32E+03 1.36E+09	2.19E+03 1.36E+09		1	0.1	Methylcholanthrene, 3- Methylene Chloride Methylene-bis(2-chloroaniline), 4,4'-	56-49-5 75-09-2 101-14-4	1.5E-01 1.6E+03 3.3E+01	3.5E-01 2.7E+03 7.7E+01	2.6E+03 1.0E+03 3.9E+04	1.0E-01 1.0E+03 2.3E+01	7.0E+03 2.3E+03 5.5E+03	5.8E+03	3.2E+03 1.6E+03				
4.6E-02 I 1.6E+00 C	1.3E-05 C 4.6E-04 C		2.0E-02 C 6.0E-04 I			1.36E+09 1.36E+09			1	0.1	Methylene-bis(N,N-dimethyl) Aniline, 4,4'- Methylenbisbenzenamine, 4,4'- Methylenediphenyl Diisocyanate	101-61-1 101-77-9 101-68-8	7.1E+01 2.0E+00	1.7E+02 4.8E+00	1.3E+06 3.6E+04	5.0E+01 1.4E+00		1.2E+08 3.6E+06	1.2E+08 3.6E+06				
		7.0E-02 H 1.5E-01 I 2.5E-02 I		V		5.00E+02 1.36E+09 1.36E+09	1.28E+04 1.36E+09 1.36E+09		1	0.1	Methylstyrene, Alpha- Metolachlor Metribuzin	98-83-9 51218-45-2 21087-64-9					8.2E+04 1.8E+05 2.9E+04	4.1E+05 6.9E+04	1.2E+05 2.1E+04	8.2E+04 1.2E+05 2.1E+04			
1.8E+01 C	5.1E-03 C	2.5E-01 P 3.0E+00 P 2.0E-04 I		V		1.36E+09 1.36E+09 1.36E+09	8.58E+05 1.38E+03		1	0.1	Metsulfuron-methyl Mineral oils Mirex	74223-84-6 8012-95-1 2385-85-5	1.8E-01	2.1E+00	1.7E-01		2.9E+05 3.5E+06 2.3E+02	6.9E+05	2.1E+05 3.5E+06 2.3E+02	2.1E+05 3.5E+06 2.3E+02			
		2.0E-03 I 5.0E-03 I 1.0E-01 I				1.36E+09 1.36E+09 1.36E+09			1	0.1	Molinate Molybdenum Monochloramine	2212-67-1 7439-98-7 10599-90-3					2.3E+03 5.8E+03 1.2E+05	5.5E+03	1.6E+03 5.8E+03 1.2E+05	1.6E+03 5.8E+03 1.2E+05			
		2.0E-03 P 2.5E-02 I 3.0E-04 X				1.36E+09 1.36E+09 1.36E+09		0.04	1	0.1	Monomethylaniline Myclobutanil N,N'-Diphenyl-1,4-benzenediamine	100-61-8 88671-89-0 74-31-7					2.3E+03 2.9E+04 3.5E+02	5.5E+03 6.9E+04 8.3E+02	1.6E+03 2.1E+04 2.5E+02	1.6E+03 2.1E+04 2.5E+02			
1.8E+00 C	0.0E+00 C	2.0E-03 I 3.0E-02 X	1.0E-01 P V			1.36E+09 1.36E+09 1.36E+09	5.70E+04 1.36E+09		1	0.1	Naled Naphtha, High Flash Aromatic (HFAN) Naphthylamine, 2-	300-76-5 64742-95-6 91-59-8	1.8E+00	4.3E+00	1.3E+00		3.5E+04	6.0E+08	2.3E+03 3.5E+04	2.3E+03 3.5E+04			
		2.6E-04 C 2.6E-04 C	1.2E-01 O 1.1E-02 C 1.4E-05 C			1.36E+09 1.36E+09 1.36E+09			1	0.1	Napropamide Nickel Acetate Nickel Carbonate	15299-99-7 373-02-4 3333-67-3			6.4E+04 6.4E+04	6.4E+04 6.4E+04	1.4E+05 1.3E+04 1.3E+04	3.3E+05 3.0E+04 3.0E+04	8.3E+04 8.3E+04 8.3E+04	9.8E+04 8.1E+03 8.1E+03			
		2.6E-04 C 2.6E-04 C 2.6E-04 C	1.1E-02 C 1.1E-02 C 1.1E-02 C	1.4E-05 C V		1.36E+09 1.36E+09 1.36E+09		0.04	1	0.1	Nickel Carbonyl Nickel Hydroxide Nickel Oxide	13463-39-3 12054-48-7 1313-99-1			6.4E+04 6.4E+04 6.4E+04	6.4E+04 6.4E+04 6.4E+04	1.3E+04 1.3E+04 1.3E+04	3.0E+04 8.3E+04 1.2E+05	8.3E+04 8.3E+04 1.2E+04	1.1E+04 1.1E+04 1.2E+04			
1.7E+00 C	4.8E-04 C	2.4E-04 I 2.6E-04 C	1.1E-02 C 2.0E-02 I 1.1E-02 C	1.4E-05 C V		1.36E+09 1.36E+09 1.36E+09		0.04	1	0.1	Nickel Refinery Dust Nickel Soluble Salts Nickel Sulfide	E715532 7440-02-0 12035-72-2	1.9E+00		6.9E+04 6.4E+04 3.5E+04	6.9E+04 6.4E+04 1.9E+00	1.3E+04 2.3E+04 1.3E+04	3.0E+04 5.4E+05 8.3E+04	8.3E+04 1.1E+04 1.1E+04	1.1E+04 2.2E+04 1.1E+04			
		2.6E-04 C 1.6E+00 I	1.1E-02 C 1.4E-05 C			1.36E+09 1.36E+09 1.36E+09			1	0.1	Nickelocene Nitrate Nitrate + Nitrite (as N)	1271-28-9 14797-55-8 E701177			6.4E+04 6.4E+04	6.4E+04	1.3E+04 1.9E+06	3.0E+04 8.3E+04	8.3E+04 8.1E+03	8.1E+03 1.9E+06			
2.0E-02 P		1.0E-01 I 1.0E-02 X 4.0E-03 P	5.0E-05 X 6.0E-03 P			1.36E+09 1.36E+09 1.36E+09			1	0.1	Nitrite Nitroaniline, 2- Nitroaniline, 4-	14797-65-0 88-74-4 100-01-6	1.6E+02	3.9E+02		1.1E+02		1.2E+05 1.2E+04 4.7E+03	2.8E+04 2.8E+04 1.1E+04	3.0E+05 3.0E+05 3.6E+07	1.2E+05 8.0E+03 3.3E+03		
		4.0E-05 I	2.0E-03 I 3.0E+03 P 7.0E-02 H	9.0E-03 I V		3.05E+03 1.36E+09 1.36E+09	7.32E+04 1.36E+09		1	0.1	Nitrobenzene Nitrocellulose Nitrofurantoin	98-95-3 9004-70-0 67-20-9			2.2E+01	2.2E+01	2.3E+03 3.5E+09 8.2E+04	2.8E+02 8.3E+09 1.9E+05	2.9E+03 2.5E+09 5.7E+04	1.3E+03 2.5E+09 5.7E+04			
1.3E+00 C 1.7E-02 P	3.7E-04 C	1.0E-04 P 1.0E-01 I				1.36E+09 1.36E+09 1.36E+09			1	0.1	Nitrofurazone Nitroglycerin Nitroguanidine	59-87-0 55-63-0 556-88-7	2.5E+00 1.9E+02	5.9E+00 4.5E+02	4.5E+04 1.4E+02	1.8E+00		1.2E+02 2.8E+02	2.8E+05	8.2E+01 8.2E+04			
		8.8E-06 P 2.7E-03 H 7.7E-03 C	5.0E-03 P 2.0E-02 I	V		1.80E+04 4.86E+03 1.36E+09	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			2.4E+01 6.0E-02 2.2E+03	2.4E+01 6.0E-02 8.5E-02			3.7E+02 1.2E+03	3.7E+02 1.2E+03			
2.7E+01 C 1.2E+02 C 5.4E+00 I 7.0E+00 I	7.7E-03 C 3.4E-02 C 1.6E-03 I 2.0E-03 C				M	1.36E+09 1.36E+09 1.36E+09			1	0.1	Nitroso-N-methylurea, N- Nitroso-di-N-butylamine, N- Nitroso-di-N-propylamine, N-	684-93-5 924-16-3 621-64-7	1.2E-01 6.1E-01 4.7E-01	2.9E-01 6.4E-02 1.1E+00	2.2E+03 4.9E+02 8.3E+03	8.5E-02 1.9E+00 3.3E-01							
2.8E+00 I 1.5E+02 I 5.1E+01 I	8.0E-04 P 4.3E-02 I 1.4E-02 I					1.36E+09 1.36E+09 1.36E+09			1	0.1	Nitrosodiethanolamine, N- Nitrosodiethylamine, N- Nitrosodimethylamine, N-	1116-54-7 55-18-5 62-75-9	1.2E+00 2.2E-02 6.4E-02	2.8E+00 5.2E-02 7.2E-02	2.1E+04 3.9E+02 3.4E-02	8.2E-01 1.5E-02 3.4E-02	9.3E+00	1.4E+01	5.7E+00				
4.9E-03 I 2.2E+01 I 6.7E+00 C	2.6E-06 C 6.3E-03 C 1.9E-03 C					1.36E+09 1.36E+09 1.36E+09			1	0.1	Nitrosodiphenylamine, N- Nitrosomethyl ethylamine, N- Nitrosomorpholine [N-]	86-30-6 10595-95-6 59-89-2	6.7E+02 1.5E-01 4.9E-01	1.6E+03 2.4E-01 1.2E+00	6.4E+06 9.1E-02 8.8E+03	4.7E+02 9.1E-02 3.4E-01							
9.4E+00 C 2.1E+00 I	2.7E-03 C 6.1E-04 I					1.36E+09 1.36E+09			1	0.1	Nitrosopiperidine [N-] Nitrosopyrrolidine, N-	100-75-4 930-55-2	3.5E-01 1.6E+00	8.2E-01 3.7E+00	6.2E+03 2.7E+04	2.4E-01 1.1E+00							

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (see FO #31); H = HEAST; F = See FAQ; W = see user guide Section 2.3.5; E = see user guide Section 2.3.6; L = see user's guide Section 5.2; M = mutagen; S = see user's guide Section 5; V = volatile; R = RBA applied (see user's guide Section 5.10); c = cancer; n = noncancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; SSL values are based on DAF=1; m = Concentration may exceed ceiling limit (see user's guide Section 5.13); s = concentration may exceed Csat (see user's guide Section 5.12)

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer		Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -y)	k _e (y)	RFD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³ -y)	k _e (y)	muta gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	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 THI=1 (mg/kg)	
2.2E-01	P			1.0E-04	X					1.36E+09		1	0.1	Nitrotoluene, m-	99-08-1					1.2E+02	2.8E+02		8.2E+01	
9.0E-04	P			9.0E-04	P		V		1.51E+03	1.36E+09	1.37E+05	1		Nitrotoluene, o-	88-72-2	1.5E+01			1.5E+01	1.1E+03			1.1E+03	
1.6E-02	P			4.0E-03	P				1.36E+09	1.36E+09		1	0.1	Nitrotoluene, p-	99-99-0	2.0E+02	4.8E+02		1.4E+02	4.7E+03	1.1E+04		3.3E+03	
				3.0E-04	X	2.0E-02	P V		6.86E+00	1.36E+09	1.04E+03	1		Nonane, n-	111-84-2					3.5E+02		9.1E+01	7.2E+01	
				1.5E-02	O					1.36E+09		1	0.1	Norflurazon	27314-13-2					1.8E+04	4.1E+04		1.2E+04	
				3.0E-03	I					1.36E+09		1	0.1	Octabromodiphenyl Ether	32536-52-0					3.5E+03	8.3E+03		2.5E+03	
				5.0E-02	I					1.36E+09		1	0.006	Octahydro-1,3,5,7-tetraazocine (HMX)	2691-41-0					5.8E+04	2.3E+06		5.7E+04	
				2.0E-03	H					1.36E+09		1	0.1	Octamethylpyrophosphoramide	152-16-9				2.9E+02	2.3E+03	5.5E+03		1.6E+03	
7.8E-03	O			1.4E-01	O					1.36E+09		1	0.1	Oryzalin	19044-88-3	4.2E+02	9.9E+02		2.9E+02	1.6E+05	3.9E+05		1.1E+05	
				5.0E-03	I					1.36E+09		1	0.1	Oxadiazon	19666-30-9					5.8E+03	1.4E+04		4.1E+03	
				2.5E-02	I					1.36E+09		1	0.1	Oxamyl	23135-22-0					2.9E+04	6.9E+04		2.1E+04	
7.3E-02	O			3.0E-02	O					1.36E+09		1	0.1	Oxyfluorfen	42874-03-3	4.5E+01	1.1E+02		3.1E+01	3.5E+04	8.3E+04		2.5E+04	
				1.3E-02	I					1.36E+09		1	0.1	Paclitaxel	76738-62-0					1.5E+04	3.6E+04		1.1E+04	
				4.5E-03	I					1.36E+09		1	0.1	Paraquat Dichloride	1910-42-5					5.3E+03	1.2E+04		3.7E+03	
				6.0E-03	H					1.36E+09		1	0.1	Parathion	56-38-2					7.0E+03	1.7E+04		4.9E+03	
				5.0E-02	H		V			1.36E+09	4.49E+04	1		Pebutate	1114-71-2					5.8E+04			5.8E+04	
				3.0E-01	O					1.36E+09		1	0.1	Pendimethalin	40487-42-1					3.5E+05	8.3E+05		2.5E+05	
				2.0E-03	I		V		3.12E-01	1.36E+09	5.13E+05	1		Pentabromodiphenyl Ether	32534-81-9					2.3E+03			2.3E+03	
				1.0E-04	I					1.36E+09		1	0.1	Pentabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-99)	60348-60-9					1.2E+02	2.8E+02		8.2E+01	
				8.0E-04	I		V			1.36E+09	8.12E+04	1		Pentachlorobenzene	608-93-5					9.3E+02			9.3E+02	
9.0E-02	P			3.0E-03	I		V		4.57E+02	1.36E+09	9.65E+03	1		Pentachloroethane	76-01-7	3.6E+01			3.6E+01	3.5E+03			3.5E+03	
2.6E-01	H			3.0E-03	I		V			1.36E+09	4.32E+05	1		Pentachloronitrobenzene	82-68-8	1.3E+01			1.3E+01	5.5E+03			5.5E+03	
4.0E-01	I	5.1E-06	C	5.0E-03	I					1.36E+09		1	0.25	Pentachlorophenol	87-86-5	8.2E+00	7.7E+00	3.3E+06	4.0E+00	5.8E+03	5.5E+03		2.8E+03	
4.0E-03	X			2.0E-03	P					1.36E+09		1	0.1	Pentaerythritol tetranitrate (PETN)	78-11-5	8.2E+02	1.9E+03		5.7E+02	2.3E+03	5.5E+03		1.6E+03	
				1.0E+00	P V				3.88E+02	1.36E+09	7.79E+02	1		Pentane, n-	109-66-0							3.4E+03	3.4E+03	
				7.0E-04	I					1.36E+09		1		Perchlorates									8.2E+02	
				7.0E-04	I					1.36E+09		1		~Ammonium Perchlorate	7790-98-9								8.2E+02	
				7.0E-04	I					1.36E+09		1		~Lithium Perchlorate	7791-03-9								8.2E+02	
				7.0E-04	I					1.36E+09		1		~Perchlorate and Perchlorate Salts	14797-73-0								8.2E+02	
				7.0E-04	I					1.36E+09		1		~Potassium Perchlorate	7778-74-7								8.2E+02	
				7.0E-04	I					1.36E+09		1		~Sodium Perchlorate	7601-89-0								8.2E+02	
				2.0E-02	P					1.36E+09		1	0.1	Perfluorobutane sulfonic acid (PFBS)	375-73-5					2.3E+04	5.5E+04		1.6E+04	
				2.0E-02	P					1.36E+09		1	0.1	Perfluorobutanesulfonate	45187-15-3					2.3E+04	5.5E+04		1.6E+04	
				5.0E-02	I					1.36E+09		1	0.1	Permethrin	52645-53-1					5.8E+04	1.4E+05		4.1E+04	
2.2E-03	C	6.3E-07	C							1.36E+09		1	0.1	Phenacetin	62-44-2	1.5E+03	3.5E+03	2.6E+07	1.0E+03					
				2.4E-01	O					1.36E+09		1	0.1	Phenmedipham	13684-63-4					2.8E+05	6.6E+05		2.0E+05	
				3.0E-01	I	2.0E-01	C			1.36E+09		1	0.1	Phenol	108-95-2					3.5E+05	8.3E+05	1.2E+09	2.5E+05	
				4.0E-03	I					1.36E+09		1	0.1	Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1					4.7E+03	1.1E+04		3.3E+03	
				5.0E-04	X					1.36E+09		1	0.1	Phenothiazine	92-84-2					5.8E+02	1.4E+03		4.1E+02	
				2.0E-04	X		V		1.29E+02	1.36E+09	7.06E+03	1		Phenyl Isothiocyanate	103-72-0					2.3E+02			2.3E+02	
				6.0E-03	I					1.36E+09		1	0.1	Phenylethylamine, m-	108-45-2					7.0E+03	1.7E+04		4.9E+03	
1.2E-01	P			4.0E-03	P					1.36E+09		1	0.1	Phenylethylamine, o-	95-54-5	2.7E+01	6.4E+01		1.9E+01	4.7E+03	1.1E+04		3.3E+03	
				1.0E-03	X					1.36E+09		1	0.1	Phenylethylamine, p-	106-50-3					1.2E+03	2.8E+03		8.2E+02	
1.9E-03	H			2.0E-04	H					1.36E+09		1	0.1	Phenylphenol, 2-	90-43-7	1.7E+03	4.0E+03		1.2E+03				1.6E+02	
						3.0E-04	I V		1.61E+03	1.36E+09	9.81E+02	1		Phorate	298-02-2					2.3E+02	5.5E+02		1.3E+00	
				2.0E-02	I					1.36E+09		1	0.1	Phosgene	75-44-5							1.3E+00	1.6E+04	
				4.9E+01	P					1.36E+09		1		Phosmet	732-11-6					2.3E+04	5.5E+04		1.6E+04	
				4.9E+01	P					1.36E+09		1		Phosphates, Inorganic									5.7E+07	
				4.9E+01	P					1.36E+09		1		~Aluminum metaphosphate	13776-88-0					5.7E+07			5.7E+07	
				4.9E+01	P					1.36E+09		1		~Ammonium polyphosphate	68333-79-9					5.7E+07			5.7E+07	
				4.9E+01	P					1.36E+09		1		~Calcium pyrophosphate	7790-76-3					5.7E+07			5.7E+07	
				4.9E+01	P					1.36E+09		1		~Diammonium phosphate	7783-28-0					5.7E+07			5.7E+07	
				4.9E+01	P					1.36E+09		1		~Dicalcium phosphate	7757-93-9					5.7E+07			5.7E+07	
				4.9E+01	P					1.36E+09		1		~Dimagnesium phosphate	7782-75-4					5.7E+07			5.7E+07	
				4.9E+01	P					1.36E+09		1		~Dipotassium phosphate	7758-11-4					5.7E+07			5.7E+07	
				4.9E+01	P					1.36E+09		1		~Disodium phosphate	7558-79-4					5.7E+07			5.7E+07	
				4.9E+01	P					1.36E+09		1		~Monoaluminum phosphate	13530-50-2					5.7E+07			5.7E+07	
				4.9E+01	P					1.36E+09		1		~Monoammonium phosphate	7722-76-1					5.7E+07			5.7E+07	
				4.9E+01	P					1.36E+09		1		~Monocalcium phosphate	7758-23-8					5.7E+07			5.7E+07	
				4.9E+01	P																			

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (see FO #31); H = HEAST; F = See FAQ; W = see user guide Section 2.3.5; E = see user guide Section 2.3.6; L = see user's guide Section 5.2; M = mutagen; S = see user's guide Section 5; V = volatile; R = RBA applied (see user's guide Section 5.10); c = cancer; n = noncancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; SSL values are based on DAF=1; m = Concentration may exceed ceiling limit (see user's guide Section 5.13); s = concentration may exceed Csat (see user's guide Section 5.12)														Contaminant									
Toxicity and Chemical-specific Information														Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -y)	k _e (y)	RFD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³ -y)	k _e (y)	muta	C _{sat}	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	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 TH=1 (mg/kg)
				4.9E+01	P					1.36E+09		1		~Trialuminum sodium tetra decahydrogenoctaorthophosphate (dihydrate)	15136-87-5					5.7E+07			5.7E+07
				4.9E+01	P					1.36E+09		1		~Tricalcium phosphate	7758-87-4					5.7E+07			5.7E+07
				4.9E+01	P					1.36E+09		1		~Trimagnesium phosphate	7757-87-1					5.7E+07			5.7E+07
				4.9E+01	P					1.36E+09		1		~Tripotassium phosphate	7778-53-2					5.7E+07			5.7E+07
				4.9E+01	P					1.36E+09		1		~Trisodium phosphate	7601-54-9					5.7E+07			5.7E+07
				3.0E-04	I	3.0E-04	I	V		1.36E+09		1		Phosphine	7803-51-2					3.5E+02		1.8E+06	3.5E+02
				4.9E+01	P	1.0E-02	I			1.36E+09		1		Phosphoric Acid	7664-38-2					5.7E+07		6.0E+07	2.9E+07
				2.0E-05	I			V		1.36E+09	6.92E+03	1		Phosphorus, White	7723-14-0					2.3E+01			2.3E+01
1.4E-02	I	2.4E-06	C	2.0E-02	I					1.36E+09		1	0.1	Phthalates						2.3E+04	5.5E+04		1.6E+04
1.9E-03	P			2.0E-01	I					1.36E+09		1	0.1	~Bis(2-ethylhexyl)phthalate	117-81-7	2.3E+02	5.5E+02	6.9E+06	1.6E+02	2.3E+05	5.5E+05		1.6E+05
				1.0E+00	I					1.36E+09		1	0.1	~Butyl Benzyl Phthalate	85-68-7	1.7E+03	4.1E+03		1.2E+03	2.3E+05	2.8E+05		1.6E+05
				1.0E-01	I					1.36E+09		1	0.1	~Butylphthalyl Butylglycolate	85-70-1					1.2E+06	2.8E+06		8.2E+05
				1.0E-01	I					1.36E+09		1	0.1	~Dibutyl Phthalate	84-74-2					1.2E+05	2.8E+05		8.2E+04
				8.0E-01	I					1.36E+09		1	0.1	~Diethyl Phthalate	84-66-2					9.3E+05	2.2E+06		6.6E+05
				1.0E-01	I			V		1.36E+09	2.13E+04	1		~Dimethylterephthalate	120-61-6					1.2E+05			1.2E+05
				1.0E-02	P					1.36E+09		1	0.1	~Octyl Phthalate, di-N-	117-84-0					1.2E+04	2.8E+04		8.2E+03
				1.0E+00	H					1.36E+09		1	0.1	~Phthalic Acid, P-	100-21-0					1.2E+06	2.8E+06		8.2E+05
				2.0E+00	I	2.0E-02	C			1.36E+09		1	0.1	~Phthalic Anhydride	85-44-9					2.3E+06	5.5E+06	1.2E+08	1.6E+06
				7.0E-02	I					1.36E+09		1	0.1	Picloram	1918-02-1					8.2E+04	1.9E+05		5.7E+04
				1.0E-04	X					1.36E+09		1	0.1	Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3					1.2E+02	2.8E+02		8.2E+01
				9.0E-04	X					1.36E+09		1	0.1	Picric Acid (2,4,6-Trinitrophenol)	88-89-1					1.1E+03	2.5E+03		7.4E+02
				7.0E-05	O					1.36E+09		1	0.1	Pinimiphos, Methyl	29232-93-7					8.2E+01	1.9E+02		5.7E+01
3.0E+01	C	8.6E-03	C	7.0E-06	H					1.36E+09		1	0.1	Polybrominated Biphenyls	59536-65-1	1.1E-01	2.6E-01	1.9E+03	7.7E-02	8.2E+00	1.9E+01		5.7E+00
				7.0E-02	S	2.0E-05	S	7.0E-05	I		V		0.14	Polychlorinated Biphenyls (PCBs)									
				2.0E+00	S	5.7E-04	S			1.36E+09	7.14E+05	1	0.14	~Aroclor 1016	12674-11-2	4.7E+01	7.9E+01	4.4E+02	2.7E+01	8.2E+01	1.4E+02		5.1E+01
				2.0E+00	S	5.7E-04	S			1.36E+09	2.04E+05	1	0.14	~Aroclor 1221	11104-28-2	1.6E+00	2.8E+00	4.4E+00	8.3E-01				
				2.0E+00	S	5.7E-04	S			1.36E+09	1.12E+05	1	0.14	~Aroclor 1232	11141-16-5	1.6E+00	2.8E+00	2.4E+00	7.2E-01				
				2.0E+00	S	5.7E-04	S			1.36E+09	5.91E+05	1	0.14	~Aroclor 1242	53469-21-9	1.6E+00	2.8E+00	1.3E+01	9.5E-01				
				2.0E+00	S	5.7E-04	S			1.36E+09	6.25E+05	1	0.14	~Aroclor 1248	12672-29-6	1.6E+00	2.8E+00	1.3E+01	9.5E-01				
				2.0E+00	S	5.7E-04	S	2.0E-05	I		V		0.14	~Aroclor 1254	11097-69-1	1.6E+00	2.8E+00	1.8E+01	9.7E-01	2.3E+01	3.9E+01		1.5E+01
				2.0E+00	S	5.7E-04	S			1.36E+09	1.31E+06	1	0.14	~Aroclor 1260	11096-82-5	1.6E+00	2.8E+00	2.8E+01	9.9E-01				
								6.0E-04	X		V		0.14	~Aroclor 5460	11126-42-4					7.0E+02	1.2E+03		4.4E+02
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	WV			1.36E+09	2.43E+06	1	0.14	~Heptachlorobiphenyl, 2,3,3',4,4',5,5'-(PCB 189)	39635-31-9	8.4E-01	1.4E+00	2.6E+01	5.2E-01	2.7E+01	4.6E+01	1.4E+04	1.7E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	WV			1.36E+09	1.58E+06	1	0.14	~Hexachlorobiphenyl, 2,3',4,4',5,5'-(PCB 167)	52663-72-6	8.4E-01	1.4E+00	1.7E+01	5.1E-01	2.7E+01	4.6E+01	9.2E+03	1.7E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	WV			1.36E+09	1.04E+06	1	0.14	~Hexachlorobiphenyl, 2,3,3',4,4',5,5'-(PCB 157)	69782-90-7	8.4E-01	1.4E+00	1.1E+01	5.0E-01	2.7E+01	4.6E+01	6.1E+03	1.7E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	WV			1.36E+09	1.11E+06	1	0.14	~Hexachlorobiphenyl, 2,3,3',4,4',5,5'-(PCB 156)	38380-08-4	8.4E-01	1.4E+00	1.2E+01	5.0E-01	2.7E+01	4.6E+01	6.5E+03	1.7E+01
3.9E+03	W	1.1E+00	W	2.3E-08	W	1.3E-06	WV			1.36E+09	1.58E+06	1	0.14	~Hexachlorobiphenyl, 3,3',4,4',5,5'-(PCB 169)	32774-16-6	8.4E-04	1.4E-03	1.7E-02	5.1E-04	2.7E-02	4.6E-02	9.2E+00	1.7E-02
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	WV			1.36E+09	7.33E+05	1	0.14	~Pentachlorobiphenyl, 2',3,4,4',5,5'-(PCB 123)	65510-44-3	8.4E-01	1.4E+00	7.9E+00	4.9E-01	2.7E+01	4.6E+01	4.3E+03	1.7E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	WV			1.36E+09	5.90E+05	1	0.14	~Pentachlorobiphenyl, 2,3',4,4',5,5'-(PCB 118)	31508-00-6	8.4E-01	1.4E+00	6.3E+00	4.9E-01	2.7E+01	4.6E+01	3.4E+03	1.7E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	WV			1.36E+09	6.01E+05	1	0.14	~Pentachlorobiphenyl, 2,3,3',4,4'-(PCB 105)	32598-14-4	8.4E-01	1.4E+00	6.5E+00	4.9E-01	2.7E+01	4.6E+01	3.5E+03	1.7E+01
3.9E+00	W	1.1E-03	W	2.3E-05	W	1.3E-03	WV			1.36E+09	1.05E+06	1	0.14	~Pentachlorobiphenyl, 2,3,4,4',5,5'-(PCB 114)	74472-37-0	8.4E-01	1.4E+00	1.1E+01	5.0E-01	2.7E+01	4.6E+01	6.1E+03	1.7E+01
1.3E+04	W	3.8E+00	W	7.0E-09	W	4.0E-07	WV			1.36E+09	7.26E+05	1	0.14	~Pentachlorobiphenyl, 3,3',4,4',5,5'-(PCB 126)	57465-28-8	2.5E-04	4.2E-04	2.3E-03	1.5E-04	8.2E-03	1.4E-02	1.3E+00	5.1E-03
2.0E+00	I	5.7E-04	I		V					1.36E+09	5.32E+05	1	0.14	~Polychlorinated Biphenyls (high risk)	1336-36-3	1.6E+00	2.8E+00	1.1E+01	9.4E-01				
4.0E-01	I	1.0E-04	I		V					1.36E+09		1	0.14	~Polychlorinated Biphenyls (low risk)	1336-36-3								
7.0E-02	I	2.0E-05	I		V					1.36E+09		1	0.14	~Polychlorinated Biphenyls (lowest risk)	1336-36-3								
1.3E+01	W	3.8E-03	W	7.0E-06	W	4.0E-04	WV			1.36E+09		1	0.14	~Tetrachlorobiphenyl, 3,3',4,4'-(PCB 77)	32598-13-3	2.5E-01	4.2E-01	4.4E+03	1.6E-01	8.2E+00	1.4E+01	2.4E+06	5.1E+00
3.9E+01	W	1.1E-02	W	2.3E-06	W	1.3E-04	WV			1.36E+09	5.09E+05	1	0.14	~Tetrachlorobiphenyl, 3,4,4',5'-(PCB 81)	70362-50-4	8.4E-02	1.4E-01	5.5E-01	4.8E-02	2.7E+00	4.6E+00	3.0E+02	1.7E+00
				6.0E-04	I					1.36E+09		1	0.1	Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9							3.6E+06	3.6E+06
				6.0E-02	I		V			1.36E+09	1.41E+05	1	0.13	Polynuclear Aromatic Hydrocarbons (PAHs)						7.0E+04	1.3E+05		4.5E+04
				3.0E-01	I		V			1.36E+09	5.23E+05	1	0.13	~Anthracene	120-12-7					3.5E+05	6.4E+05		2.3E+05
1.0E-01	E	6.0E-05	E		V		M			1.36E+09	4.41E+06	1	0.13	~Benz[a]anthracene	56-55-3	3.3E+01	5.9E+01	9.0E+02	2.1E+01				
1.2E+00	C	1.1E-04	C		V		M			1.36E+09		1	0.13	~Ben									

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (see FO #31); H = HEAST; F = See FAQ; W = see user guide Section 2.3.5; E = see user guide Section 2.3.6; L = see user's guide Section 5.2; M = mutagen; S = see user's guide Section 5; V = volatile; R = RBA applied (see user's guide Section 5.10); c = cancer; n = noncancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; SSL values are based on DAF=1; n = Concentration may exceed ceiling limit (see user's guide Section 5.13); s = concentration may exceed Csat (see user's guide Section 5.12)														Contaminant										
Toxicity and Chemical-specific Information														Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -y) ⁻¹	k _e (y)	RFD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³)	k _e (y)	muta	gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	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 THI=1 (mg/kg)
				7.5E-02	I						1.36E+09		1	0.1	Pronamide	23950-58-5					8.8E+04	2.1E+05		6.2E+04
				1.3E-02	I						1.36E+09		1	0.1	Propachlor	1918-16-7					1.5E+04	3.6E+04		1.1E+04
				5.0E-03	I						1.36E+09		1	0.1	Propamil	709-98-8					5.8E+03	1.4E+04		4.1E+03
1.9E-01	O			4.0E-02	O						1.36E+09		1	0.1	Propargite	2312-35-8	1.7E+01	4.0E+01		1.2E+01	4.7E+04	1.1E+05		3.3E+04
				2.0E-03	I		V			1.11E+05	1.36E+09	6.27E+04	1	1	Propargyl Alcohol	107-19-7					2.3E+03			2.3E+03
				2.0E-02	I						1.36E+09		1	0.1	Propazine	139-40-2					2.3E+04	5.5E+04		1.6E+04
				2.0E-02	I						1.36E+09		1	0.1	Propham	122-42-9					2.3E+04	5.5E+04		1.6E+04
				1.0E-01	O						1.36E+09		1	0.1	Propiconazole	60207-90-1					1.2E+05	2.8E+05		8.2E+04
						8.0E-03	I	V		3.26E+04	1.36E+09	8.94E+03	1	1	Propionaldehyde	123-38-6							3.1E+02	3.1E+02
				1.0E-01	X	1.0E+00	X	V		2.64E+02	1.36E+09	6.99E+03	1	1	Propyl benzene	103-65-1					1.2E+05			2.4E+04
						3.0E+00	C	V		3.49E+02	1.36E+09	7.04E+02	1	1	Propylene	115-07-1							9.3E+03	9.3E+03
				2.0E+01	P						1.36E+09		1	0.1	Propylene Glycol	57-55-6					2.3E+07	5.5E+07		1.6E+07
						2.7E-04	A				1.36E+09		1	0.1	Propylene Glycol Dinitrate	6423-43-4							1.6E+06	1.6E+06
2.4E-01	I	3.7E-06	I	7.0E-01	H	2.0E+00	I	V		1.06E+05	1.36E+09	7.83E+04	1	1	Propylene Glycol Monomethyl Ether	107-98-2					8.2E+05			3.7E+05
						3.0E-02	I	V		7.77E+04	1.36E+09	1.03E+04	1	1	Propylene Oxide	75-56-9	1.4E+01		3.4E+01	9.7E+00			1.4E+03	1.4E+03
				1.0E-03	I		V			5.30E+05	1.36E+09	5.54E+04	1	1	Pyridine	110-86-1					1.2E+03			1.2E+03
3.0E+00	I			5.0E-04	I						1.36E+09		1	0.1	Quinalphos	13593-03-8	1.1E+00	2.6E+00		7.7E-01	5.8E+02	1.4E+03		4.1E+02
											1.36E+09		1	0.1	Quinoline	91-22-5								4.1E+02
				9.0E-03	I						1.36E+09		1	0.1	Quizalofop-ethyl	76578-14-8					1.1E+04	2.5E+04		7.4E+03
						3.0E-02	A				1.36E+09		1	0.1	Refractory Ceramic Fibers (units in fibers)	E715557								
				3.0E-02	I						1.36E+09		1	0.1	Resmethrin	10453-86-8					3.5E+04	8.3E+04		2.5E+04
				5.0E-02	H		V				1.36E+09	4.65E+05	1	1	Ronnel	299-84-3	1.5E+01	3.5E+01	2.6E+05	1.0E+01	5.8E+04			5.8E+04
2.2E-01	C	6.3E-05	C	4.0E-03	I						1.36E+09		1	0.1	Rotenone	83-79-4					4.7E+03	1.1E+04		3.3E+03
											1.36E+09		1	0.1	Safrole	94-59-7								3.3E+03
				5.0E-03	I						1.36E+09		1	0.1	Selenious Acid	7783-00-8					5.8E+03			5.8E+03
				5.0E-03	I	2.0E-02	C				1.36E+09		1	0.1	Selenium	7782-49-2					5.8E+03		1.2E+08	5.8E+03
				5.0E-03	C	2.0E-02	C				1.36E+09		1	0.1	Selenium Sulfide	7446-34-6					5.8E+03		1.2E+08	5.8E+03
				1.4E-01	O						1.36E+09		1	0.1	Sethoxydim	74051-80-2					1.6E+05	3.9E+05		1.1E+05
						3.0E-03	C				1.36E+09		1	0.1	Silica (crystalline, respirable)	7631-86-9							1.8E+07	1.8E+07
				5.0E-03	I						1.36E+09		0.04	Silver	7440-22-4									5.8E+03
1.2E-01	H			5.0E-03	I						1.36E+09		1	0.1	Simazine	122-34-9	2.7E+01	6.4E+01		1.9E+01	5.8E+03	1.4E+04		4.1E+03
				1.3E-02	I						1.36E+09		1	0.1	Sodium Acifluorfen	62476-59-9					1.5E+04	3.6E+04		1.1E+04
				4.0E-03	I						1.36E+09		1	0.1	Sodium Azide	26628-22-8					4.7E+03			4.7E+03
2.7E-01	H			3.0E-02	I						1.36E+09		1	0.1	Sodium Diethyldithiocarbamate	148-18-5	1.2E+01	2.9E+01		8.5E+00	3.5E+04	8.3E+04		2.5E+04
				5.0E-02	A	1.3E-02	C				1.36E+09		1	0.1	Sodium Fluoride	7681-49-4					5.8E+04		7.7E+07	5.8E+04
				2.0E-05	I						1.36E+09		1	0.1	Sodium Fluoroacetate	62-74-8					2.3E+01	5.5E+01		1.6E+01
				1.0E-03	H						1.36E+09		1	0.1	Sodium Metavanadate	13718-26-8					1.2E+03			1.2E+03
				8.0E-04	P						1.36E+09		1	0.1	Sodium Tungstate	13472-45-2					9.3E+02			9.3E+02
				8.0E-04	A						1.36E+09		1	0.1	Sodium Tungstate Dihydrate	10213-10-2					9.3E+02			9.3E+02
2.4E-02	H			3.0E-02	I						1.36E+09		1	0.1	Strofos (Tetrachlorovinphos)	961-11-5	1.4E+02	3.2E+02		9.6E+01	3.5E+04	8.3E+04		2.5E+04
				6.0E-01	I						1.36E+09		1	0.1	Strontium, Stable	7440-24-6					7.0E+05			7.0E+05
				3.0E-04	I						1.36E+09		1	0.1	Strychnine	57-24-9					3.5E+02	8.3E+02		2.5E+02
				2.0E-01	I	1.0E+00	I	V		8.67E+02	1.36E+09	9.35E+03	1	1	Styrene	100-42-5					2.3E+05		4.1E+04	3.5E+04
				3.0E-03	P						1.36E+09		1	0.1	Styrene-Acrylonitrile (SAN) Trimer	57964-39-3					3.5E+03	8.3E+03		2.5E+03
				1.0E-03	P	2.0E-03	X				1.36E+09		1	0.1	Sulfolane	126-33-0					1.2E+03	2.8E+03	1.2E+07	8.2E+02
				8.0E-04	P						1.36E+09		1	0.1	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9					9.3E+02	2.2E+03		6.6E+02
						1.0E-03	C	V			1.36E+09		1	0.1	Sulfur Trioxide	7446-11-9							6.0E+06	6.0E+06
						1.0E-03	C				1.36E+09		1	0.1	Sulfuric Acid	7664-93-9							6.0E+06	6.0E+06
2.5E-02	I	7.1E-06	I	5.0E-02	H						1.36E+09		1	0.1	Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl	140-57-8	1.3E+02	3.1E+02	2.3E+06	9.2E+01	5.8E+04	1.4E+05		4.1E+04
				3.0E-02	H						1.36E+09		1	0.1	TCMTB	21564-17-0					3.5E+04	8.3E+04		2.5E+04
				7.0E-02	I						1.36E+09		1	0.1	Tebuthiuron	34014-18-1					8.2E+04	1.9E+05		5.7E+04
				2.0E-02	H						1.36E+09		1	0.1	Temephos	3383-96-8					2.3E+04	5.5E+04		1.6E+04
				1.3E-02	I						1.36E+09		1	0.1	Terbacil	5902-51-2					1.5E+04	3.6E+04		1.1E+04
				2.5E-05	H		V			3.09E+01	1.36E+09	2.64E+05	1	1	Terbufos	13071-79-9					2.9E+01			2.9E+01
				1.0E-03	I						1.36E+09		1	0.1	Terbutryn	886-50-0					1.2E+03	2.8E+03		8.2E+02
5.0E-03	C	1.3E-06	C	3.0E-04	I		V				1.36E+09	3.99E+03	1	0.1	Tert-Butyl Acetate	540-88-5	6.5E+02	1.5E+03	3.8E+01	3.5E+01				2.8E+02
				1.0E-04	I						1.36E+09		1	0.1	Tetrabromodiphenyl ether, 2,2',4,4'-(BDE-47)	5436-43-1					1.2E+02	2.8E+02		8.2E+01

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (see FO #31); H = HEAST; F = See FAQ; W = see user guide Section 2.3.5; E = see user guide Section 2.3.6; L = see user's guide Section 5.2; M = mutagen; S = see user's guide Section 5; V = volatile; R = RBA applied (see user's guide Section 5.10); c = cancer; n = noncancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; SSL values are based on DAF=1; m = Concentration may exceed ceiling limit (see user's guide Section 5.13); s = concentration may exceed Csat (see user's guide Section 5.12)

Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³ -y)	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³)	k _e (y)	muta gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	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 THI=1 (mg/kg)
				4.3E-02	O					1.36E+09		1	0.1	Thiufenulfuron-methyl	79277-27-3					5.0E+04	1.2E+05		3.5E+04
				1.0E-02	I					1.36E+09		1	0.1	Thiobencarb	28249-77-6					1.2E+04	2.8E+04		8.2E+03
				7.0E-02	X					1.36E+09		1	0.0075	Thiodiglycol	111-48-8					8.2E+04	2.6E+06		7.9E+04
				3.0E-04	H					1.36E+09		1	0.1	Thiofanox	39196-18-4					3.5E+02	8.3E+02		2.5E+02
1.2E-02	O			2.7E-02	O					1.36E+09		1	0.1	Thiophanate, Methyl	23564-05-8	2.8E+02	6.7E+02		2.0E+02	3.1E+04	7.4E+04		2.2E+04
				1.5E-02	O					1.36E+09		1	0.1	Thiram	137-26-8					1.8E+04	4.1E+04		1.2E+04
				6.0E-01	H					1.36E+09		1		Tin	7440-31-5					7.0E+05			7.0E+05
				8.0E-02	I	1.0E-04	A	V		1.36E+09		1		Titanium Tetrachloride	7550-45-0					9.3E+04		6.0E+05	6.0E+05
		1.1E-05	C			5.0E+00	I	V	8.18E+02	1.36E+09	4.29E+03	1		Toluene	108-88-3							9.4E+04	4.7E+04
						8.0E-06	C	V		1.36E+09	7.62E+05	1		Toluene-2,4-diisocyanate	584-84-9			8.5E+02	8.5E+02			2.7E+01	2.7E+01
1.8E-01	X			2.0E-04	X					1.36E+09		1	0.1	Toluene-2,5-diamine	95-70-5	1.8E+01	4.3E+01			2.3E+02	5.5E+02		1.6E+02
		1.1E-05	C			8.0E-06	C	V	1.71E+03	1.36E+09	6.32E+05	1		Toluene-2,6-diisocyanate	91-08-7			7.0E+02	7.0E+02			2.2E+01	2.2E+01
				5.0E-03	P					1.36E+09		1	0.1	Toluic Acid, p-	99-94-5					5.8E+03	1.4E+04		4.1E+03
1.6E-02	P	5.1E-05	C							1.36E+09		1	0.1	Toluidine, o- (Methylaniline, 2-)	95-53-4	2.0E+02	4.8E+02	3.3E+05	1.4E+02				
3.0E-02	P			4.0E-03	X					1.36E+09		1	0.1	Toluidine, p-	106-49-0	1.1E+02	2.6E+02		7.7E+01	4.7E+03	1.1E+04		3.3E+03
				3.0E+00	P			V	3.42E-01	1.36E+09	1.38E+03	1		Total Petroleum Hydrocarbons (Aliphatic High)	E1790670					3.5E+06			3.5E+06
						6.0E-01	P	V	1.41E+02	1.36E+09	8.29E+02	1		Total Petroleum Hydrocarbons (Aliphatic Low)	E1790666							2.2E+03	2.2E+03
				1.0E-02	X	1.0E-01	P	V	6.86E+00	1.36E+09	1.04E+03	1		Total Petroleum Hydrocarbons (Aliphatic Medium)	E1790668					1.2E+04	4.6E+02		4.4E+02
				4.0E-02	P					1.36E+09		1	0.1	Total Petroleum Hydrocarbons (Aromatic High)	E1790676					4.7E+04	1.1E+05		3.3E+04
				4.0E-03	P	3.0E-02	P	V	1.82E+03	1.36E+09	3.54E+03	1		Total Petroleum Hydrocarbons (Aromatic Low)	E1790672					4.7E+03		4.6E+02	4.2E+02
				4.0E-03	P	3.0E-03	P	V	1.36E+09	5.24E+04		1		Total Petroleum Hydrocarbons (Aromatic Medium)	E1790674					4.7E+03		6.9E+02	6.0E+02
1.1E+00	I	3.2E-04	I							1.36E+09		1	0.1	Toxaphene	8001-35-2	3.0E+00	7.0E+00	5.2E+04	2.1E+00	1.1E+02	2.5E+02		7.4E+01
				3.0E-05	X					1.36E+09		1	0.1	Toxaphene, Weathered	E1841606					3.5E+01	8.3E+01		2.5E+01
				7.5E-03	I					1.36E+09		1	0.1	Tralometrin	66841-25-6					8.8E+03	2.1E+04		6.2E+03
				3.0E-04	A			V		1.36E+09	3.36E+03	1		Tri-n-butyltin	688-73-3					3.5E+02			3.5E+02
				8.0E+01	X					1.36E+09		1	0.1	Triacetin	102-76-1					9.3E+07	2.2E+08		6.6E+07
				3.4E-02	O					1.36E+09		1	0.1	Triadimefon	43121-43-3					4.0E+04	9.4E+04		2.8E+04
7.2E-02	O			2.5E-02	O			V		1.36E+09	3.62E+05	1		Triallate	2303-17-5	4.6E+01			4.6E+01	2.9E+04			2.9E+04
				1.0E-02	I					1.36E+09		1	0.1	Trisulfuron	82097-50-5					1.2E+04	2.8E+04		8.2E+03
				8.0E-03	I					1.36E+09		1	0.1	Tribenuron-methyl	101200-48-0					9.3E+03	2.2E+04		6.6E+03
				5.0E-03	I			V		1.36E+09	4.83E+04	1		Tri Bromobenzene, 1,2,4-	615-54-3					5.8E+03			5.8E+03
9.0E-03	P			9.0E-03	X					1.36E+09		1	0.1	Tribromophenol, 2,4,6-	118-79-6	3.6E+02	8.6E+02		2.6E+02	1.1E+04	2.5E+04		7.4E+03
				1.0E-02	P					1.36E+09		1	0.1	Tributyl Phosphate	126-73-8					1.2E+04	2.8E+04		8.2E+03
				3.0E-04	P					1.36E+09		1	0.1	Tributyltin Compounds	E1790678					3.5E+02	8.3E+02		2.5E+02
				3.0E-04	I					1.36E+09		1	0.1	Tributyltin Oxide	56-35-9					3.5E+02	8.3E+02		2.5E+02
7.0E-02	I			3.0E+01	I	5.0E+00	P	V	9.10E+02	1.36E+09	1.29E+03	1	0.1	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1	4.7E+01	1.1E+02		3.3E+01	3.5E+07		2.8E+04	2.8E+04
2.9E-02	H			2.0E-02	I					1.36E+09		1	0.1	Trichloroacetic Acid	76-03-9	1.1E+02			3.3E+01	3.5E+07		2.8E+04	2.8E+04
7.0E-03	X			3.0E-05	X					1.36E+09		1	0.1	Trichloroaniline HCl, 2,4,6-	33663-50-2	1.1E+02	2.7E+02		7.9E+01	4.0E+04	9.4E+04		2.5E+01
				8.0E-04	X			V		1.36E+09	3.22E+04	1		Trichloroaniline, 2,4,6-	634-93-5	4.7E+02	1.1E+03		3.3E+02	3.5E+01	8.3E+01		2.5E+01
2.9E-02	P			1.0E-02	I	2.0E-03	P	V	4.04E+02	1.36E+09	2.99E+04	1		Trichlorobenzene, 1,2,4-	120-82-1	1.1E+02		1.1E+02	9.3E+02			2.6E+02	2.6E+02
				2.0E+00	I	5.0E+00	I	V	6.40E+02	1.36E+09	1.65E+03	1		Trichloroethane, 1,1,1-	71-55-6					2.3E+06		3.6E+04	3.6E+04
5.7E-02	I	1.6E-05	I	4.0E-03	I	2.0E-04	X	V	2.16E+03	1.36E+09	7.22E+03	1		Trichloroethane, 1,1,2-	79-00-5	5.7E+01		5.5E+00	5.0E+00	4.7E+03		6.3E+00	6.3E+00
4.6E-02	I	4.1E-06	I	5.0E-04	I	2.0E-03	I	V	6.92E+02	1.36E+09	2.21E+03	1		Trichloroethylene	79-01-6	7.1E+01	6.6E+00	6.0E+00		5.8E+02		1.9E+01	1.9E+01
				3.0E-01	I			V	1.23E+03	1.36E+09	1.04E+03	1		Trichlorofluoromethane	75-69-4					3.5E+05			3.5E+05
				1.0E-01	I					1.36E+09		1	0.1	Trichlorophenol, 2,4,5-	95-95-4					1.2E+05	2.8E+05		8.2E+04
1.1E-02	I	3.1E-06	I	1.0E-03	P					1.36E+09		1	0.1	Trichlorophenol, 2,4,6-	88-06-2	3.0E+02	7.0E+02	5.4E+06	2.1E+02	1.2E+03	2.8E+03		8.2E+02
				1.0E-02	I					1.36E+09		1	0.1	Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5					1.2E+04	2.8E+04		8.2E+03
				8.0E-03	I					1.36E+09		1	0.1	Trichlorophenoxypropionic acid, -2,4,5	93-72-1					9.3E+03	2.2E+04		6.6E+03
3.0E+01	I			5.0E-03	I			V	1.28E+03	1.36E+09	1.50E+04	1		Trichloropropane, 1,1,2-	598-77-6				1.1E-01	5.8E+03			5.8E+03
				4.0E-03	I	3.0E-04	I	V	1.40E+03	1.36E+09	1.57E+04	1		Trichloropropane, 1,2,3-	96-18-4	1.1E-01			1.1E-01	4.7E+03		2.1E+01	2.1E+01
				3.0E-03	X	3.0E-04	P	V	3.11E+02	1.36E+09	2.34E+03	1		Trichloropropene, 1,2,3-	96-19-5					3.5E+03		3.1E+00	3.1E+00
				2.0E-02	A					1.36E+09		1	0.1	Tricresyl Phosphate (TCP)	1330-78-5					2.3E+04	5.5E+04		1.6E+04
				3.0E-03	I					1.36E+09		1	0.1	Tridiphan	58138-08-2					3.5E+03	8.3E+03		2.5E+03
						7.0E-03	I	V	2.79E+04	1.36E+09	1.58E+04	1		Triethylamine	121-44-8							4.8E+02	4.8E+02
				2.0E+00	P					1.36E+09		1	0.1	Triethylene Glycol	112-27-6					2.3E+06	5.5E+06		1.6E+06
7.7E-03	I			7.5E-03	I	2.0E+01	P																

Key: I = IRIS; P = PPRTV; O = OPP; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (see FO #31); H = HEAST; F = See FAQ; W = see user guide Section 2.3.5; E = see user guide Section 2.3.6; L = see user's guide Section 5.2; M = mutagen; S = see user's guide Section 5; V = volatile; R = RBA applied (see user's guide Section 5.10); c = cancer; n = noncancer; * = where: n SL < 100X c SL; ** = where n SL < 10X c SL; SSL values are based on DAF=1; m = Concentration may exceed ceiling limit (see user's guide Section 5.13); s = concentration may exceed Csat (see user's guide Section 5.12)

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³ -y) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _o (mg/m ³)	k _e y	v	o	muta gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	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 THI=1 (mg/kg)	
8.3E-03	P	9.0E-03	I	7.0E-06	P	1.36E+09					0.026	1.36E+09		0.026		Vanadium Pentoxide Vanadium and Compounds	1314-62-1 7440-62-2			2.0E+03	2.0E+03	1.1E+04 5.9E+03		4.2E+04 6.0E+05	8.4E+03 5.8E+03	
		1.0E-03	I		V	1.36E+09						1.23E+05		1		Vernolate	1929-77-7					1.2E+03			1.2E+03	
		1.2E-03	O			1.36E+09								1	0.1	Vinclozolin	50471-44-8					1.4E+03	3.3E+03			9.8E+02
		1.0E+00	H	2.0E-01	I	2.75E+03	1.36E+09	4.40E+03	1					1		Vinyl Acetate	108-05-4					1.2E+06		3.9E+03	3.8E+03	
7.2E-01	I	3.2E-05	H	3.0E-03	I	2.47E+03	1.36E+09	1.37E+03	1					1		Vinyl Bromide	593-60-2			5.2E-01	5.2E-01			1.8E+01	1.8E+01	
		4.4E-06	I	1.0E-01	I	3.92E+03	1.36E+09	9.56E+02	1					1		Vinyl Chloride	75-01-4	4.5E+00		2.7E+00	1.7E+00	3.5E+03		4.2E+02	3.7E+02	
				3.0E-04	I		1.36E+09							1	0.1	Warfarin	81-81-2					3.5E+02	8.3E+02		2.5E+02	
		2.0E-01	S	1.0E-01	S	3.90E+02	1.36E+09	5.58E+03	1					1		Xylene, p-	106-42-3					2.3E+05		2.4E+03	2.4E+03	
		2.0E-01	S	1.0E-01	S	3.88E+02	1.36E+09	5.47E+03	1					1		Xylene, m-	108-38-3					2.3E+05		2.4E+03	2.4E+03	
		2.0E-01	S	1.0E-01	S	4.34E+02	1.36E+09	6.46E+03	1					1		Xylene, o-	95-47-6					2.3E+05		2.8E+03	2.8E+03	
		2.0E-01	I	1.0E-01	I	2.60E+02	1.36E+09	5.74E+03	1					1		Xylenes	1330-20-7					2.3E+05		2.5E+03	2.5E+03	
		3.0E-04	I			1.36E+09								1		Zinc Phosphide	1314-84-7					3.5E+02			3.5E+02	
		3.0E-01	I			1.36E+09								1		Zinc and Compounds	7440-66-6					3.5E+05			3.5E+05	
		5.0E-02	I			1.36E+09								1	0.1	Zinc	12122-67-7					5.8E+04	1.4E+05		4.1E+04	
		8.0E-05	X			1.36E+09								1		Zirconium	7440-67-7					9.3E+01			9.3E+01	

TR=1E-06
THQ=1.0