

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

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

Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1					
SFO (mg/kg-day) <sup>-1</sup>	ke y	IUR (ug/m <sup>3</sup> ) <sup>-1</sup>	ke y	RfD <sub>o</sub> (mg/kg-day)	ke y	RfC <sub>i</sub> (mg/m <sup>3</sup> )	ke y	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)
1.8E-02	C	5.1E-06	C	1.5E-01	I				1.0E+00	1.0E-01		1.4E+09		ALAR	1596-84-5	1.8E+02	4.3E+02	3.3E+06	1.3E+02	1.8E+05	4.1E+05		1.2E+05
8.7E-03	I			4.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Acephate	30560-19-1	3.8E+02	8.9E+02		2.6E+02	4.7E+03	1.1E+04		3.3E+03
		2.2E-06	I			9.0E-03	I	V	1.0E+00		1.1E+05	1.4E+09	8.7E+03	Acetaldehyde	75-07-0			4.9E+01			3.4E+02		3.4E+02
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Acetochlor	34256-82-1					2.3E+04	5.5E+04		1.6E+04
				9.0E-01	I	3.1E+01	A	V	1.0E+00		1.1E+05	1.4E+09	1.4E+04	Acetone	67-64-1					1.1E+06		1.8E+06	6.7E+05
						2.0E-03	X	V	1.0E+00		1.1E+05	1.4E+09	2.4E+04	Acetone Cyanohydrin	75-86-5						2.1E+02		2.1E+02
				6.0E-02	I				1.0E+00		1.3E+05	1.4E+09	1.3E+04	Acetonitrile	75-05-8							3.4E+03	3.4E+03
3.8E+00	C	1.3E-03	C						1.0E+00	1.0E-01		1.4E+09		Acetophenone	98-86-2					1.2E+05			1.2E+05
				5.0E-04	I	2.0E-05	I	V	1.0E+00		2.3E+04	1.4E+09	6.9E+03	Acetylaminofluorene, 2-	53-96-3	8.6E-01	2.0E+00	1.3E+04	6.0E-01				
5.0E-01	I	1.0E-04	I	2.0E-03	I	6.0E-03	I	M	1.0E+00	1.0E-01		1.4E+09		Acrolein	107-02-8	6.5E+00	1.5E+01	1.7E+05	4.6E+00	5.8E+02	5.5E+03	6.1E-01	6.0E-01
				5.0E-01	I	1.0E-03	I	V	1.0E+00		1.1E+05	1.4E+09	9.5E+04	Acrylamide	79-06-1					2.3E+03		3.6E+07	1.6E+03
									1.0E+00		1.1E+05	1.4E+09	9.5E+04	Acrylic Acid	79-10-7					5.8E+05	4.2E+02		4.2E+02
5.4E-01	I	6.8E-05	I	4.0E-02	A	2.0E-03	I	V	1.0E+00		1.1E+04	1.4E+09	7.7E+03	Acrylonitrile	107-13-1	6.1E+00		1.4E+00	1.1E+00	4.7E+04		6.7E+01	6.7E+01
						6.0E-03	P		1.0E+00	1.0E-01		1.4E+09		Adiponitrile	111-69-3					1.2E+04	2.8E+04		3.6E+07
5.6E-02	C			1.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Alachlor	15972-60-8	5.8E+01	1.4E+02		4.1E+01	1.2E+04	2.8E+04		8.2E+03
				1.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Aldicarb	116-06-3					1.2E+03	2.8E+03		8.2E+02
				1.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Aldicarb Sulfone	1646-88-4					1.2E+03	2.8E+03		8.2E+02
									1.0E+00	1.0E-01		1.4E+09		Aldicarb sulfoxide	1646-87-3								
1.7E+01	I	4.9E-03	I	3.0E-05	I				1.0E+00			1.4E+09	1.7E+06	Aldrin	309-00-2	1.9E-01		4.3E+00	1.8E-01	3.5E+01			3.5E+01
				2.5E-01	I				1.0E+00	1.0E-01		1.4E+09		Allyl	74223-64-6					2.9E+05	6.9E+05		2.1E+05
				5.0E-03	I	1.0E-04	X	V	1.0E+00		1.1E+05	1.4E+09	3.4E+04	Allyl Alcohol	107-18-6					5.8E+03		1.5E+01	1.5E+01
2.1E-02	C	6.0E-06	C			1.0E-03	I	V	1.0E+00		1.4E+03	1.4E+09	1.6E+03	Allyl Chloride	107-05-1	1.6E+02		3.2E+00	3.2E+00			6.9E+00	6.9E+00
				1.0E+00	P	5.0E-03	P		1.0E+00			1.4E+09		Aluminum	7429-90-5					1.2E+06		3.0E+07	1.1E+06
				4.0E-04	I				1.0E+00			1.4E+09		Aluminum Phosphide	20859-73-8					4.7E+02			4.7E+02
				3.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Amdri	67485-29-4					3.5E+02	8.3E+02		2.5E+02
2.1E+01	C	6.0E-03	C			9.0E-03	I		1.0E+00	1.0E-01		1.4E+09		Amethrin	834-12-8	1.6E-01	3.7E-01	2.8E+03	1.1E-01	1.1E+04	2.5E+04		7.4E+03
									1.0E+00	1.0E-01		1.4E+09		Aminodiphenyl, 4-	92-67-1								
				8.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Aminophenol, m-	591-27-5					9.3E+04	2.2E+05		6.6E+04
				2.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Aminophenol, p-	123-30-8					2.3E+04	5.5E+04		1.6E+04
				2.5E-03	I				1.0E+00	1.0E-01		1.4E+09		Amtraz	33089-61-1					2.9E+03	6.9E+03		2.1E+03
				2.0E-01	I	1.0E-01	I	V	1.0E+00			1.4E+09		Ammonia	7804-41-7					2.3E+05			2.3E+05
						3.0E-03	X	V	1.0E+00		1.4E+04	1.4E+09	2.6E+04	Ammonium Sulfamate	7773-06-0								3.4E+02
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I		1.0E+00	1.0E-01		1.4E+09		Amyl Alcohol, tert.	75-85-4							6.0E+06	3.4E+02
4.0E-02	P			2.0E-03	X				1.0E+00	1.0E-01		1.4E+09		Aniline	62-53-3	5.7E+02	1.4E+03	1.0E+07	4.0E+02	8.2E+03	1.9E+04		5.7E+03
				4.0E-04	I				1.5E-01			1.4E+09		Anthrquinone, 9,10-	64-65-1	8.2E+01	1.9E+02		5.7E+01	2.3E+03	5.5E+03		1.6E+03
									1.5E-01			1.4E+09		Antimony (metallic)	7440-36-0					4.7E+02			4.7E+02
				5.0E-04	H				1.5E-01			1.4E+09		Antimony Pentoxide	1314-60-9					5.8E+02			5.8E+02
				9.0E-04	H				1.5E-01			1.4E+09		Antimony Potassium Tartrate	11071-15-1					1.1E+03			1.1E+03
				4.0E-04	H				1.5E-01			1.4E+09		Antimony Tetroxide	1332-81-6					4.7E+02			4.7E+02
				2.0E-04	I				1.5E-01			1.4E+09		Antimony Trioxide	1309-64-4							1.2E+06	1.2E+06
2.5E-02	I	7.1E-06	I	1.3E-02	I				1.0E+00	1.0E-01		1.4E+09		Apollo	74115-24-5	1.3E+02	3.1E+02	2.3E+06	9.2E+01	1.5E+04	3.6E+04		1.1E+04
				5.0E-02	H				1.0E+00	1.0E-01		1.4E+09		Aramite	140-67-8					5.8E+04	1.4E+05		4.1E+04
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C		1.0E+00	3.0E-02		1.4E+09		Arsenic, Inorganic	7440-38-2	3.6E+00	1.7E+01	3.9E+03	3.0E+00	5.8E+02	2.8E+03	8.9E+04	4.8E+02
				3.5E-06	C	5.0E-05	I		1.0E+00			1.4E+09		Arsine	7784-42-1					4.1E+00		3.0E+05	4.1E+00
				9.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Assure	76578-14-8					1.1E+04	2.5E+04		7.4E+03
				5.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Asulam	3337-71-1					5.8E+04	1.4E+05		4.1E+04
2.3E-01	C			3.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Atrazine	1912-24-9	1.4E+01	3.4E+01	6.7E+04	1.0E+01	4.1E+04	9.7E+04		2.9E+04
8.8E-01	C	2.5E-04	C						1.0E+00	1.0E-01		1.4E+09		Auramine	492-80-8	3.7E+00	8.8E+00		2.6E+00				
1.1E-01	I	3.1E-05	I						1.0E+00	1.0E-01		1.4E+09		Avermectin B1	65195-55-3					4.7E+02	1.1E+03		3.3E+02
				1.0E+00	P	7.0E-06	P		1.0E+00	1.0E-01		1.4E+09	5.2E+05	Azobenzene	103-33-3	3.0E+01		2.1E+02	2.6E+01	1.2E+06	2.8E+06	4.2E+04	4.0E+04
									1.0E+00	1.0E-01		1.4E+09		Azodicarbonamide	123-77-3								
				2.0E-01	I	5.0E-04	H		7.0E-02			1.4E+09		Barium	7440-39-3					2.3E+05		3.0E+06	2.2E+05
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M	2.5E-02			1.4E+09		Barium Chromate	10294-40-3	6.5E+00		1.1E+02	6.2E+00	2.3E+04		1.2E+06	2.3E+04
				4.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Baygon	114-26-1					4.7E+03	1.1E+04		3.3E+03
				3.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Bayleton	43121-43-3					3.5E+04	8.3E+04		2.5E+04
				2.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Baythroid	68359-37-5</								

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Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day) <sup>-1</sup>	ke IUR (ug/m <sup>3</sup> -y) <sup>-1</sup>	ke RfD <sub>o</sub> (mg/kg-day)	ke RfC <sub>i</sub> (mg/m <sup>3</sup> )	ke V c mutagen	GIABS	ABS	C <sub>sat</sub> (m <sup>3</sup> /kg)	PEF	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)			
1.3E+01	I			V	1.0E+00		3.2E+02	1.4E+09	6.8E+04	Benzotrithloride	98-07-7	2.5E-01			2.5E-01							
1.7E-01	I	4.9E-05 C 2.4E-03 I	1.0E-01 P 2.0E-03 I 2.0E-03 I	P 1.0E-03 P 2.0E-05 I	V	1.0E+00 1.0E+00 7.0E-03	1.5E+03	1.4E+09 1.4E+09	2.6E+04	Benzyl Alcohol Benzyl Chloride Beryllium and compounds	100-51-6 100-44-7 7440-41-7	1.9E+01		6.4E+00 6.9E+03	4.8E+00 6.9E+03	1.2E+05 2.3E+03 2.3E+03	2.8E+05	1.1E+02 1.2E+05	8.2E+04 1.1E+02 2.3E+03			
			1.0E-04 I 9.0E-03 P 1.5E-02 I		V	1.0E+00 1.0E+00 1.0E+00		1.4E+09 1.4E+09 1.4E+09		Bidrin Bifenox Biphenrin	141-66-2 42576-02-3 82657-04-3					1.2E+02 1.1E+04 1.8E+04	2.8E+02 2.5E+04 4.1E+04		8.2E+01 7.4E+03 1.2E+04			
8.0E-03 7.0E-02	I H	1.0E-05 H	4.0E-02 I 3.0E-03 P	I 4.0E-04 X V	V	1.0E+00 1.0E+00	1.0E+03	1.4E+09 1.4E+09	3.5E+04	Biphenyl, 1,1'- Bis(2-chloro-1-methylethyl) ether Bis(2-chloroethoxy)methane	92-52-4 108-60-1 111-91-1	4.1E+02 4.7E+01		4.3E+01 2.2E+01	4.1E+02	5.8E+05 4.7E+04 3.5E+03	2.0E+02 4.1E+04	2.0E+02 4.7E+04 2.5E+03				
1.1E+00 2.2E+02	I I	3.3E-04 I 6.2E-02 I		V	V	1.0E+00 1.0E+00	5.1E+03 4.2E+03	1.4E+09 1.4E+09	4.3E+04 1.9E+03	Bis(2-chloroethyl)ether Bis(chloromethyl)ether Bisphenol A	111-44-4 542-88-1 80-05-7	3.0E+00 1.5E-02		1.6E+00 3.7E-04	1.0E+00 3.6E-04	5.8E+04 1.4E+05			4.1E+04			
			2.0E-01 I 2.0E+00 P 4.0E-02 C	I 2.0E-02 H P 1.3E-02 C	V	1.0E+00 1.0E+00 1.0E+00		1.4E+09 1.4E+09 1.4E+09		Boron And Borates Only Boron Trichloride Boron Trifluoride	7440-42-8 10294-34-5 7637-07-2					2.3E+05 2.3E+06 4.7E+04	1.2E+08 1.2E+08 7.7E+07	2.3E+05 2.3E+06 4.7E+04				
7.0E-01 2.0E+00	I X	6.0E-04 X	4.0E-03 I 6.0E-02 I	V	V	1.0E+00 1.0E+00	2.4E+03 6.8E+02	1.4E+09 1.4E+09	5.9E+03 8.4E+03	Bromate Bromo-2-chloroethane, 1- Bromobenzene	15541-45-4 107-04-0 108-86-1	4.7E+00 1.6E+00		1.2E-01 1.1E-01	4.7E+00	4.7E+03			4.7E+03			
			8.0E-03 I 4.0E-02 X V	V	V	1.0E+00 1.0E+00	4.0E+03 9.3E+02 9.2E+02	1.4E+09 1.4E+09 1.4E+09	3.8E+03 4.0E+03 9.7E+03	Bromochloromethane Bromodichloromethane Bromoform	74-97-5 75-27-4 75-25-2	5.3E+01 4.1E+02		1.3E+00 1.1E+02	1.3E+00 8.6E+01	2.3E+04 2.3E+04			2.3E+04 2.3E+04			
6.2E-02 7.9E-03	I I	3.7E-05 C 1.1E-06 I	2.0E-02 I 2.0E-02 I	V	V	1.0E+00 1.0E+00	9.3E+02 9.2E+02	1.4E+09 1.4E+09	4.0E+03 9.7E+03	Bromodichloromethane Bromoform	75-27-4 75-25-2	5.3E+01 4.1E+02		1.3E+00 1.1E+02	1.3E+00 8.6E+01	2.3E+04 2.3E+04			2.3E+04 2.3E+04			
			1.4E-03 I 5.0E-03 H 2.0E-02 I	I 5.0E-03 V V	V	1.0E+00 1.0E+00 1.0E+00	3.6E+03	1.4E+09 1.4E+09 1.4E+09	1.4E+03 1.2E+05	Bromomethane Bromophos Bromoxynil	74-83-9 2104-96-3 1689-84-5					1.6E+03 5.8E+03 2.3E+04		3.1E+01	3.0E+01 5.8E+03 1.6E+04			
3.4E+00	C	3.0E-05 I	2.0E-02 I 2.0E-03 I	V I	V	1.0E+00 1.0E+00	6.7E+02 7.6E+03	1.4E+09 1.4E+09	4.7E+05 8.7E+02 3.0E+04	Bromoxynil Octanoate Butadiene, 1,3- Butarol, N-	1689-99-2 106-99-0 11-36-3	9.6E-01		3.5E-01	2.6E-01	2.3E+04			2.3E+04			
1.9E-03	P		2.0E-01 I 2.0E+00 P 5.0E-02 I	I 3.0E+01 P V	V	1.0E+00 1.0E+00 1.0E+00	1.4E+09 2.1E+04	1.4E+09 1.4E+09	1.0E-01 2.9E+04 8.6E+04	Butyl Benzyl Phthalate Butyl alcohol, sec- Butylate	85-88-7 78-92-2 2008-41-5	1.7E+03	4.1E+03		1.2E+03	2.3E+05 2.3E+06 5.8E+04	5.5E+05	3.8E+06	1.6E+05 1.5E+06 5.8E+04			
2.0E-04 3.6E-03	C P	5.7E-08 C	3.0E-01 P 5.0E-02 P	P V	V	1.0E+00 1.0E+00	1.1E+02	1.4E+09 1.4E+09	8.1E+03	Butylated hydroxyanisole Butylated hydroxytoluene Butylbenzene, n-	25013-16-5 128-37-0 104-51-8	1.6E+04 9.1E+02	3.9E+04 2.1E+03	2.9E+08 6.4E+02	1.1E+04	3.5E+05 5.8E+04	8.3E+05		2.5E+05 5.8E+04			
			1.0E-01 X 1.0E-01 X 2.0E-02 A	V V V	V	1.0E+00 1.0E+00 1.0E+00	1.5E+02 1.8E+02	1.4E+09 1.4E+09	7.4E+03 7.4E+03	Butylbenzene, sec- Butylbenzene, tert- Cacodylic Acid	135-98-8 98-06-6 75-60-5					1.2E+05 1.2E+05 2.3E+04	5.5E+04		1.2E+05 1.2E+05 1.6E+04			
1.8E-03 1.8E-03 5.0E-01	I I C	1.0E-03 I 5.0E-04 I 1.5E-01 C	1.0E-05 A 1.0E-05 A 2.0E-02 C	I I C	A A M	2.5E-02 5.0E-02 2.5E-02	1.0E-03 1.0E-03	1.4E+09 1.4E+09		Cadmium (Diet) Cadmium (Water) Calcium Chromate	7440-43-9 7440-43-9 13765-19-0			9.3E+03 9.3E+03	9.3E+03	1.2E+03 1.2E+03	6.9E+03	6.0E+04	9.8E+02 9.8E+02 2.3E+04			
1.5E-01 2.3E-03	C C	4.3E-05 C 6.6E-07 C	5.0E-01 I 2.0E-03 I 1.3E-01 I	I I I	C	1.0E+00 1.0E+00 1.0E+00	1.4E+09 1.4E+09 1.4E+09	1.4E+09 1.4E+09		Caprolactam Captalof Captan	105-80-2 2425-06-1 133-06-2	2.2E+01 1.4E+03	5.2E+01 3.4E+03	3.9E+05 2.5E+07	1.5E+01 1.0E+03	5.8E+05 2.3E+03 1.5E+05	1.4E+06 5.5E+03 3.6E+05	1.3E+07	4.0E+05 1.6E+03 1.1E+05			
			1.0E-01 I 5.0E-03 I 1.0E-01 I	I I I	V	1.0E+00 1.0E+00 1.0E+00	7.4E+02	1.4E+09 1.4E+09 1.4E+09	1.2E+03	Carbaryl Carbofuran Carbon Disulfide	63-25-2 1563-66-2 75-15-0					1.2E+05 5.8E+03 1.2E+05	2.8E+05 1.4E+04	3.6E+03	8.2E+04 4.1E+03 3.5E+03			
7.0E-02	I	6.0E-06 I	4.0E-03 I 1.0E-02 I 1.0E-01 I	I I I	V	1.0E+00 1.0E+00 1.0E+00	4.6E+02	1.4E+09 1.4E+09 1.4E+09	1.5E+03	Carbon Tetrachloride Carbosulfan Carboxin	56-23-5 55285-14-8 5234-68-4	4.7E+01		3.1E+00	2.9E+00	4.7E+03 1.2E+04 1.2E+05	2.8E+04 2.8E+05		5.7E+02 8.2E+03 8.2E+04			
			1.0E-01 I 1.5E-02 I	V I	V	1.0E+00 1.0E+00	1.4E+09 1.4E+09	1.4E+09 1.5E+05		Ceric oxide Chloral Hydrate Chloramben	1306-38-3 302-17-0 133-90-4					1.2E+05 1.8E+04	4.1E+04		5.4E+06 1.2E+05 1.2E+04			
4.0E-01 3.5E-01 1.0E+01	H I I	1.0E-04 I 4.6E-03 C	5.0E-04 I 3.0E-04 I	I I	V	1.0E+00 1.0E+00	1.4E+09 1.4E+09	1.4E+09 1.4E+09	9.0E+05	Chloranil Chlordane Chlordecone (Kepone)	118-75-2 12789-03-6 143-50-0	8.1E+00 9.3E+00 3.3E-01	1.9E+01 5.5E+01 7.7E-01	1.1E+02 3.6E+03	5.7E+00 7.5E+00 2.3E-01	5.8E+02 3.5E+02	3.4E+03 8.3E+02	2.8E+03	4.2E+02 2.5E+02			
			7.0E-04 A 2.0E-02 I 1.0E-01 I	A I A	V	1.0E+00 1.0E+00 1.0E+00	2.8E+03	1.4E+09 1.4E+09 1.4E+09	1.2E+03	Chlorfenirphos Chlorfuran, Ethyl- Chlorine	470-90-6 90982-32-4 7782-50-5					8.2E+02 2.3E+04 1.2E+05	1.9E+03 5.5E+04	7.8E-01	5.7E+02 1.6E+04 7.8E-01			
			3.0E-02 I 3.0E-02 I	I I	V	1.0E+00 1.0E+00	1.4E+09 1.4E+09	1.4E+09 1.4E+09	1.0E+03	Chlorine Dioxide Chlorite (Sodium Salt) Chloro-1,1-difluoroethane, 1-	10049-04-4 7758-19-2 75-68-3					3.5E+04 3.5E+04		1.2E+06	3.4E+04 3.5E+04 2.3E+05			
4.6E-01 1.0E-01	H P	7.7E-05 C	2.0E-02 H 3.0E-03 X	I I	V	1.0E+00 1.0E+00	7.5E+02	1.4E+09 1.4E+09	1.1E+03	Chloro-1,3-butadiene, 2- Chloro-2-methylaniline HCl, 4- Chloro-2-methylaniline, 4-	126-99-8 3165-93-3 95-69-2	7.1E+00 3.3E+01	1.7E+01 7.7E+01	4.4E-02 2.2E+05	4.4E-02 2.3E+01	2.3E+04 3.5E+03	8.3E+03		9.4E+01 2.5E+03			

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

Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) <sup>-1</sup>	ke IUR (ug/m <sup>3</sup> -y) <sup>-1</sup>	ke (mg/kg-day)	RfD <sub>o</sub> (mg/kg-day)	ke RfC <sub>i</sub> (mg/m <sup>3</sup> -y)	ke V o c	muta-gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)		
2.7E-01	X		2.0E-03	H	3.0E-05	I			1.0E+00	2.8E+04	1.4E+09	1.9E+04	Chloroacetaldehyde, 2-Chloroacetic Acid Chloroacetophenone, 2-	107-20-0 79-11-8 532-27-4	1.2E+01			1.2E+01				2.3E+03 5.5E+03 1.8E+05	1.6E+03 1.8E+05
2.0E-01	P		4.0E-03	I	5.0E-02	P V			1.0E+00	1.0E-01	1.4E+09	Chloroaniline, p-Chlorobenzene Chlorobenzilate	106-47-8 108-90-7 510-15-6	1.6E+01 3.9E+01		1.1E+01		4.7E+03 2.3E+04 2.3E+04	1.1E+04		1.4E+03	3.3E+03 1.3E+03 1.6E+04	
1.1E-01	C	3.1E-05	2.0E-02	I					1.0E+00	1.0E-01	1.4E+09	Chlorobenzoic Acid, p-Chlorobenzotrifluoride, 4-Chlorobutane, 1-Chlorodifluoromethane Chloroethanol, 2-Chloroform	74-11-3 98-56-6 109-69-3 75-45-6 107-07-3 67-66-3	3.0E+01 7.0E+01	5.4E+05	2.1E+01		3.5E+04 8.3E+04		8.9E+03	2.5E+04 2.5E+03 4.7E+04		
3.1E-02	C	2.3E-05	1.0E-02	I	9.8E-02	A V			1.0E+00	2.5E+03	1.4E+09	2.6E+03	Chloromethane Chloromethyl Methyl Ether Chloronitrobenzene, o-Chlorotoluene, p-Chlorotoluene, p-Chlorozotocin Chlorpropylamine Chloropyridos	74-87-3 107-30-2 88-73-3 54749-90-5 101-21-3 2921-88-2	1.1E+02	1.4E+00	1.4E+00		2.3E+04 1.2E+04		1.1E+03	2.1E+05 2.1E+05 2.3E+04 1.0E+03	
2.4E+00	C	6.9E-04	3.0E-03	P	1.0E-05	X			1.0E+00	2.6E+04	1.4E+09	5.3E+03	Chloromethyl Methyl Ether Chloronitrobenzene, o-Chlorotoluene, p-Chlorotoluene, p-Chlorozotocin Chlorpropylamine Chloropyridos	107-30-2 88-73-3 106-43-4 54749-90-5 101-21-3 2921-88-2	1.4E+00 1.1E+01	2.6E+01	9.5E-02	8.9E-02		3.5E+03 8.3E+03	6.0E+04	2.4E+03	
6.3E-03	P		1.0E-03	P	6.0E-04	P			1.0E+00	1.0E-01	1.4E+09	1.4E+09	Chloronitrobenzene, p-Chlorophenol, 2-Chloropicrin	100-00-5 95-57-8 76-06-2	5.2E+02	1.2E+03		3.6E+02		1.2E+03 2.8E+03	3.6E+06	8.2E+02 5.8E+03 8.2E+00	
3.1E-03	C	8.9E-07	1.5E-02	I					1.0E+00	1.0E-01	1.4E+09	8.1E+03	Chlorothalonil Chlorotoluene, o-Chlorotoluene, p-Chlorotoluene, p-Chlorozotocin Chlorpropylamine Chloropyridos	1897-45-6 95-49-8 106-43-4 54749-90-5 101-21-3 2921-88-2	1.1E+03	2.5E+03	1.9E+07	7.4E+02		1.8E+04 2.3E+04 2.3E+04	4.1E+04	1.2E+04 2.3E+04 2.3E+04	
2.4E+02	C	6.9E-02	2.0E-01	I					1.0E+00	1.0E-01	1.4E+09	1.4E+09	Chlorotoluene, p-Chlorotoluene, p-Chlorozotocin Chlorpropylamine Chloropyridos	106-43-4 54749-90-5 101-21-3 2921-88-2	1.4E-02	3.2E-02	2.4E+02	9.6E-03		2.3E+05 1.2E+03	5.5E+05 2.8E+03	1.6E+05 8.2E+02	
5.0E-01	J	8.4E-02	3.0E-03	I	1.0E-04	I	M		2.5E-02	1.4E+09	1.4E+09	1.4E+09	Chromium(III) Insoluble Salts Chromium(VI) Chromium, Total	16065-83-1 18540-29-9 7440-47-3	6.5E+00		2.0E+02	6.3E+00		1.8E+06 3.5E+03	6.0E+05	1.8E+06 3.5E+03	
9.0E-03	P	6.2E-04	3.0E-04	P	6.0E-06	P			1.0E+00	1.0E+00	1.4E+09	1.4E+09	Cobalt Coke Oven Emissions Copper	4440-48-4 8007-45-2 7440-50-8		1.9E+03	1.9E+03		3.5E+02		3.6E+04	3.5E+02	
5.0E-02	I		6.0E-01	C					1.0E+00	1.0E-01	1.4E+09	1.4E+09	Cresol, m-Cresol, o-Cresol, p-Cresol	108-39-4 95-48-7 106-44-5					5.8E+04 5.8E+04 1.2E+05	1.4E+05 1.4E+05 2.8E+05	3.6E+09 3.6E+09 3.6E+09	4.1E+04 4.1E+04 8.2E+04	
1.9E+00	H		1.0E-01	A	6.0E-01	C			1.0E+00	1.0E-01	1.4E+09	1.4E+09	Cresol, p-chloro-m-Cresol	59-50-7 1319-77-3	1.1E+00		1.7E+00		1.2E+05 1.2E+05	2.8E+05 2.8E+05	3.6E+09	8.2E+04 8.2E+04	
2.2E-01	C	6.3E-05	2.0E-03	H					1.0E+00	1.0E-01	1.4E+09	1.4E+09	Cumene Cupferron Cyanazine	98-82-8 135-20-6 21725-46-2	1.5E+01 3.9E+00	3.5E+01 9.2E+00	2.6E+05	1.0E+01 2.7E+00		1.2E+05 2.3E+03	5.5E+03	1.1E+04	9.9E+03 1.6E+03
1.0E-03	I		1.0E-03	I					1.0E+00	1.0E+00	1.4E+09	1.4E+09	Cyanides	592-01-8 544-92-3					1.2E+03 5.8E+03			1.2E+03 5.8E+03	
6.0E-04	I	8.0E-04	1.0E-03	I	8.0E-04	S V			1.0E+00	9.7E+05	1.4E+09	3.5E+03	Cyanide (CN <sup>-</sup> )	57-12-5					7.0E+02		1.2E+01	1.2E+01	
1.0E-03	I		1.0E-03	I					1.0E+00	1.0E+00	1.4E+09	1.4E+09	Cyanogen	460-19-5					1.2E+03			1.2E+03	
9.0E-02	I		1.0E-02	I					1.0E+00	1.0E+00	1.4E+09	1.4E+09	Cyanogen Bromide	506-68-3					1.1E+05			1.1E+05	
5.0E-02	I		1.0E-02	I					1.0E+00	1.0E+00	1.4E+09	1.4E+09	Cyanogen Chloride	506-77-4					5.8E+04			5.8E+04	
6.0E-04	I	8.0E-04	1.0E-03	I	8.0E-04	I V			1.0E+00	1.0E+07	1.4E+09	5.2E+04	Hydrogen Cyanide	74-90-8					7.0E+02		1.8E+02	1.5E+02	
2.0E-03	I		1.0E-03	I					1.0E+00	1.0E+00	1.4E+09	1.4E+09	Potassium Cyanide	151-50-8					2.3E+03			2.3E+03	
5.0E-03	I		4.0E-02	I					4.0E-02	1.4E+09	1.4E+09	1.4E+09	Potassium Silver Cyanide	506-61-6					5.8E+03			5.8E+03	
1.0E-01	I		1.0E-01	I					4.0E-02	1.4E+09	1.4E+09	1.4E+09	Silver Cyanide	506-64-9					1.2E+05			1.2E+05	
1.0E-03	I		1.0E-03	I					1.0E+00	1.4E+09	1.4E+09	1.4E+09	Sodium Cyanide	143-33-9					1.2E+03			1.2E+03	
2.0E-04	P		1.0E-04	I					1.0E+00	1.4E+09	1.4E+09	1.4E+09	Thiocyanates	NA					2.3E+02			2.3E+02	
2.0E-04	X		1.0E-04	I					1.0E+00	1.4E+09	1.4E+09	1.4E+09	Thiocyanic Acid	463-56-9					2.3E+02			2.3E+02	
5.0E-02	I		1.0E-02	I					1.0E+00	1.4E+09	1.4E+09	1.4E+09	Zinc Cyanide	557-21-1					5.8E+04			5.8E+04	
2.3E-02	H		6.0E+00	I V					1.0E+00	1.2E+02	1.4E+09	1.0E+03	Cyclohexane	110-82-7	1.4E+02	3.4E+02		1.0E+02		2.7E+04	2.7E+04		
5.0E+00	I	7.0E-01	1.0E+00	P V					1.0E+00	5.1E+03	1.4E+09	4.2E+04	Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-Cyclohexanone	87-84-3 108-94-1					5.8E+06		1.3E+05	1.3E+05	
5.0E-03	P	1.0E+00	1.0E+00	X V					1.0E+00	2.8E+02	1.4E+09	1.5E+03	Cyclohexene	110-83-8					5.8E+03		6.4E+03	3.1E+03	
2.0E-01	I		1.0E-01	I					1.0E+00	2.9E+05	1.4E+09	7.5E+04	Cyclohexylamine	108-91-8					2.3E+05			2.3E+05	
5.0E-03	I		1.0E-03	I					1.0E+00	1.0E-01	1.4E+09	1.4E+09	Cyhalothrin/karate	68085-85-8					5.8E+03	1.4E+04		4.1E+03	
1.0E-02	I		1.0E-02	I					1.0E+00	1.0E-01	1.4E+09	1.4E+09	Cypermethrin	52315-07-8					1.2E+04	2.8E+04		8.2E+03	

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

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) <sup>-1</sup>	ke y <sup>-1</sup>	IUR (ug/m <sup>3</sup> -y) <sup>-1</sup>	ke y <sup>-1</sup>	RfD <sub>o</sub> (mg/kg-day)	ke y <sup>-1</sup>	RfC <sub>i</sub> (mg/m <sup>3</sup> -y)	ke y <sup>-1</sup>	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
2.4E-01	I	6.9E-05	C	7.5E-03	I				1.0E+00	1.0E-01		1.4E+09		Cyromazine	66215-27-8									
3.4E-01	I	9.7E-05	C	2.0E-04	I			V	1.0E+00	1.0E-01		1.4E+09	2.1E+08	DDD	72-54-8	1.4E+01	3.2E+01	2.4E+05	9.6E+00	8.8E+03	2.1E+04		6.2E+03	
3.4E-01	I	9.7E-05	I	5.0E-04	I				1.0E+00	3.0E-02		1.4E+09		DDE, p,p'	72-55-9	9.6E+00	2.7E+02	9.3E+00	9.6E+00	9.6E+00	9.3E+00			
7.0E-04	I			1.0E-02	I				1.0E+00	1.0E-01		1.4E+09		DDT	50-29-3	9.6E+00	7.6E+01	1.7E+05	8.5E+00	5.8E+02	4.6E+03		5.2E+02	
6.1E-02	H			3.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Dacihal	1861-32-1					1.2E+04	2.8E+04		8.2E+03	
1.2E-03	I			7.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Dalapon	75-99-0					3.5E+04	8.3E+04		2.5E+04	
6.1E-02	H			4.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'-(BDE-209)	1163-19-5	4.7E+03	1.1E+04		3.3E+03	8.2E+03	1.9E+04		5.7E+03	
8.0E-01	P	6.0E-03	P	1.0E-02	A			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Demeton	8065-48-3					4.7E+01	1.1E+02		3.3E+01	
8.4E-02	I	2.7E-05	C	2.0E-02	I			V	1.0E+00	1.0E-01		1.4E+09		Di(2-ethylhexyl)adipate	103-23-1	2.7E+03	6.4E+03		1.9E+03	7.0E+05	1.7E+06		4.9E+05	
2.0E+00	I	6.0E-04	I	9.0E-03	X			V	1.0E+00	1.0E-01		1.4E+09		Diallate	2303-16-4	5.4E+01	1.3E+02		3.8E+01	8.2E+02	1.9E+03		5.7E+02	
8.4E-02	I	2.7E-05	C	1.0E-02	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dibenzothiophene	132-65-0					1.2E+04			1.2E+04	
2.0E+00	I	6.0E-04	I	2.0E-02	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dibromo-3-chloropropane, 1,2-	96-12-8	4.1E+00		6.5E-02	6.4E-02	2.3E+02		2.8E+01	2.5E+01	
8.4E-02	I	2.7E-05	C	1.0E-02	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dibromobenzene, 1,3-	108-36-1					4.7E+02			4.7E+02	
2.0E+00	I	6.0E-04	I	1.0E-02	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dibromobenzene, 1,4-	106-37-6					1.2E+04			1.2E+04	
2.0E+00	I	6.0E-04	I	2.0E-02	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dibromochloromethane	124-48-1	3.9E+01		3.6E+00	3.3E+00	2.3E+04			2.3E+04	
2.0E+00	I	6.0E-04	I	9.0E-03	X			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dibromomethane, 1,2-	106-93-4	1.6E+00		1.8E-01	1.6E-01	1.1E+04		3.4E+02	3.3E+02	
4.2E-03	P			3.0E-04	P			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dibromomethane (Methylene Bromide)	74-95-3					1.2E+04		9.9E+01	9.8E+01	
4.2E-03	P			3.0E-02	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dibutyltin Compounds	NA					3.5E+02	8.3E+02		2.5E+02	
4.2E-03	P			3.0E-02	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dicamba	1918-00-9					3.5E+04	8.3E+04		2.5E+04	
5.0E-02	I	4.0E-03	I	9.0E-02	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichloro-2-butene, 1,4-	764-41-0			3.6E-02	3.6E-02					
5.4E-03	C	1.1E-05	C	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichloro-2-butene, cis-1,4-	1476-11-5			3.2E-02	3.2E-02					
4.5E-01	I	3.4E-04	C	7.0E-02	A			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichloro-2-butene, trans-1,4-	110-57-6			3.2E-02	3.2E-02					
5.7E-03	C	1.6E-06	C	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichloroacetic Acid	79-43-6	6.5E+01	1.5E+02		4.6E+01	4.7E+03	1.1E+04		3.3E+03	
9.1E-02	I	2.6E-05	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorobenzene, 1,2-	95-50-1					1.1E+05		1.0E+04	9.3E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorobenzene, 1,4-	106-46-7	6.1E+02		1.2E+01	1.1E+01	8.2E+04		3.7E+04	2.5E+04	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorobenzene, 1,3-	91-94-1	7.3E+00	1.7E+01	4.9E+04	5.1E+00	2.3E+05	2.5E+04		3.7E+02	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorobenzene, 1,4-	90-96-2					1.1E+04		2.5E+04	7.4E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorodifluoromethane	75-71-8					2.3E+05		3.7E+02	3.7E+02	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichloroethane, 1,1-	78-34-3	5.7E+02		1.6E+01	1.6E+01	2.3E+05			2.3E+05	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichloroethane, 1,2-	107-06-2	3.6E+01		2.2E+00	2.0E+00	7.0E+03		1.4E+02	1.4E+02	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichloroethylene, 1,1-	75-35-4					5.8E+04		1.0E+03	1.0E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichloroethylene, 1,2-cis-	156-59-2					7.0E+03		1.4E+02	1.4E+02	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichloroethylene, 1,2-trans-	156-60-5					5.8E+04		1.0E+03	1.0E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03	
1.6E+01	I	4.6E-03	I	2.0E-01	I			V	1.0E+00	1.0E-01		1.4E+09	5.2E+05	Dichlorophenyl, 2,4-	120-83-2					3.5E+03	8.3E+03	</		

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

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1						
SFO (mg/kg-day) <sup>-1</sup>	ke y	IUR (ug/m <sup>3</sup> -y) <sup>-1</sup>	ke y	RfD <sub>o</sub> (mg/kg-day)	ke y	RfC <sub>i</sub> (mg/m <sup>3</sup> -y)	ke y	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)
1.1E+01	P								1.0E+00	1.0E-01		1.4E+09		Dimethylbenzidine, 3,3'-	119-93-7	3.0E-01	7.0E-01		2.1E-01				
				1.0E-01	P	3.0E-02	I	V	1.0E+00		1.1E+05	1.4E+09	1.3E+05	Dimethylformamide	68-12-2					1.2E+05		1.7E+04	1.5E+04
5.5E+02	C	1.6E-01	C	1.0E-04	X	2.0E-06	X	V	1.0E+00		1.7E+05	1.4E+09	1.6E+05	Dimethylhydrazine, 1,1-Dimethylhydrazine, 1,2-	540-73-8	5.9E-03		1.3E-02	4.1E-03	1.2E+02		1.4E+00	1.4E+00
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Dimethylphenol, 2,4-Dimethylphenol, 2,6-Dimethylphenol, 3,4-	105-67-9 576-26-1 95-65-8				2.3E+04	5.5E+04		1.6E+04	
4.5E-02	C	1.3E-05	C					V	1.0E+00		1.1E+03	1.4E+09	1.0E+03	Dimethylvinylchloride	513-37-1		9.5E-01	9.4E-01					
				8.0E-05	X				1.0E+00	1.0E-01		1.4E+09		Dinitro-o-cresol, 4,6-Dinitro-o-cyclohexyl Phenol, 4,6-	534-52-1 131-89-5				9.3E+01	2.2E+02	2.3E+03	5.5E+03	6.6E+01 1.6E+03
				1.0E-04	P				1.0E+00	1.0E-01		1.4E+09		Dinitrobenzene, 1,2-Dinitrobenzene, 1,3-Dinitrobenzene, 1,4-	528-29-0 99-65-0 100-25-4				1.2E+02	2.8E+02	1.2E+02	2.8E+02	8.2E+01 8.2E+01 8.2E+01
6.8E-01	I			2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Dinitrophenol, 2,4-Dinitrotoluene Mixture, 2,4/2,6-Dinitrotoluene, 2,4-	51-28-5 NA 121-14-2	4.8E+00 1.1E+01	1.1E+01 2.4E+01	1.9E+05 7.4E+00	3.4E+00 7.4E+00	2.3E+03	5.4E+03		1.6E+03
3.1E-01	C	8.9E-05	C	2.0E-03	X				1.0E+00	1.0E-01		1.4E+09		Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 4-Amino-2,6-	606-20-2 36572-78-2 19406-51-0	2.2E+00	5.2E+00	1.5E+00	2.3E+03	5.4E+03		2.5E+02 2.3E+03 2.3E+03	
1.5E+00	P			3.0E-04	X				1.0E+00	9.9E-02		1.4E+09		Dinitrotoluene, 2,6-Dinitrotoluene, 2-Amino-4,6-Dinitrotoluene, 4-Amino-2,6-	606-20-2 36572-78-2 19406-51-0	2.2E+00	5.2E+00	1.5E+00	3.5E+02	8.4E+02		2.5E+02 2.3E+03 2.3E+03	
4.5E-01	X			9.0E-04	S				1.0E+00	1.0E-01		1.4E+09		Dinitrotoluene, Technical grade	25321-14-6	7.3E+00	1.7E+01		5.1E+00	1.1E+03	2.5E+03		7.4E+02
1.0E-01	I	5.0E-06	I	1.0E-03	I	3.0E-02	I	V	1.0E+00	1.0E-01	1.2E+05	1.4E+09	4.0E+04	Dinoseb	88-85-7					1.2E+03	2.8E+03		8.2E+02
				3.0E-02	I				1.0E+00			1.4E+09		Dioxane, 1,4-Dioxins	123-91-1	3.3E+01	9.7E+01	2.4E+01		3.5E+04		5.2E+03	4.5E+03
6.2E+03	I	1.3E+00	I						1.0E+00	3.0E-02		1.4E+09		--Hexachlorodibenzo-p-dioxin, Mixture --TCDD, 2,3,7,8-	NA 1746-01-6	5.3E-04 2.5E-05	4.2E-03 2.0E-04	1.3E+01 6.3E-04	4.7E-04 2.2E-05	8.2E-04	6.4E-03	3.4E-01	7.2E-04
1.3E+05	C	3.8E+01	C	7.0E-10	I	4.0E-08	C	V	1.0E+00	1.0E-01		1.4E+09	2.0E+06	Diphenamid	957-51-7					3.5E+04	8.3E+04		2.5E+04
				8.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Diphenyl Sulfone	127-63-9					9.3E+02	2.2E+03		6.6E+02
				2.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Diphenylamine	122-39-4					2.9E+04	6.9E+04		2.1E+04
8.0E-01	I	2.2E-04	I						1.0E+00	1.0E-01		1.4E+09		Diphenylhydrazine, 1,2-Diquat	122-86-7 85-00-7	4.1E+00	9.7E+00	7.6E+04	2.9E+00	2.6E+03	6.1E+03		1.8E+03
7.1E+00	C	1.4E-01	C	2.2E-03	I				1.0E+00	1.0E-01		1.4E+09		Direct Black 38	1937-37-7	4.6E-01	1.1E+00	1.2E+02	3.2E-01				
7.4E+00	C	1.4E-01	C						1.0E+00	1.0E-01		1.4E+09		Direct Blue 6	2602-46-2	4.4E-01	1.0E+00	1.2E+02	3.1E-01				
6.7E+00	C	1.4E-01	C						1.0E+00	1.0E-01		1.4E+09		Direct Brown 95	16071-86-6	4.9E-01	1.2E+00	1.2E+02	3.4E-01				
				4.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Disulfoton	298-04-4					4.7E+01	1.1E+02		3.3E+01
				1.0E-02	I			V	1.0E+00			1.4E+09	4.5E+04	Dithiane, 1,4-Diuron	505-29-3 330-54-1					1.2E+04			1.2E+04
				2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Iodine	20439-10-3					2.3E+03	5.5E+03		1.6E+03
				4.0E-03	I				1.0E+00	1.0E-01		1.4E+09		EPTC	759-94-4					4.7E+03	1.1E+04		3.3E+03
				2.5E-02	I			V	1.0E+00			1.2E+05		Endosulfan	115-29-7					2.9E+04			2.9E+04
				6.0E-03	I			V	1.0E+00			4.1E+05		Endothal	145-73-3					7.0E+03			7.0E+03
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Endrin	72-20-8					2.3E+04	5.5E+04		1.6E+04
9.9E-03	I	1.2E-06	I	3.0E-04	I	1.0E-03	I	V	1.0E+00	1.0E-01		1.4E+09		Epiclorohydrin	106-89-8	3.3E+02		1.9E+02	1.2E+02	3.5E+02	8.3E+02		2.5E+02
				6.0E-03	P	1.0E-03	I	V	1.0E+00		1.1E+04	1.4E+09	1.9E+04	Epoxybutane, 1,2-Ethephon	106-88-7 16672-87-0					7.0E+03		8.3E+01	8.2E+01
				2.0E-02	I				1.0E+00		1.5E+04	1.4E+09	7.7E+03	Ethion	563-12-2					6.7E+02			6.7E+02
				5.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Ethoxyethanol Acetate, 2-Ethyl Chloride (Chloroethane)	111-15-9 75-00-3					5.8E+03	1.4E+04		4.1E+03
				5.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Ethyl Acetate	563-12-2					5.8E+02	1.4E+03		4.1E+02
4.8E-02	H			1.0E-01	P	6.0E-02	P	V	1.0E+00		3.1E+04	1.4E+09	6.2E+04	Ethyl Acrylate	111-15-9	6.8E+01			6.8E+01	1.2E+05		1.6E+04	1.4E+04
				9.0E-02	P	2.0E-01	I	V	1.0E+00		1.1E+05	1.4E+09	9.8E+04	Ethyl Chloride (Chloroethane)	111-15-9					1.1E+05			8.6E+04
				9.0E-01	I	7.0E-02	P	V	1.0E+00		1.1E+04	1.4E+09	8.6E+03	Ethyl Ether	141-78-6					1.1E+06			2.6E+03
				5.0E-03	P	8.0E-03	P	V	1.0E+00		2.5E+03	1.4E+09	6.3E+03	Ethyl Methacrylate	140-88-5					5.8E+03			2.1E+02
									1.0E+00		2.1E+03	1.4E+09	1.3E+03	Ethyl-p-nitrophenyl Phosphonate	75-00-3					1.1E+05			5.7E+04
				2.0E-01	I			V	1.0E+00		1.0E+04	1.4E+09	3.1E+03	Ethylbenzene	60-29-7					2.3E+05			2.3E+05
				9.0E-02	H	3.0E-01	P	V	1.0E+00		1.1E+03	1.4E+09	5.8E+03	Ethylene Cyanohydrin	97-63-2					1.1E+05			7.1E+03
1.1E-02	C	2.5E-06	C	1.0E-05	I	1.0E+00	I	V	1.0E+00	1.0E-01		1.4E+09		Ethylene Glycol Monobutyl Ether	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	1.0E+00		4.8E+02	1.4E+09	5.7E+03	Ethylene Glycol	100-41-4					1.2E+05			2.0E+04
				7.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Ethylene Glycol Monobutyl Ether	109-78-4					8.2E+04	1.9E+05		5.7E+04
				9.0E-02	P			V	1.0E+00		1.9E+05	1.4E+09	1.8E+05	Ethylene Glycol	107-15-3					1.1E+05			1.1E+05
				2.0E+00	I	4.0E-01	C		1.0E+00	1.0E-01		1.4E+09		Ethylene Glycol Monobutyl Ether	107-21-1					2.3E+06	5.5E+06	2.4E+09	1.6E+06
				1.0E-01	I	1.6E+00	I		1.0E+00	1.0E-01		1.4E+09		Ethylene Oxide	111-78-2					1.2E+05	2.8E+05	9.5E+09	8.2E+04
3.1E-01	C	8.8E-05	C						1.0E+00		1.2E+05	1.4E+09	6.1E+03	Ethylene Thiourea	75-21-8	1.1E+01		8.5E-01	7.9E-01				8.0E+02
4.5E-02	C	1.3E-05	C	8.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Ethyleneimine	96-45-7	7.3E+01	1.7E+02	1.3E+06	5.1E+01	9.3E+01	2.2E+02		6.6E+01
6.5E+01	C	1.9E-02	C					V	1.0E+00		1.5E+05	1.4E+09	2.4E+04	Ethylphthalyl Ethyl Glycolate	151-56-4	5.0E-02		1.5E-02	1.2E-02				
				3.0E+00	I				1.0E+00	1.0E-01		1.4E+09		Express	84-72-0					3.5E+06	8.3E+06		2.5E+06
				8.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Fenamiphos	101200-48-0					9.3E+03	2.2E+04		6.6E+03
				2.5E-04	I				1.0E+00	1.0E-01		1.4E+09		Fenamiphos	22224-92-6					2.9E+			

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

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) <sup>-1</sup>	ke y	IUR (ug/m <sup>3</sup> -y) <sup>-1</sup>	ke y	RfD <sub>o</sub> (mg/kg-day)	ke y	RfC <sub>i</sub> (mg/m <sup>3</sup> -y)	ke y	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
				2.5E-02	I				1.0E+00	1.0E-01				Fenpropathrin	39515-41-8					2.9E+04	6.9E+04		2.1E+04	
				1.3E-02	I				1.0E+00	1.0E-01				Fluometuron	2164-17-2					1.5E+04	3.6E+04		1.1E+04	
				4.0E-02	C	1.3E-02	C		1.0E+00					Fluoride	16984-48-8					4.7E+04		7.7E+07	4.7E+04	
				6.0E-02	I	1.3E-02	C		1.0E+00					Fluorine (Soluble Fluoride)	7782-41-4					7.0E+04		7.7E+07	7.0E+04	
				8.0E-02	I				1.0E+00	1.0E-01				Fluridone	59756-60-4					9.3E+04	2.2E+05		6.6E+04	
				2.0E-02	I				1.0E+00	1.0E-01				Flurprimidol	56425-91-3					2.3E+04	5.5E+04		1.6E+04	
				6.0E-02	I				1.0E+00	1.0E-01				Flutolanil	66332-96-5					7.0E+04	1.7E+05		4.9E+04	
				1.0E-02	I				1.0E+00	1.0E-01				Fluvalinate	69409-94-5					1.2E+04	2.8E+04		8.2E+03	
				1.0E-01	I				1.0E+00	1.0E-01				Folpet	133-07-3					1.2E+05	2.8E+05		8.2E+04	
				1.9E-01	I				1.0E+00	1.0E-01				Fomesafen	72178-02-0				1.2E+01					
						2.0E-03	I		1.0E+00	1.0E-01				Fonofos	944-22-9	9.3E+02	2.2E+03			2.3E+03	5.5E+03		1.6E+03	
				1.3E-05	I	2.0E-01	I	9.8E-03	A	V				Formaldehyde	50-00-0			7.3E+01	7.3E+01	2.3E+05	2.3E+05	3.3E+03	3.3E+03	
				9.0E-01	P	3.0E-04	X	V	1.0E+00		1.1E+05	1.4E+09	9.3E+04	Formic Acid	64-18-6					1.1E+06		1.2E+02	1.2E+02	
				3.0E+00	I				1.0E+00	1.0E-01				Fosetyl-AL	39148-24-8					3.5E+06	8.3E+06		2.5E+06	
														Furans										
				1.0E-03	X		V		1.0E+00	3.0E-02		1.4E+09	2.0E+05	Dibenzofuran	132-64-9					1.2E+03	9.2E+03		1.0E+03	
				1.0E-03	I		V		1.0E+00	3.0E-02	6.2E+03	1.4E+09	2.6E+03	Furan	110-00-9					1.2E+03	9.2E+03		1.0E+03	
				9.0E-01	I	2.0E+00	I	V	1.0E+00	3.0E-02	1.7E+05	1.4E+09	1.2E+04	Tetrahydrofuran	109-99-9					1.1E+06	8.3E+06	1.1E+05	9.6E+04	
				3.8E+00	H				1.0E+00	1.0E-01		1.4E+09		Furazolidone	67-45-8	8.6E-01	2.0E+00		6.0E-01					
				1.5E+00	C	4.3E-04	C		1.0E+00		1.0E+04	1.4E+09	4.9E+04	Furfural	98-01-1					3.5E+03		1.1E+04	2.6E+03	
				3.0E-02	I	8.6E-06	C		1.0E+00	1.0E-01		1.4E+09		Furium	531-82-8	2.2E+00	5.2E+00	3.9E+04	1.5E+00					
									1.0E+00	1.0E-01		1.4E+09		Furmecyclox	60568-05-0	1.1E+02	2.6E+02	1.9E+06	7.7E+01					
				4.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Glufosinate, Ammonium	77182-82-2					4.7E+02	1.1E+03		3.3E+02	
						8.0E-05	C		1.0E+00	1.0E-01		1.4E+09		Glutaraldehyde	111-30-8					2.3E+05	2.3E+05	3.3E+03	3.3E+03	
				4.0E-04	I	1.0E-03	H	V	1.0E+00		1.1E+05	1.4E+09	7.3E+04	Glycidyl	785-34-4					4.7E+02		3.2E+02	1.9E+02	
				1.0E-01	I				1.0E+00	1.0E-01		1.4E+09		Glyphosate	1071-83-6					1.2E+05	2.8E+05		8.2E+04	
				3.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Goal	428(4)-03-3					3.5E+03	8.3E+03		2.5E+03	
				1.0E-02	X		V		1.0E+00		1.4E+09	1.5E+05		Guandine	113-00-8					1.2E+04			1.2E+04	
				2.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Guandine Chloride	50-01-1					2.3E+04	5.5E+04		1.6E+04	
				3.0E-03	A	1.0E-02	A		1.0E+00	1.0E-01		1.4E+09		Guthion	86-50-0					3.5E+03	8.3E+03	6.0E+07	2.5E+03	
				5.0E-05	I				1.0E+00	1.0E-01		1.4E+09		Haloxyp, Methyl	69906-40-2					5.8E+01	1.4E+02		4.1E+01	
				1.3E-02	I				1.0E+00	1.0E-01		1.4E+09		Harmony	79277-27-3					1.5E+04	3.6E+04		1.1E+04	
				5.0E-04	I		V		1.0E+00		4.8E+05		Heptachlor	76-44-8	7.3E-01		4.5E+00	6.3E-01		5.8E+02			5.8E+02	
				1.3E-05	I				1.0E+00		8.4E+05		Heptachlor Epoxide	1024-57-3					1.5E+01				1.5E+01	
				2.0E-03	I		V		1.0E+00		3.8E+05		Hexabromobenzene	87-82-1	3.6E-01		4.0E+00	3.3E-01		2.3E+03			2.3E+03	
				2.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BUL-153)	68631-49-2					2.3E+02	5.5E+02		1.6E+02	
				1.6E+00	I	4.6E-04	I	8.0E-04	I	V		1.0E+00	1.4E+09	6.8E+04	Hexachlorobenzene	118-74-1	2.0E+00	1.8E+00	9.6E-01	9.3E+02			9.3E+02	
				7.8E-02	I	2.2E-05	I	1.0E-03	P	V		1.0E+00	1.7E+01	1.4E+09	1.1E+04	Hexachlorobutadiene	87-68-3	4.2E+01	6.0E+00	5.3E+00	1.2E+03			1.2E+03
				6.3E+00	I	1.8E-03	I	8.0E-03	A			1.0E+00	1.0E-01	1.4E+09	Hexachlorocyclohexane, Alpha	519-84-6	5.2E-01	1.2E+00	9.3E+03	3.6E-01	9.3E+03	2.2E+04	6.6E+03	
				1.8E+00	I	5.3E-04	I		1.0E+00	1.0E-01		1.4E+09		Hexachlorocyclohexane, Beta-	319-85-7	1.8E+00	4.3E+00	3.1E+04	1.3E+00					
				1.1E+00	C	3.1E-04	C	3.0E-04	I		1.0E+00	4.0E-02	1.4E+09	Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	3.0E+00	1.8E+01	5.4E+04	2.5E+00	3.5E+02	2.1E+03		3.0E+02	
				1.8E+00	I	5.1E-04	C		1.0E+00	1.0E-01		1.4E+09		Hexachlorocyclohexane, Technical	608-73-1	1.8E+00	4.3E+00	3.3E+04	1.3E+00					
				6.0E-03	I	2.0E-04	I	V	1.0E+00		1.6E+01	1.4E+09	8.5E+03	Hexachlorocyclopentadiene	77-47-4					7.0E+03		7.5E+00	7.5E+00	
				7.0E-04	I	3.0E-02	I	V	1.0E+00		1.4E+09	8.0E+03		Hexachloroethane	67-72-1	8.2E+01		8.9E+00	8.0E+00	8.2E+02		1.1E+03	4.6E+02	
				3.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Hexachlorophene	70-30-4					3.5E+02	8.3E+02		2.5E+02	
				1.1E-01	I				1.0E+00	1.5E-02		1.4E+09		Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	3.0E+01	4.7E+02		2.8E+01	3.5E+03	5.5E+04		3.3E+03	
				4.0E-04	P				1.0E+00		5.2E+03	1.4E+09	3.0E+05	Hexamethylene Diisocyanate, 1,6-	822-06-0					4.7E+02	1.1E+03	1.3E+01	1.3E+01	
				6.0E-02	H	7.0E-01	I	V	1.0E+00		1.4E+02	1.4E+09	8.3E+02	Hexamethylphosphoramide	680-31-9					4.7E+02	1.1E+03		3.3E+02	
				2.0E+00	P				1.0E+00	1.0E-01		1.4E+09		Hexane, N-	110-54-3					7.0E+04		2.5E+03	2.5E+03	
				5.0E-03	I	3.0E-02	I	V	1.0E+00		3.3E+03	1.4E+09	1.3E+04	Hexanedioic Acid	124-04-9					2.3E+06	5.5E+06		1.6E+06	
				3.3E-02	I				1.0E+00	1.0E-01		1.4E+09		Hexanone, 2-	591-78-6					5.8E+03		1.7E+03	1.3E+03	
				3.0E+00	I	4.9E-03	I		1.0E+00	1.0E-01		1.4E+09		Hexazinone	51235-04-2					3.9E+04	9.1E+04		2.7E+04	
				3.0E+00	I	4.9E-03	I		1.0E+00			1.4E+09		Hydrazine	302-01-2	1.1E+00		3.4E+03	1.1E+00			1.8E+05	1.8E+05	
									1.0E+00			1.4E+09		Hydrazine Sulfate	10034-93-2	1.1E+00		3.4E+03	1.1E+00					
						2.0E-02	I	V	1.0E+00			1.4E+09		Hydrogen Chloride	7647-01-0							1.2E+08	1.2E+08	
						1.4E-02	C	V	1.0E+00			1.4E+09		Hydrogen Fluoride	7664-39-3					4.7E+04		8.3E+07	4.7E+04	
						2.0E-03	I	V	1.															

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Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) <sup>-1</sup>	ke y <sup>-1</sup>	IUR (ug/m <sup>3</sup> ) <sup>-1</sup>	ke y	RfD <sub>o</sub> (mg/kg-day)	ke y	RfC <sub>i</sub> (mg/m <sup>3</sup> )	ke y	ke y	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)
9.5E-04	I			2.0E-01 1.5E-02	I	2.0E+00	C			1.0E+00	1.0E-01		1.4E+09	4.2E+05	Isophorone Isopropanol	78-59-1 33820-53-0	3.4E+03	8.1E+03		2.4E+03	2.3E+05 1.8E+04	5.5E+05	1.2E+10	1.6E+05 1.8E+04
				2.0E+00 1.0E-01 5.0E-02	P	2.0E-01	P	V		1.0E+00	1.0E+00	1.1E+05	1.4E+09	2.8E+04	Isopropanol Isopropyl Methyl Phosphonic Acid Isoxaben	67-63-0 1832-54-8 82558-50-7					2.3E+06 1.2E+05 5.8E+04		2.4E+04	2.4E+04 8.2E+04 4.1E+04
				3.0E-01 7.5E-02 2.0E-03	A	3.0E-01	A	V		1.0E+00	1.0E-01		1.4E+09		JP-7 Kerb Lactofen	NA 23950-58-5 77501-63-4					8.8E+04 2.3E+03	2.1E+05 5.5E+03	1.8E+09	1.8E+09 6.2E+04 1.6E+03
5.0E-01 8.5E-03	C C	1.5E-01 1.2E-05	C C	2.0E-02	C	2.0E-04	C		M	2.5E-02			1.4E+09		<b>Lead Compounds</b> --Lead Chromate --Lead Phosphate	7758-97-6 7446-27-7	6.5E+00 3.8E+02		1.1E+02 1.4E+06	6.2E+00 3.8E+02	2.3E+04		1.2E+06	2.3E+04
2.8E-01 8.5E-03	C C	8.0E-05 1.2E-05	C C							1.0E+00	1.0E-01		1.4E+09		--Lead acetate --Lead and Compounds --Lead subacetate	301-04-2 7439-92-1 1335-32-6	1.2E+01 3.8E+02	2.8E+01	2.1E+05 1.4E+06	8.2E+00 2.7E+02				8.0E+02
				1.0E-07 2.0E-03 2.0E-03	I		V			1.0E+00	1.0E-01	2.4E+00	1.4E+09	1.9E+03	--Tetraethyl Lead Linuron Lithium	78-00-2 330-55-2 7439-93-2					1.2E-01 2.3E+03 2.3E+03	5.5E+03		1.6E+03 2.3E+03
				2.0E-01 5.0E-04 1.0E-02	I					1.0E+00	1.0E-01		1.4E+09		Londax MCPA MCPB	83055-99-6 94-74-6 94-81-5					2.3E+05 5.8E+02 1.2E+04	5.5E+05 1.4E+03 2.8E+04		1.6E+05 4.1E+02 8.2E+03
				1.0E-03 2.0E-02 1.0E-01	I					1.0E+00	1.0E-01		1.4E+09		MCPB Malathion Maleic Anhydride	93-65-2 121-75-5 108-31-6					1.2E+03 2.3E+04 1.2E+05	2.8E+03 5.5E+04 2.8E+05	4.2E+06	8.2E+02 1.6E+04 8.0E+04
				5.0E-01 1.0E-04 3.0E-02	I P H					1.0E+00	1.0E-01		1.4E+09		Maleic Hydrzade Malononitrile Mancozeb	123-33-1 109-77-3 8018-01-7					5.8E+05 1.2E+02 3.5E+04	1.4E+06 2.8E+02 8.3E+04		4.1E+05 8.2E+01 2.5E+04
				5.0E-03 1.4E-01 2.4E-02	I S	5.0E-05 5.0E-05	I I			1.0E+00	1.0E-01		1.4E+09		Maneb Manganese (Diet) Manganese (Non-diet)	12427-38-2 7439-96-5 7439-96-5					5.8E+03 2.8E+04	1.4E+04	3.0E+05	4.1E+03 2.6E+04
				9.0E-05 3.0E-02	H I					1.0E+00	1.0E-01		1.4E+09		Mephostolan Mepiquat Chloride	950-107-7 24307-26-4					1.1E+02 3.5E+04	2.5E+02 8.3E+04		7.4E+01 2.5E+04
				3.0E-04 1.0E-04	I I	3.0E-04 3.0E-04	S I	V		7.0E-02		3.1E+00	1.4E+09	3.0E+04	--Mercuric Chloride (and other Mercury salts) --Mercury (elemental) --Methyl Mercury	7487-94-7 7439-97-6 22967-92-6					3.5E+02 1.2E+02	1.8E+06 4.0E+01		3.5E+02 1.2E+02
				8.0E-05 3.0E-05 3.0E-05	I I I		V			1.0E+00	1.0E-01		1.4E+09	1.9E+06	--Phenylmercuric Acetate Merphos Merphos Oxide	62-38-4 160-50-5 88-48-8					9.3E+01 3.5E+01 3.5E+01	2.2E+02 8.3E+01		6.6E+01 3.5E+01 2.5E+01
				6.0E-02 1.0E-04 5.0E-05	I I I	3.0E-02	P P V			1.0E+00	1.0E-01	4.6E+03	1.4E+09	6.8E+03	Metabul Methacrylonitrile Methamidophos	57837-19-1 126-98-7 10265-92-6					7.0E+04 1.2E+02 5.8E+01	1.7E+05 1.4E+02	8.9E+02	4.9E+04 1.0E+02 4.1E+01
				2.0E+00 1.0E-03 2.5E-02	I I I	2.0E+01	I I V			1.0E+00	1.0E-01	1.1E+05	1.4E+09	2.9E+04	Methanol Methidathion Methomyl	67-56-1 950-37-8 16752-77-5					2.3E+06 1.2E+03 2.9E+04	2.8E+03 6.9E+04	2.5E+06	1.2E+06 8.2E+02 2.1E+04
4.9E-02	C	1.4E-05	C	5.0E-03 8.0E-03	I P	1.0E-03	P P V			1.0E+00	1.0E-01	1.4E+09	1.4E+09	1.2E+05	Methoxy-5-nitroaniline, 2-nitroxychlor Methoxyethanol Acetate, 2-	99-59-2 72-43-5 110-49-6	6.7E+01	1.6E+02	1.2E+06	4.7E+01	5.8E+03 9.3E+03	1.4E+04	5.4E+02	4.1E+03 5.1E+02
				5.0E-03 1.0E+00 3.0E-02	P X H	2.0E-02	I V P V			1.0E+00	1.0E-01	1.1E+05	1.4E+09	1.0E+05	Methoxyethanol, 2- Methyl Acetate Methyl Acrylate	109-86-4 79-20-9 96-33-3					5.8E+03 1.2E+06 3.5E+04	8.8E+03	6.1E+02	3.5E+03 1.2E+06 6.0E+02
				6.0E-01 1.0E-03 8.0E-02	I X H	5.0E+00	I P O I V			1.0E+00	1.0E+00	2.8E+04	1.4E+09	1.2E+04	Methyl Ethyl Ketone (2-Butanone) Methyl Hydrazine Methyl Isobutyl Ketone (4-methyl-2-pentanone)	78-93-3 60-34-4 108-10-1			1.9E+00	1.9E+00	7.0E+05 1.2E+03 9.3E+04	2.7E+05	1.4E+01	1.9E+05 1.4E+01 5.6E+04
				1.4E+00 2.5E-04	I I	7.0E-01	I V			1.0E+00	1.0E-01	2.4E+03	1.4E+09	6.3E+03	Methyl Isocyanate Methyl Methacrylate Methyl Parathion	624-83-9 80-62-6 298-00-0					1.6E+06 2.9E+02	6.9E+02	1.9E+04	1.9E+04 2.1E+02
				6.0E-02 9.9E-02	X C	4.0E-02	H V			1.0E+00	1.0E-01	3.9E+02	1.4E+09	1.2E+04	Methyl Phosphonic Acid Methyl Styrene (Mixed Isomers) Methyl methanesulfonate	993-13-5 25013-15-4 66-27-3	3.3E+01	7.8E+01	6.0E+05	2.3E+01	7.0E+04 7.0E+03	1.7E+05	2.1E+03	4.9E+04 1.6E+03
1.8E-03 9.0E-03	C P	2.6E-07 2.4E-03	C C	3.0E+00 2.0E-02	I X	3.0E+00	I V			1.0E+00	1.0E-01	8.9E+03	1.4E+09	4.9E+03	Methyl tert-Butyl Ether (MTBE) Methyl-1,4-benzenediamine dihydrochloride, 2- Methyl-5-Nitroaniline, 2-	1634-04-4 615-45-2 99-55-8	1.8E+03 3.6E+02	2.3E+02	2.1E+02	2.6E+02	3.5E+02 2.3E+04	8.3E+02 5.5E+04	6.4E+04	6.4E+04 2.5E+02 1.6E+04
8.3E+00 1.3E-01	C C	2.4E-03 3.7E-05	C C							1.0E+00	1.0E-01		1.4E+09		Methyl-N-nitro-N-nitrosoguanidine, N- Methylaniline Hydrochloride, 2-	70-25-7 636-21-5	3.9E+01 2.5E+01	9.3E-01	6.9E+03 4.5E+05	2.8E-01				

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

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

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1								
SFO (mg/kg-day) <sup>-1</sup>	ke y <sup>-1</sup>	IUR (ug/m <sup>3</sup> -y) <sup>-1</sup>	ke y <sup>-1</sup>	RfD <sub>o</sub> (mg/kg-day)	ke y <sup>-1</sup>	RfC <sub>o</sub> (mg/m <sup>3</sup> -y)	ke y <sup>-1</sup>	ke y <sup>-1</sup>	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
				1.0E-02	A					1.0E+00	1.0E-01		1.4E+09		Methylarsonic acid	124-58-3					1.2E+04	2.8E+04		8.2E+03	
1.0E-01	X			2.0E-04	X					1.0E+00	1.0E-01		1.4E+09		Methylbenzene, 1,4-diamine monohydrochloride, 2-	74612-12-7					2.3E+02	5.5E+02		1.6E+02	
2.2E+01	C	6.3E-03	C	3.0E-04	X				M	1.0E+00	1.0E-01		1.4E+09		Methylbenzene-1,4-diamine sulfate, 2-	615-50-9	3.3E+01	7.7E+01		2.3E+01	3.5E+02	8.3E+02		2.5E+02	
2.0E-03	I	1.0E-08	I	6.0E-03	I	6.0E-01	I	V	M	1.0E+00		3.3E+03	1.4E+09	2.2E+03	Methylcholanthrene, 3-	56-49-5	1.5E-01	3.5E-01	2.6E+03	1.0E-01					
1.0E-01	P	4.3E-04	C	2.0E-03	P				M	1.0E+00	1.0E-01		1.4E+09		Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	3.3E+01	7.7E+01	3.9E+04	2.3E+01	2.3E+03	5.5E+03	5.8E+03	3.2E+03	
4.6E-02	I	1.3E-05	C							1.0E+00	1.0E-01		1.4E+09		Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	7.1E+01	1.7E+02	1.3E+06	5.0E+01				1.6E+03	
1.6E+00	C	4.6E-04	C			2.0E-02	C			1.0E+00	1.0E-01		1.4E+09		Methylenediphenyl Diisocyanate	101-77-9	2.0E+00	4.8E+00	3.6E+04	1.4E+00			1.2E+08	1.2E+08	
				7.0E-02	H	6.0E-04	I		V	1.0E+00	1.0E-01		1.4E+09	1.3E+04	Methylstyrene, Alpha-	101-68-8					8.2E+04		3.6E+06	8.2E+04	
				1.5E-01	I					1.0E+00	1.0E-01		1.4E+09		Metolachlor	51218-45-2					1.8E+05	4.1E+05		1.2E+05	
				2.5E-02	I					1.0E+00	1.0E-01		1.4E+09		Metribuzin	21087-64-9					2.9E+04	6.9E+04		2.1E+04	
				3.0E+00	P				V	1.0E+00		3.4E-01	1.4E+09	1.4E+03	Mineral oils	8012-95-1					3.5E+06			3.5E+06	
1.8E+01	C	5.1E-03	C	2.0E-04	I			V		1.0E+00			1.4E+09	8.6E+05	Mirex	2385-85-5	1.8E-01		2.1E+00	1.7E-01				2.3E+02	
				2.0E-03	I					1.0E+00	1.0E-01		1.4E+09		Molinate	2212-67-1					2.3E+03	5.5E+03		1.6E+03	
				5.0E-03	I					1.0E+00			1.4E+09		Molybdenum	7439-98-7					5.8E+03			5.8E+03	
				1.0E-01	I					1.0E+00			1.4E+09		Monochloramine	10599-90-3					1.2E+05			1.2E+05	
				2.0E-03	P					1.0E+00	1.0E-01		1.4E+09		Monomethylaniline	100-61-8					2.3E+03	5.5E+03		1.6E+03	
				3.0E-04	X					1.0E+00	1.0E-01		1.4E+09		N,N-Diphenyl-1,4-benzenediamine	74-31-7					3.5E+02	8.3E+02		2.5E+02	
1.8E+00	C	0.0E+00	C	2.0E-03	I			V		1.0E+00			1.4E+09	5.7E+04	Naled	300-76-5					2.3E+03			2.3E+03	
				3.0E-02	X	1.0E-01	P	V		1.0E+00			1.4E+09		Naphtha, High Flash Aromatic (HFAN)	64742-95-6	1.8E+00	4.3E+00		1.3E+00	3.5E+04		6.0E+08		3.5E+04
				1.0E-01	I					1.0E+00	1.0E-01		1.4E+09		Napropamide	15299-99-7					1.2E+05	2.8E+05		8.2E+04	
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1.0E+00	1.0E-01		1.4E+09		Nickel Acetate	373-02-4		6.4E+04	6.4E+04	6.4E+04	1.3E+04	3.0E+04	8.3E+04		8.1E+03
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1.0E+00	1.0E-01		1.4E+09		Nickel Carbonate	3333-87-3		6.4E+04	6.4E+04	6.4E+04	1.3E+04	3.0E+04	8.3E+04		8.1E+03
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1.0E+00			1.4E+09		Nickel Carbonyl	13463-39-3		6.4E+04	6.4E+04	6.4E+04	1.3E+04			8.3E+04	1.1E+04
				2.6E-04	C	1.1E-02	C	1.4E-05	C	4.0E-02			1.4E+09		Nickel Hydroxide	12054-48-7		6.4E+04	6.4E+04	6.4E+04	1.3E+04			8.3E+04	1.1E+04
				2.6E-04	C	1.1E-02	C	2.0E-05	C	4.0E-02			1.4E+09		Nickel Oxide	1313-99-1		6.4E+04	6.4E+04	6.4E+04	1.3E+04			1.2E+05	1.2E+04
				2.4E-04	I	1.1E-02	C	1.4E-05	C	4.0E-02			1.4E+09		Nickel Refinery Dust	NA		6.9E+04	6.9E+04	6.9E+04	1.3E+04			8.3E+04	1.1E+04
				2.6E-04	C	2.0E-02	I	9.0E-05	A	4.0E-02			1.4E+09		Nickel Soluble Salts	7440-02-0	1.9E+00	6.4E+04	6.4E+04	6.4E+04	2.3E+04			4.5E+05	2.2E+04
1.7E+00	C	4.8E-04	I	1.1E-02	C	1.4E-05	C			4.0E-02			1.4E+09		Nickel Sulfide	12035-72-2		3.5E+04	1.9E+00	1.9E+00	1.3E+04			8.3E+04	1.1E+04
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1.0E+00	1.0E-01		1.4E+09		Nickelocene	1271-28-9		6.4E+04	6.4E+04	6.4E+04	1.3E+04	3.0E+04	8.3E+04		8.1E+03
				1.6E+00	I					1.0E+00			1.4E+09		Nitrate	14797-55-8					1.9E+06			1.9E+06	
										1.0E+00			1.4E+09		Nitrate + Nitrite (as N)	NA									
				1.0E-01	I					1.0E+00			1.4E+09		Nitrite	14797-85-0					1.2E+05				1.2E+05
				1.0E-02	X	5.0E-05	X			1.0E+00	1.0E-01		1.4E+09		Nitroamine, 2-	88-74-4					1.2E+04	2.8E+04	3.0E+05		8.0E+03
				4.0E-03	P	6.0E-03	P			1.0E+00	1.0E-01		1.4E+09		Nitroamine, 4-	100-01-6	1.6E+02	3.9E+02		1.1E+02	4.7E+03	1.1E+04	3.6E+07		3.3E+03
				4.0E-05	I	2.0E-03	I	9.0E-03	I	V		3.1E+03	1.4E+09	7.3E+04	Nitrobenzene	98-95-3			2.2E+01	2.2E+01	2.3E+03			2.9E+03	1.3E+03
				3.0E+03	P					1.0E+00	1.0E-01		1.4E+09		Nitrocellulose	9004-70-0					3.5E+09	8.3E+09			2.5E+09
				7.0E-02	H					1.0E+00	1.0E-01		1.4E+09		Nitrofurantoin	61-20-9					8.2E+04	1.9E+05			5.7E+04
1.3E+00	C	3.7E-04	C							1.0E+00	1.0E-01		1.4E+09		Nitrofurazone	59-87-0	2.5E+00	5.9E+00	4.5E+04	1.8E+00				8.2E+01	
1.7E-02	P			1.0E-04	P					1.0E+00	1.0E-01		1.4E+09		Nitroglycerin	55-63-0	1.9E+02	4.5E+02		1.4E+02	1.2E+02	2.8E+02			8.2E+01
				1.0E-01	I					1.0E+00	1.0E-01		1.4E+09		Nitroguanidine	556-88-7					1.2E+05	2.8E+05			8.2E+04
				8.8E-06	P	5.0E-03	P	V		1.0E+00		1.8E+04	1.4E+09	1.7E+04	Nitromethane	75-52-5			2.4E+01	2.4E+01				3.7E+02	3.7E+02
2.7E+01	C	7.7E-03	C	2.7E-03	H	2.0E-02	I	V		1.0E+00		4.9E+03	1.4E+09	1.3E+04	Nitropropane, 2	79-48-9	1.2E-01	2.9E-01	2.2E+03	8.5E-02				1.2E+03	1.2E+03
1.2E+02	C	3.4E-02	C						M	1.0E+00	1.0E-01		1.4E+09		Nitroso-N-ethylurea, N-	759-73-9			6.0E-02	6.0E-02					
5.4E+00	I	1.6E-03	I						V	1.0E+00			1.4E+09	2.4E+05	Nitroso-N-methylurea, N-	684-93-5	2.7E-02	6.4E-02	4.9E+02	1.9E-02					
7.0E+00	I	2.0E-03	C							1.0E+00	1.0E-01		1.4E+09		Nitroso-di-N-butylamine, N-	924-16-3	6.1E-01		1.9E+00	4.6E-01					
										1.0E+00	1.0E-01		1.4E+09		Nitroso-di-N-propylamine, N-	621-64-7	4.7E-01	1.1E+00	8.3E+03	3.3E-01					
2.8E+00	I	8.0E-04	C							1.0E+00	1.0E-01		1.4E+09		Nitrosodiethanolamine, N-	1116-54-7	1.2E+00	2.8E+00	2.1E+04	8.2E-01					
1.5E+02	I	4.3E-02	I						M	1.0E+00	1.0E-01		1.4E+09		Nitrosodiethylamine, N-	55-18-5	2.2E-02	5.2E-02	3.9E+02	1.5E-02					
5.1E+01	I	1.4E-02	I	8.0E-06	P	4.0E-05	X	V	M	1.0E+00		2.4E+05	1.4E+09	8.2E+04	Nitrosodimethylamine, N-	62-75-9	6.4E-02		7.2E-02	3.4E-02	9.3E+00		1.4E+01	5.7E+00	
4.9E-03	I	2.6E-06	C							1.0E+00	1.0E-01		1.4E+09		Nitrosodiphenylamine, N-	86-30-6	6.7E+02	1.6E+03	6.4E+06	4.7E+02					
2.2E+01	I	6.3E-03	C						V	1.0E+00		1.1E+05	1.4E+09	1.2E+05	Nitrosomethylethylamine, N-	10595-95-6	1.5E-01		2.4E-01	9.1E-02					
6.7E+00	C	1.9E-03	C																						



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

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1						
SFO (mg/kg-day) <sup>-1</sup>	ke y	IUR (ug/m <sup>3</sup> -y) <sup>-1</sup>	ke y	RfD <sub>o</sub> (mg/kg-day)	ke y	RfC <sub>i</sub> (mg/m <sup>3</sup> )	ke y	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)
				5.0E-02	I				1.0E+00	6.0E-03		1.4E+09		Octahydro-1,3,5,7-tetra-nitro-1,3,5,7-tetrazocine (HMX)	2691-41-0					5.8E+04	2.3E+06		5.7E+04
				2.0E-03	H				1.0E+00	1.0E-01		1.4E+09		Octamethylpyrophosphoramide	152-16-9					2.3E+03	5.5E+03		1.6E+03
				5.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Oryzalin	19044-88-3					5.8E+04	1.4E+05		4.1E+04
				5.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Oxadiazon	19666-30-9					5.8E+03	1.4E+04		4.1E+03
				2.5E-02	I				1.0E+00	1.0E-01		1.4E+09		Oxamyl	23135-22-0					2.9E+04	6.9E+04		2.1E+04
				1.3E-02	I				1.0E+00	1.0E-01		1.4E+09		Paclitaxel	76738-62-0					1.5E+04	3.6E+04		1.1E+04
				4.5E-03	I				1.0E+00	1.0E-01		1.4E+09		Paraquat Dichloride	1910-42-5					5.3E+03	1.2E+04		3.7E+03
				6.0E-03	H				1.0E+00	1.0E-01		1.4E+09		Parathion	56-38-2					7.0E+03	1.7E+04		4.9E+03
				5.0E-02	H			V	1.0E+00			4.5E+04		Pebulate	1114-71-2					5.8E+04			5.8E+04
				4.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Pendimethalin	40487-42-1					4.7E+04	1.1E+05		3.3E+04
				2.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Pentabromodiphenyl Ether	32534-81-9					2.3E+03	5.5E+03		1.6E+03
				1.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Pentabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-99)	60348-60-9					1.2E+02	2.8E+02		8.2E+01
				8.0E-04	I			V	1.0E+00			8.1E+04		Pentachlorobenzene	608-93-5					9.3E+02			9.3E+02
9.0E-02	P							V	1.0E+00		4.5E+02	1.4E+09	9.7E+03	Pentachloroethane	76-01-7	3.6E+01			3.6E+01				
2.6E-01	H			3.0E-03	I			V	1.0E+00			4.3E+05	Pentachloronitrobenzene	82-68-8	1.3E+01			1.3E+01		3.5E+03			3.5E+03
4.0E-01	I	5.1E-06	C	5.0E-03	I				1.0E+00	2.5E-01		1.4E+09	Pentachlorophenol	87-86-5	8.2E+00	7.7E+00	3.3E+06	4.0E+00	5.8E+03	5.5E+03		2.8E+03	
4.0E-03	X			2.0E-03	P				1.0E+00	1.0E-01		1.4E+09	Pentaerythritol tetranitrate (PETN)	78-11-5	8.2E+02	1.9E+03		5.7E+02	2.3E+03	5.5E+03		1.6E+03	
						1.0E+00	P	V	1.0E+00		3.9E+02	1.4E+09	7.8E+02	Pentane, n-	109-68-0							3.4E+03	3.4E+03
				7.0E-04	I				1.0E+00			1.4E+09		Perchlorates									
				7.0E-04	I				1.0E+00			1.4E+09		--Ammonium Perchlorate	7790-98-9					8.2E+02			8.2E+02
				7.0E-04	I				1.0E+00			1.4E+09		--Lithium Perchlorate	7791-03-9					8.2E+02			8.2E+02
				7.0E-04	I				1.0E+00			1.4E+09		--Perchlorate and Perchlorate Salts	14797-73-0					8.2E+02			8.2E+02
				7.0E-04	I				1.0E+00			1.4E+09		--Potassium Perchlorate	7778-74-7					8.2E+02			8.2E+02
				7.0E-04	I				1.0E+00			1.4E+09		--Sodium Perchlorate	1601-89-0					8.2E+02			8.2E+02
				2.0E-02	P			V	1.0E+00			1.4E+09	1.3E+05	Perfluorobutane Sulfonate	375-73-5					2.3E+04			2.3E+04
2.2E-03	C	6.3E-07	C	5.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Permethrin	52645-53-1	1.5E+03	3.5E+03	2.6E+07	1.0E+03	5.8E+04	1.4E+05		4.1E+04
				2.5E-01	I				1.0E+00	1.0E-01		1.4E+09		Phenmedipham	13684-63-4					2.9E+05	6.9E+05		2.1E+05
				3.0E-01	I	2.0E-01	C		1.0E+00	1.0E-01		1.4E+09		Phenol	108-95-2					3.5E+05	8.3E+05	1.2E+09	2.5E+05
				5.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Phenothiazine	92-84-2					5.8E+02	1.4E+03		4.1E+02
				6.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Phenylethylamine, m-	108-45-2					7.0E+03	1.7E+04		4.9E+03
				1.9E-01	H				1.0E+00	1.0E-01		1.4E+09		Phenylethylamine, o-	95-54-5	7.0E+01	1.6E+02		4.9E+01				
				1.9E-01	H				1.0E+00	1.0E-01		1.4E+09		Phenylethylamine, p-	106-50-3					2.2E+05	5.2E+05		1.6E+05
1.9E-03	H			2.0E-04	H				1.0E+00	1.0E-01		1.4E+09		Phenylphenol, 2-	90-43-7	1.7E+03	4.0E+03		1.2E+03				
				3.0E-04	I	V			1.0E+00	1.0E-01	1.6E+03	1.4E+09	9.8E+02	Phorate	298-02-2					2.3E+02	5.5E+02	1.3E+00	1.6E+02
				2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Phosgene	75-44-5								1.3E+00
				4.9E+01	P				1.0E+00			1.4E+09		Phosmet	732-11-6					2.3E+04	5.5E+04		1.6E+04
				4.9E+01	P				1.0E+00			1.4E+09		Phosphates, Inorganic									
				4.9E+01	P				1.0E+00			1.4E+09		--Aluminum metaphosphate	13176-88-0					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Ammonium polyphosphate	68333-79-9					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Calcium pyrophosphate	7790-76-3					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Diammonium phosphate	7783-28-0					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Dicalcium phosphate	7757-93-9					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Dimagnesium phosphate	7782-75-4					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Dipotassium phosphate	7758-11-4					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Disodium phosphate	7558-79-4					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monoaluminum phosphate	13530-50-2					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monoammonium phosphate	7722-76-1					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monocalcium phosphate	7758-23-8					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monomagnesium phosphate	7757-86-0					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monopotassium phosphate	7778-77-0					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Monosodium phosphate	7558-80-7					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Polyphosphoric acid	8017-16-1					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Potassium triphosphate	13845-36-8					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium acid pyrophosphate	7758-16-9					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium aluminum phosphate (acidic)	7785-88-8					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium aluminum phosphate (anhydrous)	10279-59-1					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium aluminum phosphate (tetrahydrate)	10305-76-7					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium hexametaphosphate	10124-56-8					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium polyphosphate	68915-31-1					5.7E+07			5.7E+07
				4.9E+01	P				1.0E+00			1.4E+09		--Sodium trimetaphosphate	7785-84-4					5.7E+07			

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

Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) <sup>-1</sup>	ke y	IUR (ug/m <sup>3</sup> -y) <sup>-1</sup>	ke y	RfD <sub>o</sub> (mg/kg-day)	ke y	RfC <sub>i</sub> (mg/m <sup>3</sup> )	ke y	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)
4.9E+01	P			4.9E+01	P				1.0E+00			1.4E+09		-Trialuminum sodium tetra decahydrogenoctaorthophosphate (dihydrate)	15136-87-5					5.7E+07			5.7E+07
4.9E+01	P			4.9E+01	P				1.0E+00			1.4E+09		-Tricalcium phosphate	7758-87-4					5.7E+07			5.7E+07
4.9E+01	P			4.9E+01	P				1.0E+00			1.4E+09		-Trimagnesium phosphate	7757-87-1					5.7E+07			5.7E+07
4.9E+01	P			4.9E+01	P				1.0E+00			1.4E+09		-Tripotassium phosphate	7778-53-2					5.7E+07			5.7E+07
4.9E+01	P			4.9E+01	P				1.0E+00			1.4E+09		-Trisodium phosphate	7801-54-9					5.7E+07			5.7E+07
3.0E-04	I	3.0E-04	I	3.0E-04	I	V			1.0E+00			1.4E+09		Phosphine	7803-51-2					3.5E+02		1.8E+06	3.5E+02
4.9E+01	P			4.9E+01	P				1.0E+00			1.4E+09		Phosphoric Acid	7664-38-2					5.7E+07		6.0E+07	2.9E+07
2.0E-05	I			2.0E-05	I	V			1.0E+00			6.9E+03		Phosphorus, White	7723-14-0					2.3E+01			2.3E+01
1.4E-02	I	2.4E-06	C	2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		<b>Phthalates</b>									
				1.0E+00	I				1.0E+00	1.0E-01		1.4E+09		-Bis(2-ethylhexyl)phthalate	117-81-7	2.3E+02	5.5E+02	6.9E+06	1.6E+02	2.3E+04	5.5E+04		1.6E+04
				1.0E-01	I				1.0E+00	1.0E-01		1.4E+09		-Butylphthalyl Butylglycolate	85-70-1					1.2E+06	2.8E+06		8.2E+05
				1.0E-01	I				1.0E+00	1.0E-01		1.4E+09		-Diethyl Phthalate	84-74-2					1.2E+05	2.8E+05		8.2E+04
				8.0E-01	I				1.0E+00	1.0E-01		1.4E+09		-Diethyl Phthalate	84-66-2					9.3E+05	2.2E+06		6.6E+05
				1.0E-01	I	V			1.0E+00			2.1E+04		-Dimethylterephthalate	120-61-6					1.2E+05			1.2E+05
				1.0E-02	P				1.0E+00	1.0E-01		1.4E+09		-Octyl Phthalate, di-N-	117-84-0					1.2E+04	2.8E+04		8.2E+03
				1.0E+00	H				1.0E+00	1.0E-01		1.4E+09		-Phthalic Acid, P-	100-21-0					1.2E+06	2.8E+06		8.2E+05
				2.0E+00	I	2.0E-02	C		1.0E+00	1.0E-01		1.4E+09		-Phthalic Anhydride	85-44-9					2.3E+06	5.5E+06	1.2E+08	1.6E+06
				7.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Picloram	1918-02-1					8.2E+04	1.9E+05		5.7E+04
				1.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3					1.2E+02	2.8E+02		8.2E+01
				1.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Pirimphos, Methyl	29232-93-7					1.2E+04	2.8E+04		8.2E+03
3.0E+01	C	8.6E-03	C	7.0E-06	H				1.0E+00	1.0E-01		1.4E+09		Polybrominated Biphenyls	59536-65-1	1.1E-01	2.6E-01	1.9E+03	7.7E-02	8.2E+00	1.9E+01		5.7E+00
7.0E-02	S	2.0E-05	S	7.0E-05	I		V		1.0E+00	1.4E-01		1.4E+09	5.9E+05	<b>Polychlorinated Biphenyls (PCBs)</b>									
									1.0E+00	1.4E-01		1.4E+09	1.1E+05	-Aroclor 1016	12674-11-2	4.7E+01	7.9E+01	3.6E+02	2.7E+01	8.2E+01	1.4E+02		5.1E+01
2.0E+00	S	5.7E-04	S				V		1.0E+00	1.4E-01		1.4E+09	1.1E+05	-Aroclor 1221	11104-28-2	1.6E+00	2.8E+00	2.4E+00	7.2E-01				
2.0E+00	S	5.7E-04	S				V		1.0E+00	1.4E-01		1.4E+09	1.1E+05	-Aroclor 1232	11141-16-5	1.6E+00	2.8E+00	2.4E+00	7.2E-01				
2.0E+00	S	5.7E-04	S				V		1.0E+00	1.4E-01		1.4E+09	7.9E+05	-Aroclor 1242	53469-21-9	1.6E+00	2.8E+00	1.7E+01	9.7E-01				
2.0E+00	S	5.7E-04	S				V		1.0E+00	1.4E-01		1.4E+09	5.1E+05	-Aroclor 1248	12872-29-6	1.6E+00	2.8E+00	1.1E+01	9.4E-01				
2.0E+00	S	5.7E-04	S	2.0E-05	I		V		1.0E+00	1.4E-01		1.4E+09	8.4E+05	-Aroclor 1254	11097-69-1	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				V		1.0E+00	1.4E-01		1.4E+09	1.3E+06	-Aroclor 1260	11098-82-5	1.6E+00	2.8E+00	2.8E+01	9.9E-01				
				6.0E-04	X		V		1.0E+00	1.4E-01		1.4E+09	7.2E+05	-Aroclor 5460	11126-42-4					7.0E+02	1.2E+03		4.4E+02
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1.0E+00	1.4E-01		1.4E+09	2.0E+06	-Heptachlorobiphenyl, 2,3,3',4,4',5,5'-(PCB 189)	39635-81-9	8.4E-01	1.4E+00	2.2E+01	5.1E-01	2.7E+01	4.6E+01	1.2E+04	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1.0E+00	1.4E-01		1.4E+09	1.4E+06	-Hexachlorobiphenyl, 2,3',4,4',5,5'-(PCB 167)	52663-72-6	8.4E-01	1.4E+00	1.5E+01	5.1E-01	2.7E+01	4.6E+01	8.4E+03	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1.0E+00	1.4E-01		1.4E+09	1.5E+06	-Hexachlorobiphenyl, 2,3,3',4,4',5,5'-(PCB 157)	69782-90-7	8.4E-01	1.4E+00	1.6E+01	5.1E-01	2.7E+01	4.6E+01	8.5E+03	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1.0E+00	1.4E-01		1.4E+09	1.5E+06	-Hexachlorobiphenyl, 2,3,3',4,4',5-(PCB 156)	38380-08-4	8.4E-01	1.4E+00	1.7E+01	5.1E-01	2.7E+01	4.6E+01	9.0E+03	1.7E+01
3.9E+03	E	1.1E+00	E	2.3E-08	E	1.3E-06	E	V	1.0E+00	1.4E-01		1.4E+09	1.4E+06	-Hexachlorobiphenyl, 3,3',4,4',5,5'-(PCB 169)	32174-16-6	8.4E-04	1.4E-03	1.5E-02	5.1E-04	2.7E-02	4.6E-02	8.4E+00	1.7E-02
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1.0E+00	1.4E-01		1.4E+09	1.0E+06	-Pentachlorobiphenyl, 2,3,4,4',5-(PCB 123)	65510-44-3	8.4E-01	1.4E+00	1.1E+01	5.0E-01	2.7E+01	4.6E+01	6.0E+03	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1.0E+00	1.4E-01		1.4E+09	8.3E+05	-Pentachlorobiphenyl, 2,3',4',4',5-(PCB 118)	31508-00-6	8.4E-01	1.4E+00	8.9E+00	5.0E-01	2.7E+01	4.6E+01	6.0E+03	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1.0E+00	1.4E-01		1.4E+09	8.5E+05	-Pentachlorobiphenyl, 2,3,3',4,4'-(PCB 105)	32598-14-4	8.4E-01	1.4E+00	9.1E+00	5.0E-01	2.7E+01	4.6E+01	4.9E+03	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	1.0E+00	1.4E-01		1.4E+09	1.0E+06	-Pentachlorobiphenyl, 2,3,4,4',5-(PCB 114)	44472-37-0	8.4E-01	1.4E+00	1.1E+01	5.0E-01	2.7E+01	4.6E+01	6.0E+03	1.7E+01
1.3E+04	E	3.8E+00	E	7.0E-09	E	4.0E-07	E	V	1.0E+00	1.4E-01		1.4E+09	1.0E+06	-Pentachlorobiphenyl, 3,3',4,4',5-(PCB 126)	57465-28-8	2.5E-04	4.2E-04	3.3E-03	1.5E-04	8.2E-03	1.4E-02	1.8E+00	5.1E-03
2.0E+00	I	5.7E-04	I				V		1.0E+00	1.4E-01		1.4E+09	7.9E+05	-Polychlorinated Biphenyls (high risk)	1336-36-3	1.6E+00	2.8E+00	1.7E+01	9.7E-01				
4.0E-01	I	1.0E-04	I				V		1.0E+00	1.4E-01		1.4E+09	1.4E+05	-Polychlorinated Biphenyls (low risk)	1336-36-3								
7.0E-02	I	2.0E-05	I				V		1.0E+00	1.4E-01		1.4E+09	1.4E+05	-Polychlorinated Biphenyls (lowest risk)	1336-36-3								
1.3E+01	E	3.8E-03	E	7.0E-06	E	4.0E-04	E		1.0E+00	1.4E-01		1.4E+09		-Tetrachlorobiphenyl, 3,3',4,4'-(PCB 77)	32598-13-3	2.5E-01	4.2E-01	4.4E+03	1.6E-01	8.2E+00	1.4E+01	2.4E+06	5.1E+00
3.9E+01	E	1.1E-02	E	2.3E-06	E	1.3E-04	E	V	1.0E+00	1.4E-01		1.4E+09	7.3E+05	-Tetrachlorobiphenyl, 3,4,4',5-(PCB 81)	70382-50-4	8.4E-02	1.4E-01	7.8E-01	4.9E-02	2.7E+00	4.6E+00	4.2E+02	1.7E+00
				6.0E-04	I		V		1.0E+00	1.0E-01		1.4E+09		Polymers: Methylenedioxybenzylidene Diisocyanate (MDI)	9016-87-9								3.6E+06
				6.0E-02	I		V		1.0E+00	1.3E-01		1.4E+09	1.4E+05	<b>Polynuclear Aromatic Hydrocarbons (PAHs)</b>									
				3.0E-01	I		V		1.0E+00	1.3E-01		1.4E+09	5.2E+05	-Acenaphthene	83-32-9					3.0E+04	1.3E+05		4.5E+04
7.3E-01	E	1.1E-04	C				V	M	1.0E+00	1.3E-01		1.4E+09	4.4E+06	-Anthracene	120-12-7					7.5E+05	6.4E+05		2.3E+05
							V	M	1.0E+00	1.3E-01		1.4E+09	4.4E+06	-Benz[a]anthracene	56-55-3	4.5E+00	8.1E+00	4.9E+02	2.9E+00				
1.2E+00	C	1.1E-04	C				V		1.0E														

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Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) <sup>-1</sup>	ke (y)	IUR (ug/m <sup>3</sup> ) <sup>-1</sup>	ke (y)	RfD <sub>o</sub> (mg/kg-day)	ke (y)	RfC <sub>i</sub> (mg/m <sup>3</sup> )	ke (y)	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)
3.4E-05	C	2.0E-02	I	3.0E-03	I	V	1.0E+00	1.3E-01	1.0E+00	1.3E-01	1.4E+09	4.6E+04	-Naphthalene	91-20-3		1.7E+01	1.7E+01		2.3E+04	4.2E+04	6.1E+02	5.9E+02	
1.2E+00	C	1.1E-04	C	3.0E-02	I	V	1.0E+00	1.3E-01	1.0E+00	1.3E-01	1.4E+09	2.4E+06	-Nitropyrene, 4- -Pyrene Potassium Perfluorobutane Sulfonate	57835-92-4 129-00-0 29420-49-3	2.7E+00	5.0E+00	1.5E+05	1.8E+00	3.5E+04	6.4E+04		2.3E+04	1.6E+04
1.5E-01	I	9.0E-03	I	6.0E-03	H	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	4.2E+05	Prochloraz Profuralin Prometon	67747-09-5 26399-36-0 1610-18-0	2.2E+01	5.2E+01		1.5E+01	1.1E+04	2.5E+04		7.4E+03	7.0E+03
4.0E-03	I	1.3E-02	I	5.0E-03	I		1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Prometryn Propachlor Propanil	7287-19-6 1918-16-7 709-98-8					4.7E+03	1.1E+04		3.3E+03	1.1E+04
2.0E-02	I	2.0E-03	I	2.0E-02	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	6.3E+04	Propargite Propargyl Alcohol Propazine	2312-35-8 107-19-7 139-40-2					2.3E+04	5.5E+04		1.6E+04	2.3E+03
2.0E-02	I	1.3E-02	I	8.0E-03	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	8.9E+03	Propam Propiconazole Propionaldehyde	122-42-9 60207-90-1 123-38-6					2.3E+04	5.5E+04		1.6E+04	1.1E+04
1.0E-01	X	1.0E+00	X	3.0E+00	C	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	2.6E+02	7.0E+03	Propyl benzene Propylene Propylene Glycol	103-65-1 115-07-1 57-55-6					1.2E+05	3.1E+04		2.4E+04	9.3E+03
2.4E-01	I	3.7E-06	I	3.0E-02	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	7.8E+04	1.0E+04	Propylene Glycol Dinitrate Propylene Glycol Monoethyl Ether Propylene Glycol Monomethyl Ether	6423-43-4 1569-02-4 107-98-2					8.2E+05	8.2E+05		1.6E+06	8.2E+05
2.4E-01	I	3.7E-06	I	3.0E-02	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	7.8E+04	1.0E+04	Propylene Oxide Pursult Pydrin	75-56-9 81335-77-5 51630-58-1	1.4E+01		3.4E+01	9.7E+00	2.9E+05	2.9E+04		1.4E+03	2.1E+05
3.0E+00	I	1.0E-03	I	5.0E-04	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	5.3E+05	5.5E+04	Pyridine Quinalphos Quindline	110-86-1 3593-03-8 91-22-5	1.1E+00	2.6E+00		7.7E-01	1.2E+03	5.8E+02	1.4E+03		1.2E+03
2.2E-01	C	6.3E-05	C	3.0E-02	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	4.7E+05	Refractory Ceramic Fibers Resmethrin Ronnel	NA 10463-86-8 299-84-3					3.5E+04	8.3E+04		1.8E+08	2.5E+04
2.2E-01	C	6.3E-05	C	4.0E-03	I		1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Rotenone Safrole Savay	83-79-4 94-59-7 78587-05-0	1.5E+01	3.5E+01	2.6E+05	1.0E+01	4.7E+03	1.1E+04		3.3E+03	2.1E+04
5.0E-03	I	5.0E-03	I	5.0E-03	I	C	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Selenious Acid Selenium Selenium Sulfide	7783-00-8 7782-49-2 446-34-6					5.8E+03	5.8E+03		1.2E+08	5.8E+03
1.2E-01	H	5.0E-03	I	1.3E-02	I		1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Sethoxydim Silica (crystalline, respirable) Silver	74051-80-2 7631-86-9 1440-22-4					1.1E+05	2.5E+05		1.8E+07	5.8E+03
5.0E-01	C	1.5E-01	C	2.0E-02	C	M	2.5E-02	1.0E+00	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Simazine Sodium Acifluorfen Sodium Azide	122-34-9 62476-59-9 26628-22-8	2.7E+01	6.4E+01		1.9E+01	5.8E+03	1.4E+04		1.1E+04	4.7E+03
2.7E-01	H	5.0E-02	A	3.0E-02	I	C	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Sodium Dichromate Sodium Diethyldithiocarbamate Sodium Fluoride	10509-01-0 148-18-5 7681-49-4	6.5E+00	1.1E+02		6.2E+00	2.3E+04	3.5E+04	8.3E+04	1.2E+06	2.3E+04
2.4E-02	H	3.0E-02	I	2.0E-05	I		1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Sodium Fluoroacetate Sodium Metavanadate Stirofos (Tetrachlorovinphos)	62-74-8 13718-26-8 961-11-5	1.4E+02	3.2E+02		9.6E+01	2.3E+01	5.5E+01		1.6E+01	1.2E+03
5.0E-01	C	1.5E-01	C	2.0E-02	C	M	2.5E-02	1.0E+00	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Strontium Chromate Strontium, Stable Strychnine	7789-06-2 7440-24-6 57-24-9	6.5E+00		1.1E+02	6.2E+00	2.3E+04	7.0E+05	8.3E+02	1.2E+06	2.3E+04
3.0E-03	P	1.0E-03	P	2.0E-01	I	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	8.7E+02	9.4E+03	Styrene Styrene-Acrylonitrile (SAN) Trimer Sulfolane	100-42-5 NA 126-33-0					2.3E+05	8.3E+03	4.1E+04	3.5E+04	
8.0E-04	P	1.0E-03	C	1.0E-03	C	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Sulfonylbis(4-chlorobenzene), 1,1'- Sulfur Trioxide Sulfuric Acid	80-07-9 7446-11-9 7664-93-9					9.3E+02	2.2E+03		6.0E+06	6.0E+06
3.0E-02	H	7.0E-02	I	3.0E-02	I		1.0E+00	1.0E-01	1.0E+00	1.0E-01	1.4E+09	1.4E+09	Systhane TCMTB Tebuthiuron	88671-89-0 21564-17-0 34014-18-1					2.9E+04	6.9E+04		2.1E+04	2.5E+04
2.0E-02	H	1.3E-02	I	2.5E-05	H	V	1.0E+00	1.0E-01	1.0E+00	1.0E-01	3.1E+01	2.6E+05	Temephos Terbacil Terbufos	3383-96-8 5902-51-2 13071-79-9					2.3E+04	5.5E+04		1.8E+04	1.1E+04

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

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Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1								
SFO (mg/kg-day) <sup>-1</sup>	ke y	IUR (ug/m <sup>3</sup> -y) <sup>-1</sup>	ke y	RfD <sub>o</sub> (mg/kg-day)	ke y	RfC <sub>i</sub> (mg/m <sup>3</sup> -y)	ke y	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
				1.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Terbutryn	886-50-0					1.2E+03	2.8E+03		8.2E+02	
				1.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47)	5436-43-1					1.2E+02	2.8E+02		8.2E+01	
				3.0E-04	I			V	1.0E+00			1.4E+09	5.1E+04	Tetrachlorobenzene, 1,2,4,5-	95-94-3					3.5E+02			3.5E+02	
2.6E-02	I	7.4E-06	I	3.0E-02	I			V	1.0E+00		6.8E+02	1.4E+09	5.7E+03	Tetrachloroethane, 1,1,1,2-	630-20-6	1.3E+02		9.4E+00	8.8E+00	3.5E+04			3.5E+04	
2.0E-01	I	5.8E-05	C	2.0E-02	I			V	1.0E+00		1.9E+03	1.4E+09	1.5E+04	Tetrachloroethane, 1,1,1,2-	79-34-5	1.6E+01		3.2E+00	2.7E+00	2.3E+04			2.3E+04	
2.1E-03	I	2.6E-07	I	6.0E-03	I	4.0E-02	I	V	1.0E+00		1.7E+02	1.4E+09	2.4E+03	Tetrachloroethylene	127-18-4	1.6E+03		1.1E+02	1.0E+02	7.0E+03		4.1E+02	3.9E+02	
2.0E+01	H			3.0E-02	I			V	1.0E+00	1.0E-01		1.4E+09		Tetrachlorophenol, 2,3,4,6-	58-90-2				1.6E-01		8.3E+04		2.5E+04	
				5.0E-04	I			V	1.0E+00	1.0E-01		1.4E+09	1.1E+05	Tetrachlorolouene, p- alpha, alpha- Tetraethyl Dithiopyrophosphate	5216-25-1 3689-24-5				1.6E-01		5.8E+02	1.4E+03		4.1E+02
				8.0E+01	I			V	1.0E+00		1.1E+03	1.4E+09	1.2E+03	Tetrafluoroethane, 1,1,1,2-	811-97-2							4.3E+05		4.3E+05
				2.0E-03	P				1.0E+00	6.5E-04		1.4E+09		Tetryl (Trinitrophenylmethylnitramine)	479-45-8					2.3E+03	8.5E+05		2.3E+03	
				7.0E-06	X				1.0E+00			1.4E+09		Thallium (I) Nitrate	10102-45-1					8.2E+00			8.2E+00	
				1.0E-05	X				1.0E+00			1.4E+09		Thallium (Soluble Salts)	7440-28-0					1.2E+01			1.2E+01	
				6.0E-06	X			V	1.0E+00	1.0E-01		1.4E+09		Thallium Acetate	563-68-8					7.0E+00	1.7E+01		4.9E+00	
				2.0E-05	X				1.0E+00	1.0E-01		1.4E+09		Thallium Carbonate	6533-73-9					2.3E+01	5.5E+01		1.6E+01	
				6.0E-06	X				1.0E+00			1.4E+09		Thallium Chloride	7791-12-0					7.0E+00			7.0E+00	
				2.0E-05	X				1.0E+00			1.4E+09		Thallium Sulfate	7446-18-6					2.3E+01			2.3E+01	
				1.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Thiobencarb	28249-77-6					1.2E+04	2.8E+04		8.2E+03	
				7.0E-02	X				1.0E+00	7.5E-03		1.4E+09		Thiodiglycol	111-48-8					8.2E+04	2.6E+06		7.9E+04	
				3.0E-04	H				1.0E+00	1.0E-01		1.4E+09		Thiofanox	39196-18-4					3.5E+02	8.3E+02		2.5E+02	
				8.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Thiophanate, Methyl	23564-05-8					9.3E+04	2.2E+05		6.6E+04	
				5.0E-03	I				1.0E+00	1.0E-01		1.4E+09		Thiram	137-26-8					5.8E+03	1.4E+04		4.1E+03	
				6.0E-01	H				1.0E+00			1.4E+09		Tin	7440-31-5					7.0E+05			7.0E+05	
				1.0E-04	A	V			1.0E+00			1.4E+09		Titanium tetrachloride	7550-45-0							6.0E+05	6.0E+05	
				8.0E-02	I	5.0E+00	I	V	1.0E+00		8.2E+02	1.4E+09	4.3E+03	Toluene	108-88-3					9.3E+04		9.4E+04	4.7E+04	
1.8E-01	X			2.0E-04	X				1.0E+00	1.0E-01		1.4E+09		Toluene, 2,6-diamine	95-70-5	1.8E+01	4.3E+01		1.3E+01	2.3E+02	5.5E+02		1.6E+02	
3.0E-02	P			4.0E-03	X				1.0E+00	1.0E-01		1.4E+09		Toluene, p-	106-49-0	1.1E+02	2.6E+02		7.7E+01	4.7E+03	1.1E+04		3.3E+03	
				3.0E+00	P			V	1.0E+00		3.4E-01	1.4E+09	1.1E+03	Total Petroleum Hydrocarbons (Aliphatic High)	NA					3.5E+06			3.5E+06	
				6.0E-01	P	V			1.0E+00		1.4E+02	1.4E+09	8.3E+02	Total Petroleum Hydrocarbons (Aliphatic Low)	NA						2.2E+03		2.2E+03	
				1.0E-02	X	1.0E-01	P	V	1.0E+00		6.9E+00	1.4E+09	1.0E+03	Total Petroleum Hydrocarbons (Aliphatic Medium)	NA					1.2E+04		4.6E+02	4.4E+02	
				4.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Total Petroleum Hydrocarbons (Aromatic High)	NA					4.7E+04	1.1E+05		3.3E+04	
				4.0E-03	P	3.0E-02	P	V	1.0E+00		1.8E+03	1.4E+09	3.5E+03	Total Petroleum Hydrocarbons (Aromatic Low)	NA					4.7E+03		4.6E+02	4.2E+02	
				4.0E-03	P	3.0E-03	P	V	1.0E+00		1.4E+09	5.2E+04		Total Petroleum Hydrocarbons (Aromatic Medium)	NA				4.7E+03		6.9E+02		6.0E+02	
1.1E+00	I	3.2E-04	I						1.0E+00	1.0E-01		1.4E+09		Toxaphene	8001-35-2	3.0E+00	7.0E+00	5.2E+04	2.1E+00					
				7.5E-03	I				1.0E+00	1.0E-01		1.4E+09		Tralometrin	60841-25-6					8.8E+03	2.1E+04		6.2E+03	
				3.0E-04	A			V	1.0E+00			1.4E+09	3.4E+03	Tr-n-butyltin	888-73-3					3.5E+02			3.5E+02	
				8.0E+01	X				1.0E+00	1.0E-01		1.4E+09		Triacetin	102-76-1					9.3E+07	2.2E+08		6.6E+07	
				1.3E-02	I			V	1.0E+00			1.4E+09	3.6E+05	Triallate	2303-17-5					1.5E+04			1.5E+04	
				1.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Trinasturon	82097-50-5					1.2E+04	2.8E+04		8.2E+03	
				5.0E-03	I			V	1.0E+00			1.4E+09	4.5E+04	Tribromobenzene, 1,2,4-	615-54-3				2.6E+02	5.8E+03			5.8E+03	
9.0E-03	P			1.0E-02	P				1.0E+00	1.0E-01		1.4E+09		Tributyl Phosphate	126-73-8	3.6E+02	8.6E+02			1.2E+04	2.8E+04		8.2E+03	
				3.0E-04	P				1.0E+00	1.0E-01		1.4E+09		Tributyltin Compounds	NA					3.5E+02	8.3E+02		2.5E+02	
				3.0E-04	I				1.0E+00	1.0E-01		1.4E+09		Tributyltin Oxide	56-35-9					3.5E+02	8.3E+02		2.5E+02	
7.0E-02	I			3.0E+01	I	3.0E+01	H	V	1.0E+00		9.1E+02	1.4E+09	1.3E+03	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1 76-03-9	4.7E+01	1.1E+02		3.3E+01	3.5E+07		1.7E+05	1.7E+05	
2.9E-02	H			2.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Trichloroacetic Acid	76-03-9	4.7E+01	1.1E+02		3.3E+01	2.3E+04	5.5E+04		1.6E+04	
7.0E-03	X			3.0E-05	X				1.0E+00	1.0E-01		1.4E+09		Trichloroaniline HCl, 2,4,6-	33663-50-2	1.1E+02	2.7E+02		7.9E+01	3.5E+01	8.3E+01		2.5E+01	
				8.0E-04	X				1.0E+00			1.4E+09	3.2E+04	Trichloroaniline, 2,4,6-	634-93-5 87-61-6	4.7E+02	1.1E+03		3.3E+02	9.3E+02			9.3E+02	
2.9E-02	P			1.0E-02	I	2.0E-03	P	V	1.0E+00		4.0E+02	1.4E+09	3.0E+04	Trichlorobenzene, 1,2,4-	120-82-1	1.1E+02			1.1E+02	1.2E+04		2.6E+02	2.6E+02	
				2.0E+00	I	5.0E+00	I	V	1.0E+00		6.4E+02	1.4E+09	1.7E+03	Trichloroethane, 1,1,1-	71-55-6					2.3E+06			3.6E+04	
5.7E-02	I	1.6E-05	I	4.0E-03	I	2.0E-04	X	V	1.0E+00		2.2E+03	1.4E+09	7.2E+03	Trichloroethane, 1,1,2-	79-00-5	5.7E+01		5.5E+00	5.0E+00	4.7E+03			6.3E+00	
4.6E-02	I	4.1E-06	I	5.0E-04	I	2.0E-03	I	V	1.0E+00		6.9E+02	1.4E+09	2.2E+03	Trichloroethylene	79-01-6	7.1E+01		6.6E+00	6.0E+00	5.8E+02		1.9E+01	1.9E+01	
				3.0E-01	I	7.0E-01	H	V	1.0E+00		1.2E+03	1.4E+09	1.0E+03	Trichlorofluoromethane	75-69-4					3.5E+05		3.2E+03	3.1E+03	
				1.0E-01	I				1.0E+00	1.0E-01		1.4E+09		Trichlorophenol, 2,4,5-	95-95-4					1.2E+05	2.8E+05		8.2E+04	
1.1E-02	I	3.1E-06	I	1.0E-03	P				1.0E+00	1.0E-01		1.4E+09		Trichlorophenol, 2,4,6-	88-06-2	3.0E+02	7.0E+02	5.4E+06	2.1E+02	1.2E+03	2.8E+03		8.2E+02	
				1.0E-02	I				1.0E+00	1.0E-01		1.4E+09		Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5					1.2E+04	2.8E+04		8.2E+03	
				8.0E-03	I																			

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

Toxicity and Chemical-specific Information											Contaminant		Carcinogenic Target Risk (TR) = 1E-06				Noncancer Hazard Index (HI) = 1								
SFO (mg/kg-day) <sup>-1</sup>	Key	IUR (ug/m <sup>3</sup> ) <sup>-1</sup>	Key	RfD <sub>o</sub> (mg/kg-day)	Key	RfC <sub>i</sub> (mg/m <sup>3</sup> )	Key	Vol	muta- gen	GIABS	ABS	C <sub>sat</sub> (mg/kg)	PEF (m <sup>3</sup> /kg)	VF (m <sup>3</sup> /kg)	Analyte	CAS No.	Ingestion SL TR=1.0E-6 (mg/kg)	Dermal SL TR=1.0E-6 (mg/kg)	Inhalation SL TR=1.0E-6 (mg/kg)	Carcinogenic SL TR=1.0E-6 (mg/kg)	Ingestion SL HQ=1 (mg/kg)	Dermal SL HQ=1 (mg/kg)	Inhalation SL HQ=1 (mg/kg)	Noncarcinogenic SL HI=1 (mg/kg)	
7.7E-03 2.0E-02	I P			7.5E-03 1.0E-02	I P			V		1.0E+00 1.0E+00	1.0E-01	1.4E+09 1.4E+09	5.1E+05 1.4E+09		Trifluralin Trimethyl Phosphate	1582-09-8 512-56-1	4.2E+02 1.6E+02			4.2E+02 1.1E+02	8.8E+03 1.2E+04			2.8E+04	8.8E+03 8.2E+03
				5.0E-03 7.0E-03 1.0E-02	P V X			V		1.0E+00 1.0E+00 1.0E+00		2.9E+02 2.2E+02 1.8E+02	1.4E+09 1.4E+09 1.4E+09	9.4E+03 7.9E+03 6.6E+03	Trimethylbenzene, 1,2,3- Trimethylbenzene, 1,2,4- Trimethylbenzene, 1,3,5-	526-73-8 95-63-6 108-67-8					1.2E+04		2.1E+02 2.4E+02	2.1E+02 2.4E+02	
3.0E-02	I			3.0E-02 5.0E-04 2.0E-02	I I P					1.0E+00 1.0E+00 1.0E+00	1.9E-02 3.2E-02 1.0E-01	1.4E+09 1.4E+09 1.4E+09			Trinitrobenzene, 1,3,5- Trinitrotoluene, 2,4,6- Triphenylphosphine Oxide	99-35-4 118-96-7 791-28-6		1.1E+02 8.0E+02		9.6E+01	3.5E+04 5.8E+02 2.3E+04	4.4E+05 4.3E+03 5.5E+04		3.2E+04 5.1E+02 1.6E+04	
2.3E+00	C	6.6E-04	C	2.0E-02 1.0E-02	A X			V		1.0E+00 1.0E+00	1.0E-01 1.0E-01	1.4E+09 1.4E+09	9.0E+05		Tris(1,3-Dichloro-2-propyl) Phosphate Tris(1-chloro-2-propyl)phosphate Tris(2,3-dibromopropyl)phosphate	13674-87-8 13674-84-5 126-72-1	1.4E+00		1.7E+01	1.3E+00	2.3E+04 1.2E+04	5.5E+04 2.8E+04		1.6E+04 8.2E+03	
2.0E-02 3.2E-03	P P			7.0E-03 1.0E-01 3.0E-03	P P I	7.0E-06 4.0E-05				1.0E+00 1.0E+00 1.0E+00	1.0E-01 1.0E-01	1.4E+09 1.4E+09 1.4E+09			Tris(2-chloroethyl)phosphate Tris(2-ethylhexyl)phosphate Uranium (Soluble Salts)	115-96-3 78-42-2 NA	1.6E+02 1.0E+03	3.9E+02 2.4E+03		1.1E+02 7.2E+02	8.2E+03 1.2E+05 3.5E+03	1.9E+04 2.8E+05		5.7E+03 8.2E+04 3.5E+03	
1.0E+00	C	2.9E-04 8.3E-03	C P	9.0E-03 5.0E-03	I S	7.0E-06 1.0E-04			M	1.0E+00 2.6E-02 2.6E-02	1.0E-01	1.4E+09 1.4E+09 1.4E+09			Urethane Vanadium Pentoxide Vanadium and Compounds	51-79-6 1314-62-1 7440-62-2	3.3E+00		5.7E+04 2.0E+03	2.3E+00 2.0E+03	1.1E+04 5.9E+03		4.2E+04 6.0E+05	8.4E+03 5.8E+03	
				1.0E-03 2.5E-02 1.0E+00	I I H			V		1.0E+00 1.0E+00 1.0E+00	1.0E-01	1.4E+09 1.4E+09 2.8E+03	1.2E+05 1.4E+09 4.4E+03		Vermolate Vinclozolin Vinyl Acetate	1929-77-7 50471-44-8 109-05-4					1.2E+03 2.9E+04 1.2E+06	6.9E+04		1.2E+03 2.1E+04 3.8E+03	
7.2E-01	I	4.4E-06	I	3.0E-03 3.0E-04	I I	1.0E-01		V M		1.0E+00 1.0E+00	1.0E-01	3.4E+03 3.9E+03	1.4E+09 1.4E+09	1.4E+03 9.6E+02	Vinyl Bromide Vinyl Chloride Warfarin	593-60-2 75-01-4 81-81-2	4.5E+00		5.2E-01 2.7E+00	5.2E-01 1.7E+00	3.5E+03 3.5E+02		1.8E+01 4.2E+02 8.3E+02	1.8E+01 3.7E+02 2.5E+02	
				2.0E-01 2.0E-01 2.0E-01	S S S	1.0E-01		S V V		1.0E+00 1.0E+00 1.0E+00		3.9E+02 3.9E+02 4.3E+02	1.4E+09 1.4E+09 6.5E+03	5.6E+03 5.5E+03	Xylene, p- Xylene, m- Xylene, o-	106-42-3 108-38-3 95-47-6				2.3E+05 2.3E+05 2.3E+05		2.4E+03 2.4E+03 2.8E+03	2.4E+03 2.4E+03 2.8E+03		
				2.0E-01 3.0E-04 3.0E-01	I I I	1.0E-01		V		1.0E+00 1.0E+00 1.0E+00		2.6E+02 1.4E+09 1.4E+09	6.5E+03		Xylenes Zinc Phosphide Zinc and Compounds	1330-20-7 1314-84-7 7440-66-6				2.3E+05 3.5E+02 3.5E+05		2.8E+03	2.8E+03	2.8E+03 3.5E+02 3.5E+05	
				5.0E-02 8.0E-05	I X					1.0E+00 1.0E+00	1.0E-01	1.4E+09 1.4E+09			Zinc Zirconium	12122-67-7 7440-67-7				5.8E+04 9.3E+01	1.4E+05		4.1E+04 9.3E+01		