

Regional Removal Management Level (RML) Summary Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
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 _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
8.70E-03	I			4.00E-03	I					1	0.1	Acephate	30560-19-1	2.50E+02	n	3.30E+03	n	8.00E+01	n	
		2.20E-06	I			9.00E-03	I	V		1	1.07E+05	Acetaldehyde	75-07-0	8.20E+01	n	3.40E+02	n	1.90E+01	n	
				2.00E-02	I					1	0.1	Acetochlor	34256-82-1	1.30E+03	n	1.60E+04	n	3.50E+02	n	
				9.00E-01	I	3.10E+01	A	V		1	1.14E+05	Acetone	67-64-1	6.10E+04	n	6.70E+05	nms	1.40E+04	n	
						2.00E-03	X			1	0.1	Acetone Cyanohydrin	75-86-5	2.80E+06	nm	1.20E+07	nm			
						6.00E-02	I	V		1	1.28E+05	Acetonitrile	75-05-8	8.10E+02	n	3.40E+03	n	1.30E+02	n	
3.80E+00	C	1.30E-03	C	1.00E-01	I			V		1	2.52E+03	Acetophenone	98-86-2	7.80E+03	ns	1.20E+05	nms	1.90E+03	n	
				5.00E-04	I	2.00E-05	I	V		1	0.1	Acetylaminofluorene, 2-Acrolein	53-96-3	1.40E+01	c	6.00E+01	c	1.60E+00	c	
										1	2.27E+04		107-02-8	1.40E-01	n	6.00E-01	n	4.20E-02	n	
5.00E-01	I	1.00E-04	I	2.00E-03	I	6.00E-03	I	M		1	0.1	Acrylamide	79-06-1	2.40E+01	c**	4.60E+02	c**	5.00E+00	c**	
				5.00E-01	I	1.00E-03	I	V		1	1.09E+05	Acrylic Acid	79-10-7	9.90E+01	n	4.20E+02	n	2.10E+00	n	
5.40E-01	I	6.80E-05	I	4.00E-02	A	2.00E-03	I	V		1	1.13E+04	Acrylonitrile	107-13-1	1.60E+01	n	6.70E+01	n	4.10E+00	n	
						6.00E-03	P			1	0.1	Adiponitrile	111-69-3	8.50E+06	nm	3.60E+07	nm			
5.60E-02	C			1.00E-02	I					1	0.1	Alachlor	15972-60-8	6.30E+02	n	4.10E+03	c**	1.10E+02	c**	2.0E+00
				1.00E-03	I					1	0.1	Aldicarb	116-06-3	6.30E+01	n	8.20E+02	n	2.00E+01	n	3.0E+00
				1.00E-03	I					1	0.1	Aldicarb Sulfone	1646-88-4	6.30E+01	n	8.20E+02	n	2.00E+01	n	2.0E+00
1.70E+01	I	4.90E-03	I	3.00E-05	I			V		1	0.1	Aldicarb sulfoxide	1646-87-3							4.0E+00
										1	0.1	Aldrin	309-00-2	2.30E+00	n	1.80E+01	c**	9.20E-02	c**	
2.10E-02	C	6.00E-06	C	5.00E-03	I	1.00E-04	X	V		1	1.11E+05	Allyl Alcohol	107-18-6	3.50E+00	n	1.50E+01	n	2.10E-01	n	
				1.00E+00	P	5.00E-03	I	V		1	1.42E+03	Allyl Chloride	107-05-1	1.70E+00	n	6.90E+00	n	2.10E+00	n	
										1		Aluminum	7429-90-5	7.70E+04	n	1.10E+06	nm	2.00E+04	n	
2.10E+01	C	6.00E-03	C	4.00E-04	I					1		Aluminum Phosphide	20859-73-8	3.10E+01	n	4.70E+02	n	8.00E+00	n	
				9.00E-03	I					1	0.1	Ametryn	834-12-8	5.70E+02	n	7.40E+03	n	1.50E+02	n	
										1	0.1	Aminobiphenyl, 4-	92-67-1	2.60E+00	c	1.10E+01	c	3.00E-01	c	
				8.00E-02	P					1	0.1	Aminophenol, m-	591-27-5	5.10E+03	n	6.60E+04	n	1.60E+03	n	
				2.00E-02	P					1	0.1	Aminophenol, p-	123-30-8	1.30E+03	n	1.60E+04	n	4.00E+02	n	
				2.50E-03	I					1	0.1	Amitraz	33089-61-1	1.60E+02	n	2.10E+03	n	8.20E+00	n	
				1.00E-01	I	V				1		Ammonia	7664-41-7							
				2.00E-01	I					1		Ammonium Sulfamate	7773-06-0	1.60E+04	n	2.30E+05	nm	4.00E+03	n	
						3.00E-03	X	V		1	1.37E+04	Amyl Alcohol, tert-	75-85-4	8.20E+01	n	3.40E+02	n	6.30E+00	n	
5.70E-03	I	1.60E-06	C	7.00E-03	P	1.00E-03	I			1	0.1	Aniline	62-53-3	4.40E+02	n	5.70E+03	n	1.40E+02	n	
4.00E-02	P			2.00E-03	X					1	0.1	Anthraquinone, 9,10-	84-65-1	1.30E+02	n	1.60E+03	n	3.00E+01	n	
				4.00E-04	I				0.15			Antimony (metallic)	7440-36-0	3.10E+01	n	4.70E+02	n	7.80E+00	n	6.0E+00
				5.00E-04	H				0.15			Antimony Pentoxide	1314-60-9	3.90E+01	n	5.80E+02	n	9.70E+00	n	
				4.00E-04	H				0.15			Antimony Tetroxide	1332-81-6	3.10E+01	n	4.70E+02	n	7.80E+00	n	
						2.00E-04	I		0.15			Antimony Trioxide	1309-64-4	2.80E+05	nm	1.20E+06	nm			
1.50E+00	I	4.30E-03	I	3.00E-04	I	1.50E-05	C			1	0.03	Arsenic, Inorganic	7440-38-2	3.50E+01	nR	3.00E+02	c**R	5.20E+00	c**	1.0E+01
				3.50E-06	C	5.00E-05	I			1		Arsine	7784-42-1	2.70E-01	n	4.10E+00	n	7.00E-02	n	
				5.00E-02	I					1	0.1	Asulam	3337-71-1	3.20E+03	n	4.10E+04	n	1.00E+03	n	
2.30E-01	C			3.50E-02	I					1	0.1	Atrazine	1912-24-9	2.40E+02	c**	1.00E+03	c*	3.00E+01	c*	3.0E+00
8.80E-01	C	2.50E-04	C							1	0.1	Auramine	492-80-8	6.20E+01	c	2.60E+02	c	6.70E+00	c	
				4.00E-04	I					1	0.1	Avermectin B1	65195-55-3	2.50E+01	n	3.30E+02	n	8.00E+00	n	
1.10E-01	I	3.10E-05	I	3.00E-03	A	1.00E-02	A			1	0.1	Azinphos-methyl	86-50-0	1.90E+02	n	2.50E+03	n	5.60E+01	n	
								V		1		Azobenzene	103-33-3	5.60E+02	c	2.60E+03	c	1.20E+01	c	
				1.00E+00	P	7.00E-06	P			1	0.1	Azodicarbonamide	123-77-3	8.60E+03	n	4.00E+04	n	2.00E+04	n	
5.00E-01	C	1.50E-01	C	2.00E-01	I	5.00E-04	H		0.07			Barium	7440-39-3	1.50E+04	n	2.20E+05	nm	3.80E+03	n	2.0E+03
				2.00E-02	C	2.00E-04	C	M	0.025			Barium Chromate	10294-40-3	3.00E+01	c*	6.20E+02	c*	4.10E+00	c*	
				3.00E-01	I			V		1		Benfluralin	1861-40-1	2.30E+04	n	3.50E+05	nm	1.70E+03	n	
				5.00E-02	I					1	0.1	Benomyl	17804-35-2	3.20E+03	n	4.10E+04	n	9.70E+02	n	
				2.00E-01	I					1	0.1	Bensulfuron-methyl	83055-99-6	1.30E+04	n	1.60E+05	nm	3.90E+03	n	
				3.00E-02	I					1	0.1	Bentazon	25057-89-0	1.90E+03	n	2.50E+04	n	5.70E+02	n	
4.00E-03	P			1.00E-01	I			V		1	1.16E+03	Benzaldehyde	100-52-7	7.80E+03	ns	8.20E+04	c**s	1.90E+03	c**	
5.50E-02	I	7.80E-06	I	4.00E-03	I	3.00E-02	I	V		1	1.82E+03	Benzene	71-43-2	8.20E+01	n	4.20E+02	n	3.30E+01	n	5.0E+00
1.00E-01	X			3.00E-04	X					1	0.1	Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1	1.90E+01	n	2.50E+02	n	6.00E+00	n	
				1.00E-03	P			V		1	1.26E+03	Benzenethiol	108-98-5	7.80E+01	n	1.20E+03	n	1.70E+01	n	

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Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
2.30E+02	I	6.70E-02	I	3.00E-03 4.00E+00	I			M	1	0.1		Benzidine Benzoic Acid	92-87-5 65-85-0	5.30E-02 2.50E+05	c nm	1.00E+00 3.30E+06	c nm	1.10E-02 7.50E+04	c n	
1.30E+01	I							V	1	0.1	3.24E+02	Benzotrchloride	98-07-7	5.30E+00	c	2.50E+01	c	3.00E-01	c	
1.70E-01	I	4.90E-05	C	1.00E-01 2.00E-03	P			P	1		1.46E+03	Benzyl Alcohol Benzyl Chloride	100-51-6 100-44-7	6.30E+03 2.30E+01	n n	8.20E+04 1.10E+02	n n	2.00E+03 2.00E+00	n n	
		2.40E-03	I	2.00E-03 9.00E-03 1.50E-02	I			2.00E-05	I		0.007	Beryllium and compounds Bifenox Biphenrin	7440-41-7 42576-02-3 82657-04-3	1.60E+02 5.70E+02 9.50E+02	n n n	2.30E+03 7.40E+03 1.20E+04	n n n	2.50E+01 1.00E+02 3.00E+02	n n n	4.0E+00
8.00E-03	I			5.00E-01 4.00E-02 3.00E-03	I			4.00E-04	X	V	1	Biphenyl, 1,1'- Bis(2-chloro-1-methylethyl) ether Bis(2-chloroethoxy)methane	92-52-4 108-60-1 111-91-1	4.70E+01 3.10E+03 1.90E+02	n ns n	2.00E+02 4.70E+04 2.50E+03	n ns n	8.30E-01 7.10E+02 5.90E+01	n n n	
1.10E+00	I	3.30E-04	I					V	1		5.05E+03	Bis(2-chloroethyl)ether	111-44-4	2.30E+01	c	1.00E+02	c	1.40E+00	c	
2.20E+02	I	6.20E-02	I					V	1		4.22E+03	Bis(chloromethyl)ether	542-88-1	8.30E-03	c	3.60E-02	c	7.20E-03	c	
				5.00E-02	I				1	0.1		Bisphenol A	80-05-7	3.20E+03	n	4.10E+04	n	7.70E+02	n	
				2.00E-01 2.00E+00 4.00E-02	I P			2.00E-02 2.00E-02	H P		1	Boron And Borates Only Boron Trichloride Boron Trifluoride	7440-42-8 10294-34-5 7637-07-2	1.60E+02 1.60E+05 3.10E+03	n nm n	2.30E+05 2.30E+06 4.70E+04	nm nm n	4.00E+03 4.20E+01 2.60E+01	n n n	
7.00E-01	I			4.00E-03	I				1			Bromate	15541-45-4	9.90E+01	c**	4.70E+02	c*	1.10E+01	c**	1.0E+01
2.00E+00	X	6.00E-04	X					V	1		2.38E+03	Bromo-2-chloroethane, 1- Bromobenzene	107-04-0 108-86-1	2.60E+00 2.90E+02	c n	1.10E+01 1.80E+03	c ns	7.40E-01 6.20E+01	c n	
				8.00E-03	I			4.00E-02	X	V	1	Bromochloromethane Bromodichloromethane Bromoform	74-97-5 75-27-4 75-25-2	1.50E+02 2.90E+01 1.60E+03	n c*	6.30E+02 1.30E+02	n c	8.30E+01 1.30E+01	n c*	8.0E+01(F) 8.0E+01(F)
6.20E-02	I	3.70E-05	C	2.00E-02	I			V	1		9.32E+02	Bromomethane	74-83-9	6.80E+00	n	3.00E+01	n	7.50E+00	n	
7.90E-03	I	1.10E-06	I	2.00E-02	I			V	1		9.15E+02	Bromophos Bromoxynil	2104-96-3 1689-84-5	3.90E+02 1.30E+03	n n	5.80E+03 1.60E+04	n n	3.50E+01 3.30E+02	n n	
				1.40E-03 5.00E-03 2.00E-02	I H I			5.00E-03	I	V	1	Bromoxynil Octanoate	1689-99-2	6.80E+00	n	3.00E+01	n	7.50E+00	n	
3.40E+00	C	3.00E-05	I	2.00E-02	I			V	1		6.67E+02	Butadiene, 1,3- Butanol, N-	106-99-0 71-36-3	1.60E+03 1.80E+00	n n	2.30E+04 7.60E+00	n nms	1.40E+02 1.80E+00	n c**	
				1.00E-01	I			V	1		7.64E+03	Butyl alcohol, sec- Butylate	78-92-2 2008-41-5	1.30E+05 3.90E+03	nms n	1.50E+06 5.80E+04	nms n	2.40E+04 4.60E+02	n n	
2.00E-04	C	5.70E-08	C	5.00E-02	I			V	1	0.1	2.13E+04	Butylated hydroxyanisole	25013-16-5	2.70E+05	cm	1.10E+06	cm	1.50E+04	c	
3.60E-03	P			3.00E-01 5.00E-02 1.00E-01	P P X			V	1	0.1	1.08E+02	Butylated hydroxytoluene Butylbenzene, n- Butylbenzene, sec-	128-37-0 104-51-8 135-98-8	1.50E+04 3.90E+03 7.80E+03	c** ns ns	6.40E+04 5.80E+04 1.20E+05	c** ns nms	3.40E+02 1.00E+03 2.00E+03	c** n n	
				1.00E-01 2.00E-02 1.00E-03	X A I			V	1	0.1	1.83E+02	Butylbenzene, tert- Cacodylic Acid Cadmium (Diet)	98-06-6 75-60-5 7440-43-9	7.80E+03 1.30E+03 7.10E+01	ns n n	1.20E+05 1.60E+04 9.80E+02	nms n n	6.90E+02 4.00E+02	n n	
5.00E-01	C	1.50E-01	C	5.00E-04 2.00E-02	I C			1.00E-05 2.00E-04	A C		0.05 0.025	Cadmium (Water) Calcium Chromate Caprolactam	7440-43-9 13765-19-0 105-60-2	3.00E+01 3.10E+04	c* n	6.20E+02 4.00E+05	c* n	4.10E+00 9.90E+03	c* n	
1.50E-01	C	4.30E-05	C	2.00E-03	I				1	0.1		Captafol	2425-06-1	1.30E+02	n	1.50E+03	c**	3.20E+01	n	
2.30E-03	C	6.60E-07	C	1.30E-01	I				1	0.1		Captan	133-06-2	8.20E+03	n	1.00E+05	c**	2.40E+03	n	
				1.00E-01	I				1	0.1		Carbaryl	63-25-2	6.30E+03	n	8.20E+04	n	1.80E+03	n	
				5.00E-03	I				1	0.1		Carbofuran	1563-66-2	3.20E+02	n	4.10E+03	n	9.40E+01	n	4.0E+01
7.00E-02	I	6.00E-06	I	1.00E-01 4.00E-03	I I			7.00E-01 1.00E-01	I I	V	1	Carbon Disulfide Carbon Tetrachloride	75-15-0 56-23-5	7.70E+02 6.50E+01	ns c**	3.50E+03 2.90E+02	ns c**	8.10E+02 4.60E+01	n c**	5.0E+00
				1.00E-01	P			V	1		5.89E+03	Carbonyl Sulfide	463-58-1	6.70E+01	n	2.80E+02	n	2.10E+02	n	
				1.00E-02 1.00E-01	I I				1 1	0.1 0.1		Carbosulfan Carboxin	55285-14-8 5234-68-4	6.30E+02 6.30E+03	n n	8.20E+03 8.20E+04	n n	5.10E+01 1.90E+03	n n	
				9.00E-04	I				1			Ceric oxide	1306-38-3	1.30E+06	nm	5.40E+06	nm			
				1.00E-01 1.50E-02	I I			V	1 1	0.1 0.1		Chloral hydrate Chloramben	302-17-0 133-90-4	7.80E+03 9.50E+02	n n	1.20E+05 1.20E+04	nm n	2.00E+03 2.90E+02	n n	
4.00E-01	H								1	0.1		Chloranil	118-75-2	1.30E+02	c	5.70E+02	c	1.80E+01	c	
3.50E-01	I	1.00E-04	I	5.00E-04	I			7.00E-04	I	V	1	Chlordane	12789-03-6	3.50E+01	n	4.50E+02	n	7.40E-01	n	2.0E+00

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1.00E+01	I	4.60E-03	C	3.00E-04	I				1	0.1		Chlordecone (Kepone)	143-50-0	5.40E+00	c**	2.30E+01	c*	3.50E-01	c**	
				7.00E-04	A				1	0.1		Chlorfenvinphos	470-90-6	4.40E+01	n	5.70E+02	n	1.10E+01	n	
				2.00E-02	I				1	0.1		Chlorimuron, Ethyl- Chlorine	90982-32-4 7782-50-5	1.30E+03 1.80E-01	n n	1.60E+04 7.80E-01	n n	3.90E+02 3.00E-01	n n	
				3.00E-02	I	2.00E-04	I	V	1			Chlorine Dioxide	10049-04-4	2.30E+03	n	3.40E+04	n	4.20E-01	n	
				3.00E-02	I				1			Chlorite (Sodium Salt)	7758-19-2	2.30E+03	n	3.50E+04	n	6.00E+02	n	1.0E+03
				5.00E+01	I	V			1		1.15E+03	Chloro-1,1-difluoroethane, 1- Chloro-1,3-butadiene, 2- Chloro-2-methylaniline HCl, 4- Chloro-2-methylaniline, 4-	75-68-3 126-99-8 3165-93-3 95-69-2	5.40E+04 1.00E+00 1.20E+02 1.90E+02	ns c* c n	2.30E+05 4.40E+00 5.00E+02 2.30E+03	nms c* c c**	1.00E+05 1.90E+00 1.70E+01 5.40E+01	n c* c n	
2.70E-01	X							V	1		1.18E+04	Chloroacetaldehyde, 2- Chloroacetic Acid Chloroacetophenone, 2-	107-20-0 79-11-8 532-27-4	2.60E+02	c	1.20E+03	c	2.90E+01	c	6.0E+01
				3.00E-05	I				1	0.1		Chloroaniline, p- Chlorobenzene Chlorobenzilate	106-47-8 108-90-7 510-15-6	2.50E+02 2.80E+02 4.90E+02	n n c**	1.10E+03 1.30E+03 2.10E+03	c** ns c**	3.70E+01 7.80E+01 3.10E+01	c** n c**	1.0E+02
				3.00E-02	X				1	0.1		Chlorobenzoic Acid, p- Chlorobenzotrifluoride, 4- Chlorobutane, 1-	74-11-3 98-56-6 109-69-3	1.90E+03 2.10E+02 3.10E+03	n n ns	2.50E+04 2.50E+03 4.70E+04	n ns ns	5.10E+02 3.50E+01 6.40E+02	n n n	
				5.00E+01	I	V			1		1.68E+03	Chlorodifluoromethane	75-45-6	4.90E+04	ns	2.10E+05	nms	1.00E+05	n	
3.10E-02	C	2.30E-05	I	1.00E-02	I	9.80E-02	A	V	1		2.54E+03	Chloroethanol, 2- Chloroform	107-07-3 67-66-3	1.60E+03 3.20E+01	n c**	2.30E+04 1.40E+02	n c**	4.00E+02 2.20E+01	n c**	8.0E+01(F)
				9.00E-02	I	V			1		1.32E+03	Chloromethane	74-87-3	1.10E+02	n	4.60E+02	n	1.90E+02	n	
2.40E+00	C	6.90E-04	C					V	1		9.32E+03	Chloromethyl Methyl Ether	107-30-2	2.00E+00	c	8.90E+00	c	6.50E-01	c	
3.00E-01	P			3.00E-03	P	1.00E-05	X		1	0.1		Chloronitrobenzene, o-	88-73-3	1.80E+02	c**	7.70E+02	c**	2.40E+01	c**	
6.00E-02	P			7.00E-04	P	2.00E-03	P		1	0.1		Chloronitrobenzene, p- Chlorophenol, 2- Chloropicrin	100-00-5 95-57-8 76-06-2	4.40E+01 3.90E+02 2.00E+00	n n n	5.70E+02 5.80E+03 8.20E+00	n n n	1.30E+01 9.10E+01 8.30E-01	n n n	
3.10E-03	C	8.90E-07	C	1.50E-02	I				1	0.1		Chlorothalonil	1897-45-6	9.50E+02	n	1.20E+04	n	2.60E+02	n	
				2.00E-02	I			V	1		9.07E+02	Chlorotoluene, o- Chlorotoluene, p-	95-49-8 106-43-4	1.60E+03 1.60E+03	ns ns	2.30E+04 2.30E+04	ns ns	2.40E+02 2.50E+02	n n	
2.40E+02	C	6.90E-02	C						1	0.1		Chlorozotocin	54749-90-5	2.30E-01	c	9.60E-01	c	3.20E-02	c	
				2.00E-01	I				1	0.1		Chlorpropham	101-21-3	1.30E+04	n	1.60E+05	nm	2.80E+03	n	
				1.00E-03	A				1	0.1		Chlorpyrifos	2921-88-2	6.30E+01	n	8.20E+02	n	8.40E+00	n	
				1.00E-02	H				1	0.1		Chlorpyrifos Methyl	5598-13-0	6.30E+02	n	8.20E+03	n	1.20E+02	n	
				5.00E-02	I				1	0.1		Chlorsulfuron	64902-72-3	3.20E+03	n	4.10E+04	n	9.90E+02	n	
				1.00E-02	I				1	0.1		Chlorthal-dimethyl	1861-32-1	6.30E+02	n	8.20E+03	n	1.20E+02	n	
				8.00E-04	H				1	0.1		Chlorthiophos	60238-56-4	5.10E+01	n	6.60E+02	n	2.80E+00	n	
5.00E-01	J	8.40E-02	S	1.50E+00	I	1.00E-04	I	M	0.013	0.025		Chromium(III), Insoluble Salts Chromium(VI)	16065-83-1 18540-29-9	1.20E+05 3.00E+01	nm c**	1.80E+06 6.30E+02	nm c**	2.20E+04 3.50E+00	n c*	
				1.30E-02	I				1	0.1		Chromium, Total	7440-47-3	8.20E+02	n	1.10E+04	n	2.30E+02	n	1.0E+02
				9.00E-03	P	3.00E-04	P	6.00E-06	P			Clofentezine	74115-24-5	2.30E+01	n	3.50E+02	n	6.00E+00	n	
				6.20E-04	I			V	M			Cobalt	7440-48-4							
				4.00E-02	H				1			Coke Oven Emissions	8007-45-2							
				5.00E-02	I	6.00E-01	C		1	0.1		Copper	7440-50-8	3.10E+03	n	4.70E+04	n	8.00E+02	n	1.3E+03
				1.00E-01	A	6.00E-01	C		1	0.1		Cresol, m-	108-39-4	3.20E+03	n	4.10E+04	n	9.30E+02	n	
				5.00E-02	I	6.00E-01	C		1	0.1		Cresol, o-	95-48-7	3.20E+03	n	4.10E+04	n	9.30E+02	n	
				1.00E-01	A	6.00E-01	C		1	0.1		Cresol, p-	106-44-5	6.30E+03	n	8.20E+04	n	1.90E+03	n	
				1.00E-01	A				1	0.1		Cresol, p-chloro-m-	59-50-7	6.30E+03	n	8.20E+04	n	1.40E+03	n	
1.90E+00	H			1.00E-01	A	6.00E-01	C		1	0.1		Cresols	1319-77-3	6.30E+03	n	8.20E+04	n	1.50E+03	n	
				1.00E-03	P			V	1		1.66E+04	Crotonaldehyde, trans-	123-73-9	3.70E+01	c**	1.70E+02	c**	4.00E+00	c**	
				1.00E-01	I	4.00E-01	I	V	1		2.68E+02	Cumene	98-82-8	1.90E+03	ns	9.90E+03	ns	4.50E+02	n	
2.20E-01	C	6.30E-05	C						1	0.1		Cupferron	135-20-6	2.50E+02	c	1.00E+03	c	3.50E+01	c	
8.40E-01	H			2.00E-03	H				1	0.1		Cyanazine	21725-46-2	6.50E+01	c**	2.70E+02	c**	8.80E+00	c**	

Regional Removal Management Level (RML) Summary Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
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 _i (mg/m ³)	k e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
											Cyanides									
		1.00E-03	I							1		~Calcium Cyanide	592-01-8	7.80E+01	n	1.20E+03	n	2.00E+01	n	
		5.00E-03	I							1		~Copper Cyanide	544-92-3	3.90E+02	n	5.80E+03	n	1.00E+02	n	
		6.00E-04	I	8.00E-04	S	V				1	9.54E+05	~Cyanide (CN-)	57-12-5	2.30E+01	n	1.50E+02	n	1.50E+00	n	2.0E+02
		1.00E-03	I					V		1		~Cyanogen	460-19-5	7.80E+01	n	1.20E+03	n	2.00E+01	n	
		9.00E-02	I					V		1		~Cyanogen Bromide	506-68-3	7.00E+03	n	1.10E+05	nm	1.80E+03	n	
		5.00E-02	I					V		1		~Cyanogen Chloride	506-77-4	3.90E+03	n	5.80E+04	n	1.00E+03	n	
		6.00E-04	I	8.00E-04	I	V				1	1.00E+07	~Hydrogen Cyanide	74-90-8	2.30E+01	n	1.50E+02	n	1.50E+00	n	
		2.00E-03	I							1		~Potassium Cyanide	151-50-8	1.60E+02	n	2.30E+03	n	4.00E+01	n	
		5.00E-03	I							0.04		~Potassium Silver Cyanide	506-61-6	3.90E+02	n	5.80E+03	n	8.20E+01	n	
		1.00E-01	I							0.04		~Silver Cyanide	506-64-9	7.80E+03	n	1.20E+05	nm	1.80E+03	n	
		1.00E-03	I							1		~Sodium Cyanide	143-33-9	7.80E+01	n	1.20E+03	n	2.00E+01	n	2.0E+02
		2.00E-04	P							1		~Thiocyanates	NA	1.60E+01	n	2.30E+02	n	4.00E+00	n	
		2.00E-04	X					V		1		~Thiocyanic Acid	463-56-9	1.60E+01	n	2.30E+02	n	4.00E+00	n	
		5.00E-02	I							1		~Zinc Cyanide	557-21-1	3.90E+03	n	5.80E+04	n	1.00E+03	n	
				6.00E+00	I	V				1	1.17E+02	Cyclohexane	110-82-7	6.50E+03	ns	2.70E+04	ns	1.30E+04	n	
2.30E-02	H			5.00E+00	I	7.00E-01	P	V		1	5.11E+03	Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	2.40E+03	c	1.00E+04	c	2.40E+02	c	
				5.00E-03	P	1.00E+00	X	V		1	2.83E+02	Cyclohexanone	108-94-1	2.80E+04	ns	1.30E+05	nms	1.40E+03	n	
										1		Cyclohexene	110-83-8	3.10E+02	ns	3.10E+03	ns	7.00E+01	n	
		2.00E-01	I					V		1	2.93E+05	Cyclohexylamine	108-91-8	1.60E+04	n	2.30E+05	nm	3.80E+03	n	
		2.50E-02	I							1	0.1	Cyfluthrin	68359-37-5	1.60E+03	n	2.10E+04	n	1.20E+02	n	
		5.00E-03	I							1	0.1	Cyhalothrin	68085-85-8	3.20E+02	n	4.10E+03	n	1.00E+02	n	
		1.00E-02	I							1	0.1	Cypermethrin	52315-07-8	6.30E+02	n	8.20E+03	n	2.00E+02	n	
		7.50E-03	I							1	0.1	Cyromazine	66215-27-8	4.70E+02	n	6.20E+03	n	1.50E+02	n	
2.40E-01	I	6.90E-05	C							1	0.1	DDD	72-54-8	2.30E+02	c	9.60E+02	c	3.20E+00	c	
3.40E-01	I	9.70E-05	C					V		1		DDE, p,p'-	72-55-9	2.00E+02	c	9.30E+02	c	4.60E+00	c	
3.40E-01	I	9.70E-05	I	5.00E-04	I					1	0.03	DDT	50-29-3	3.70E+01	n	5.20E+02	n	1.00E+01	n	
				3.00E-02	I					1	0.1	Dalapon	75-99-0	1.90E+03	n	2.50E+04	n	6.00E+02	n	2.0E+02
1.80E-02	C	5.10E-06	C	1.50E-01	I					1	0.1	Daminozide	1596-84-5	3.00E+03	c**	1.30E+04	c**	4.30E+02	c**	
7.00E-04	I			7.00E-03	I					1	0.1	Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5	4.40E+02	n	5.70E+03	n	1.40E+02	n	
				4.00E-05	I					1	0.1	Demeton	8065-48-3	2.50E+00	n	3.30E+01	n	4.20E-01	n	
1.20E-03	I			6.00E-01	I					1	0.1	Di(2-ethylhexyl)adipate	103-23-1	3.80E+04	n	1.90E+05	c**m	6.50E+03	c**	4.0E+02
6.10E-02	H			7.00E-04	A					1	0.1	Diallate	2303-16-4	8.90E+02	c	3.80E+03	c	5.40E+01	c	
										1	0.1	Diazinon	333-41-5	4.40E+01	n	5.70E+02	n	1.00E+01	n	
		1.00E-02	X					V		1		Dibenzothiophene	132-65-0	7.80E+02	n	1.20E+04	n	6.50E+01	n	
8.00E-01	P	6.00E-03	P	2.00E-04	P	2.00E-04	I	V	M	1	9.79E+02	Dibromo-3-chloropropane, 1,2-	96-12-8	5.30E-01	c**	6.40E+00	c**	3.30E-02	c*	2.0E-01
				4.00E-04	X			V		1	1.59E+02	Dibromobenzene, 1,3-	108-36-1	3.10E+01	n	4.70E+02	ns	5.30E+00	n	
		1.00E-02	I					V		1		Dibromobenzene, 1,4-	106-37-6	7.80E+02	n	1.20E+04	n	1.30E+02	n	
8.40E-02	I			2.00E-02	I			V		1	8.02E+02	Dibromochloromethane	124-48-1	8.30E+02	c**s	3.90E+03	c**s	8.70E+01	c**	8.0E+01(F)
2.00E+00	I	6.00E-04	I	9.00E-03	I	9.00E-03	I	V		1	1.34E+03	Dibromoethane, 1,2-	106-93-4	3.60E+00	c*	1.60E+01	c*	7.50E-01	c*	5.0E-02
				4.00E-03	X	V				1	2.82E+03	Dibromomethane (Methylene Bromide)	74-95-3	2.40E+01	n	9.90E+01	n	8.30E+00	n	
		3.00E-04	P							1	0.1	Dibutyltin Compounds	NA	1.90E+01	n	2.50E+02	n	6.00E+00	n	
		3.00E-02	I							1	0.1	Dicamba	1918-00-9	1.90E+03	n	2.50E+04	n	5.70E+02	n	
4.20E-03	P							V		1	5.54E+02	Dichloro-2-butene, 1,4-	764-41-0	2.10E-01	c	9.40E-01	c	1.30E-01	c	
4.20E-03	P							V		1	5.19E+02	Dichloro-2-butene, cis-1,4-	1476-11-5	7.40E-01	c	3.20E+00	c	1.30E-01	c	
4.20E-03	P							V		1	7.60E+02	Dichloro-2-butene, trans-1,4-	110-57-6	7.40E-01	c	3.20E+00	c	1.30E-01	c	
5.00E-02	I			4.00E-03	I					1	0.1	Dichloroacetic Acid	79-43-6	2.50E+02	n	3.30E+03	n	7.90E+01	n	6.0E+01
				9.00E-02	I	2.00E-01	H	V		1	3.76E+02	Dichlorobenzene, 1,2-	95-50-1	1.80E+03	ns	9.30E+03	ns	3.00E+02	n	6.0E+02
5.40E-03	C	1.10E-05	C	7.00E-02	A	8.00E-01	I	V		1		Dichlorobenzene, 1,4-	106-46-7	2.60E+02	c*	1.10E+03	c*	4.80E+01	c*	7.5E+01
4.50E-01	I	3.40E-04	C							1	0.1	Dichlorobenzidine, 3,3'-	91-94-1	1.20E+02	c	5.10E+02	c	1.30E+01	c	
				9.00E-03	X					1	0.1	Dichlorobenzophenone, 4,4'-	90-98-2	5.70E+02	n	7.40E+03	n	7.80E+01	n	
				2.00E-01	I	1.00E-01	X	V		1	8.45E+02	Dichlorodifluoromethane	75-71-8	8.70E+01	n	3.70E+02	n	2.00E+02	n	
5.70E-03	C	1.60E-06	C	2.00E-01	P			V		1	1.69E+03	Dichloroethane, 1,1-	75-34-3	3.60E+02	c*	1.60E+03	c	2.80E+02	c*	
9.10E-02	I	2.60E-05	I	6.00E-03	X	7.00E-03	P	V		1	2.98E+03	Dichloroethane, 1,2-	107-06-2	3.10E+01	n	1.40E+02	n	1.30E+01	n	5.0E+00

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Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
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 _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
				5.00E-02	I	2.00E-01	I	V			1.19E+03	Dichloroethylene, 1,1-	75-35-4	2.30E+02	n	1.00E+03	n	2.80E+02	n	7.0E+00
				2.00E-03	I			V			2.37E+03	Dichloroethylene, 1,2-cis-	156-59-2	1.60E+02	n	2.30E+03	n	3.60E+01	n	7.0E+01
				2.00E-02	I			V			1.85E+03	Dichloroethylene, 1,2-trans-	156-60-5	1.60E+03	n	2.30E+04	ns	3.60E+02	n	1.0E+02
				3.00E-03	I					0.1		Dichlorophenol, 2,4-	120-83-2	1.90E+02	n	2.50E+03	n	4.60E+01	n	
				1.00E-02	I						0.05	Dichlorophenoxy Acetic Acid, 2,4-	94-75-7	7.00E+02	n	9.60E+03	n	1.70E+02	n	7.0E+01
				8.00E-03	I						0.1	Dichlorophenoxy)butyric Acid, 4-(2,4-	94-82-6	5.10E+02	n	6.60E+03	n	1.20E+02	n	
3.60E-02	C	1.00E-05	C	9.00E-02	A	4.00E-03	I	V			1.36E+03	Dichloropropane, 1,2-	78-87-5	1.60E+01	n	6.60E+01	n	8.30E+00	n	5.0E+00
				2.00E-02	P			V			1.49E+03	Dichloropropane, 1,3-	142-28-9	1.60E+03	ns	2.30E+04	ns	3.70E+02	n	
				3.00E-03	I						0.1	Dichloropropanol, 2,3-	616-23-9	1.90E+02	n	2.50E+03	n	5.90E+01	n	
1.00E-01	I	4.00E-06	I	3.00E-02	I	2.00E-02	I	V			1.57E+03	Dichloropropene, 1,3-	542-75-6	7.20E+01	n	3.10E+02	n	3.90E+01	n	
2.90E-01	I	8.30E-05	C	5.00E-04	I	5.00E-04	I				0.1	Dichlorvos	62-73-7	3.20E+01	n	4.10E+02	n	9.90E+00	n	
				1.00E-04	I						0.1	Dicrotophos	141-66-2	6.30E+00	n	8.20E+01	n	2.00E+00	n	
				8.00E-02	P	3.00E-04	X	V			2.56E+02	Dicyclopentadiene	77-73-6	1.30E+00	n	5.40E+00	n	6.30E-01	n	
1.60E+01	I	4.60E-03	I	5.00E-05	I						0.1	Dieldrin	60-57-1	3.20E+00	n	1.40E+01	c**	1.80E-01	c**	
		3.00E-04	C			5.00E-03	I				0.1	Diesel Engine Exhaust	NA							
				2.00E-03	P	2.00E-04	P				0.1	Diethanolamine	111-42-2	1.30E+02	n	1.60E+03	n	4.00E+01	n	
				3.00E-02	P	1.00E-04	P				0.1	Diethylene Glycol Monobutyl Ether	112-34-5	1.90E+03	n	2.40E+04	n	6.00E+02	n	
				6.00E-02	P	3.00E-04	P				0.1	Diethylene Glycol Monoethyl Ether	111-90-0	3.80E+03	n	4.80E+04	n	1.20E+03	n	
				1.00E-03	P			V			1.12E+05	Diethylformamide	617-84-5	7.80E+01	n	1.20E+03	n	2.00E+01	n	
3.50E+02	C	1.00E-01	C								0.1	Diethylstilbestrol	56-53-1	1.60E-01	c	6.60E-01	c	5.10E-03	c	
				8.00E-02	I						0.1	Difenzoquat	43222-48-6	5.10E+03	n	6.60E+04	n	1.60E+03	n	
				2.00E-02	I						0.1	Diflubenzuron	35367-38-5	1.30E+03	n	1.60E+04	n	2.90E+02	n	
4.40E-02	C	1.30E-05	C			4.00E+01	I	V			1.43E+03	Diffuoroethane, 1,1-	75-37-6	4.80E+04	ns	2.00E+05	nms	8.30E+04	n	
								V			1	Dihydrosafrole	94-58-6	9.90E+02	c	4.50E+03	c	3.00E+01	c	
						7.00E-01	P	V			2.26E+03	Diisopropyl Ether	108-20-3	2.20E+03	n	9.40E+03	ns	1.50E+03	n	
				8.00E-02	I			V			5.30E+02	Diisopropyl Methylphosphonate	1445-75-6	6.30E+03	ns	9.30E+04	ns	1.60E+03	n	
				2.00E-02	I						0.1	Dimethipin	55290-64-7	1.30E+03	n	1.60E+04	n	4.00E+02	n	
				2.00E-04	I						0.1	Dimethoate	60-51-5	1.30E+01	n	1.60E+02	n	4.00E+00	n	
1.60E+00	P										0.1	Dimethoxybenzidine, 3,3'-	119-90-4	3.40E+01	c	1.40E+02	c	4.70E+00	c	
1.70E-03	P			6.00E-02	P						0.1	Dimethyl methylphosphonate	756-79-6	3.80E+03	n	4.90E+04	n	1.20E+03	n	
4.60E+00	C	1.30E-03	C								0.1	Dimethylamino azobenzene [p]	60-11-7	1.20E+01	c	5.00E+01	c	5.00E-01	c	
5.80E-01	H										0.1	Dimethylaniline HCl, 2,4-	21436-96-4	9.40E+01	c	4.00E+02	c	1.30E+01	c	
2.00E-01	P			2.00E-03	X						0.1	Dimethylaniline, 2,4-	95-68-1	1.30E+02	n	1.10E+03	c**	3.70E+01	c**	
				2.00E-03	I			V			8.30E+02	Dimethylaniline, N,N-	121-69-7	1.60E+02	n	2.30E+03	ns	3.50E+01	n	
1.10E+01	P										0.1	Dimethylbenzidine, 3,3'-	119-93-7	4.90E+00	c	2.10E+01	c	6.50E-01	c	
				1.00E-01	P	3.00E-02	I	V			1.06E+05	Dimethylformamide	68-12-2	2.60E+03	n	1.50E+04	n	6.10E+01	n	
				1.00E-04	X	2.00E-06	X	V			1.72E+05	Dimethylhydrazine, 1,1-	57-14-7	5.70E-02	n	2.40E-01	n	4.20E-03	n	
5.50E+02	C	1.60E-01	C					V			1.89E+05	Dimethylhydrazine, 1,2-	540-73-8	8.80E-02	c	4.10E-01	c	2.80E-03	c	
				2.00E-02	I						0.1	Dimethylphenol, 2,4-	105-67-9	1.30E+03	n	1.60E+04	n	3.60E+02	n	
				6.00E-04	I						0.1	Dimethylphenol, 2,6-	576-26-1	3.80E+01	n	4.90E+02	n	1.10E+01	n	
4.50E-02	C	1.30E-05	C								0.1	Dimethylphenol, 3,4-	95-65-8	6.30E+01	n	8.20E+02	n	1.80E+01	n	
				1.00E-03	I			V			4.73E+02	Dimethylvinylchloride	513-37-1	1.10E+02	c	4.80E+02	cs	3.30E+01	c	
				8.00E-05	X						0.1	Dinitro-o-cresol, 4,6-	534-52-1	5.10E+00	n	6.60E+01	n	1.50E+00	n	
				2.00E-03	I						0.1	Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5	1.30E+02	n	1.60E+03	n	2.30E+01	n	
				1.00E-04	P						0.1	Dinitrobenzene, 1,2-	528-29-0	6.30E+00	n	8.20E+01	n	1.90E+00	n	
				1.00E-04	I						0.1	Dinitrobenzene, 1,3-	99-65-0	6.30E+00	n	8.20E+01	n	2.00E+00	n	
6.80E-01	I			1.00E-04	P						0.1	Dinitrobenzene, 1,4-	100-25-4	6.30E+00	n	8.20E+01	n	2.00E+00	n	
				2.00E-03	I						0.1	Dinitrophenol, 2,4-	51-28-5	1.30E+02	n	1.60E+03	n	3.90E+01	n	
											0.1	Dinitrotoluene Mixture, 2,4/2,6-	NA	8.00E+01	c	3.40E+02	c	1.10E+01	c	
3.10E-01	C	8.90E-05	C	2.00E-03	I						0.102	Dinitrotoluene, 2,4-	121-14-2	1.30E+02	n	7.40E+02	c**	2.40E+01	c**	
1.50E+00	P			3.00E-04	X						0.099	Dinitrotoluene, 2,6-	606-20-2	1.90E+01	n	1.50E+02	c**	4.90E+00	c**	
				2.00E-03	S						0.006	Dinitrotoluene, 2-Amino-4,6-	35572-78-2	1.50E+02	n	2.30E+03	n	3.90E+01	n	
4.50E-01	X			2.00E-03	S						0.009	Dinitrotoluene, 4-Amino-2,6-	19406-51-0	1.50E+02	n	2.30E+03	n	3.90E+01	n	
				9.00E-04	X						0.1	Dinitrotoluene, Technical grade	25321-14-6	5.70E+01	n	5.10E+02	c**	1.00E+01	c**	

Regional Removal Management Level (RML) Summary Table (TR=1E-04, HQ=1) May 2016

Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
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 _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
				1.00E-03	I					1	0.1	Dinoseb	88-85-7	6.30E+01	n	8.20E+02	n	1.50E+01	n	7.0E+00
1.00E-01	I	5.00E-06	I	3.00E-02	I	3.00E-02	I	V		1	1.16E+05	Dioxane, 1,4-	123-91-1	5.30E+02	c**	2.40E+03	c**	4.60E+01	c**	
6.20E+03	I	1.30E+00	I							1	0.03	Dioxins ~Hexachlorodibenzo-p-dioxin, Mixture	NA	1.00E-02	c	4.70E-02	c	1.30E-03	c	
1.30E+05	C	3.80E+01	C	7.00E-10	I	4.00E-08	C	V		1	0.03	~TCDD, 2,3,7,8-	1746-01-6	5.10E-05	n	7.20E-04	n	1.20E-05	c**	3.0E-05
				3.00E-02	I					1	0.1	Diphenamid	957-51-7	1.90E+03	n	2.50E+04	n	5.30E+02	n	
				8.00E-04	X					1	0.1	Diphenyl Sulfone	127-63-9	5.10E+01	n	6.60E+02	n	1.50E+01	n	
8.00E-01	I	2.20E-04	I	2.50E-02	I					1	0.1	Diphenylamine	122-39-4	1.60E+03	n	2.10E+04	n	3.10E+02	n	
				2.20E-03	I					1	0.1	Diphenylhydrazine, 1,2-	122-66-7	6.80E+01	c	2.90E+02	c	7.80E+00	c	
										1	0.1	Diquat	85-00-7	1.40E+02	n	1.80E+03	n	4.40E+01	n	2.0E+01
7.10E+00	C	1.40E-01	C							1	0.1	Direct Black 38	1937-37-7	7.60E+00	c	3.20E+01	c	1.10E+00	c	
7.40E+00	C	1.40E-01	C							1	0.1	Direct Blue 6	2602-46-2	7.30E+00	c	3.10E+01	c	1.10E+00	c	
6.70E+00	C	1.40E-01	C							1	0.1	Direct Brown 95	16071-86-6	8.10E+00	c	3.40E+01	c	1.20E+00	c	
				4.00E-05	I					1	0.1	Disulfoton	298-04-4	2.50E+00	n	3.30E+01	n	5.00E-01	n	
				1.00E-02	I			V		1		Dithiane, 1,4-	505-29-3	7.80E+02	n	1.20E+04	n	2.00E+02	n	
				2.00E-03	I					1	0.1	Diuron	330-54-1	1.30E+02	n	1.60E+03	n	3.60E+01	n	
				4.00E-03	I					1	0.1	Dodine	2439-10-3	2.50E+02	n	3.30E+03	n	8.00E+01	n	
				2.50E-02	I			V		1		EPTC	759-94-4	2.00E+03	n	2.90E+04	n	3.80E+02	n	
				6.00E-03	I			V		1		Endosulfan	115-29-7	4.70E+02	n	7.00E+03	n	1.00E+02	n	
9.90E-03	I	1.20E-06	I	2.00E-02	I	3.00E-04	I			1	0.1	Endothall	145-73-3	1.30E+03	n	1.60E+04	n	3.80E+02	n	1.0E+02
				6.00E-03	P	1.00E-03	I	V		1	1.05E+04	Endrin	72-20-8	1.90E+01	n	2.50E+02	n	2.30E+00	n	2.0E+00
										1		Epichlorohydrin	106-89-8	1.90E+01	n	8.20E+01	n	2.00E+00	n	
				2.00E-02	I			V		1	1.53E+04	Epoxybutane, 1,2-	106-88-7	1.60E+02	n	6.70E+02	n	4.20E+01	n	
				4.00E-02	P					1	0.1	Ethanol, 2-(2-methoxyethoxy)-	111-77-3	2.50E+03	n	3.30E+04	n	8.00E+02	n	
				5.00E-03	I					1	0.1	Ethephon	16672-87-0	3.20E+02	n	4.10E+03	n	1.00E+02	n	
				5.00E-04	I					1	0.1	Ethion	563-12-2	3.20E+01	n	4.10E+02	n	4.30E+00	n	
				1.00E-01	P	6.00E-02	P	V		1	2.38E+04	Ethoxyethanol Acetate, 2-	111-15-9	2.60E+03	n	1.40E+04	n	1.20E+02	n	
				9.00E-02	P	2.00E-01	I	V		1	1.06E+05	Ethoxyethanol, 2-	110-80-5	5.20E+03	n	4.70E+04	n	3.40E+02	n	
				9.00E-01	I	7.00E-02	P	V		1	1.08E+04	Ethyl Acetate	141-78-6	6.20E+02	n	2.60E+03	n	1.40E+02	n	
				5.00E-03	P	8.00E-03	P	V		1	2.50E+03	Ethyl Acrylate	140-88-5	4.70E+01	n	2.10E+02	n	1.40E+01	n	
				1.00E+01	I	V				1	2.12E+03	Ethyl Chloride (Chloroethane)	75-00-3	1.40E+04	ns	5.70E+04	ns	2.10E+04	n	
				2.00E-01	I			V		1	1.01E+04	Ethyl Ether	60-29-7	1.60E+04	ns	2.30E+05	nms	3.90E+03	n	
				1.00E-05	I	3.00E-01	P	V		1	1.10E+03	Ethyl Methacrylate	97-63-2	1.80E+03	ns	7.60E+03	ns	6.30E+02	n	
1.10E-02	C	2.50E-06	C	1.00E-01	I	1.00E+00	I	V		1	4.80E+02	Ethyl-p-nitrophenyl Phosphonate	2104-64-5	6.30E-01	n	8.20E+00	n	8.90E-02	n	
				7.00E-02	P					1	0.1	Ethylbenzene	100-41-4	5.80E+02	c**s	2.50E+03	c**s	1.50E+02	c**	7.0E+02
				9.00E-02	P			V		1	1.89E+05	Ethylene Cyanohydrin	109-78-4	4.40E+03	n	5.70E+04	n	1.40E+03	n	
										1		Ethylene Diamine	107-15-3	7.00E+03	n	1.10E+05	nm	1.80E+03	n	
3.10E-01	C	8.80E-05	C	2.00E+00	I	4.00E-01	C			1	0.1	Ethylene Glycol	107-21-1	1.30E+05	nm	1.60E+06	nm	4.00E+04	n	
4.50E-02	C	1.30E-05	C	1.00E-01	I	1.60E+00	I			1	0.1	Ethylene Glycol Monobutyl Ether	111-76-2	6.30E+03	n	8.20E+04	n	2.00E+03	n	
6.50E+01	C	1.90E-02	C	8.00E-05	I	3.00E-02	C	V		1	1.21E+05	Ethylene Oxide	75-21-8	1.80E+01	c*	7.90E+01	c*	5.10E+00	c*	
										1	0.1	Ethylene Thiourea	96-45-7	5.10E+00	n	6.60E+01	n	1.60E+00	n	
				3.00E+00	I			V		1	1.54E+05	Ethyleneimine	151-56-4	2.70E-01	c	1.20E+00	c	2.40E-02	c	
										1	0.1	Ethylphthalyl Ethyl Glycolate	84-72-0	1.90E+05	nm	2.50E+06	nm	5.80E+04	n	
				2.50E-04	I					1	0.1	Fenamiphos	22224-92-6	1.60E+01	n	2.10E+02	n	4.40E+00	n	
				2.50E-02	I					1	0.1	Fenpropathrin	39515-41-8	1.60E+03	n	2.10E+04	n	6.40E+01	n	
				2.50E-02	I					1	0.1	Fenvalerate	51630-58-1	1.60E+03	n	2.10E+04	n	5.00E+02	n	
				1.30E-02	I					1	0.1	Fluometuron	2164-17-2	8.20E+02	n	1.10E+04	n	2.40E+02	n	
				4.00E-02	C	1.30E-02	C			1		Fluoride	16984-48-8	3.10E+03	n	4.70E+04	n	8.00E+02	n	
				6.00E-02	I	1.30E-02	C			1		Fluorine (Soluble Fluoride)	7782-41-4	4.70E+03	n	7.00E+04	n	1.20E+03	n	4.0E+03
				8.00E-02	I					1	0.1	Fluridone	59756-60-4	5.10E+03	n	6.60E+04	n	1.40E+03	n	
				2.00E-02	I					1	0.1	Flurprimidol	56425-91-3	1.30E+03	n	1.60E+04	n	3.40E+02	n	
				7.00E-04	I					1	0.1	Flusilazole	85509-19-9	4.40E+01	n	5.70E+02	n	1.10E+01	n	
				6.00E-02	I					1	0.1	Flutolanil	66332-96-5	3.80E+03	n	4.90E+04	n	9.50E+02	n	
				1.00E-02	I					1	0.1	Fluvalinate	69409-94-5	6.30E+02	n	8.20E+03	n	2.00E+02	n	

Regional Removal Management Level (RML) Summary Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
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 _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
3.50E-03	I			1.00E-01	I				1	0.1		Folpet	133-07-3	6.30E+03	n	6.60E+04	c**	1.80E+03	n	
1.90E-01	I			2.00E-03	I				1	0.1		Fomesafen	72178-02-0	2.90E+02	c	1.20E+03	c	3.90E+01	c	
		1.30E-05	I	2.00E-01	I	9.80E-03	A	V	1		4.24E+04	Fonofos	944-22-9	1.30E+02	n	1.60E+03	n	2.40E+01	n	
				9.00E-01	P	3.00E-04	X	V	1		1.06E+05	Formaldehyde	50-00-0	7.60E+02	n	3.30E+03	n	2.00E+01	n	
				3.00E+00	I				1	0.1		Formic Acid	64-18-6	2.90E+01	n	1.20E+02	n	6.30E-01	n	
									1	0.1		Fosetyl-AL	39148-24-8	1.90E+05	nm	2.50E+06	nm	6.00E+04	n	
									1	0.03		Furans								
				1.00E-03	X			V	1	0.03		~Dibenzofuran	132-64-9	7.30E+01	n	1.00E+03	n	7.90E+00	n	
				1.00E-03	I			V	1	0.03	6.22E+03	~Furan	110-00-9	7.30E+01	n	1.00E+03	n	1.90E+01	n	
				9.00E-01	I	2.00E+00	I	V	1	0.03	1.65E+05	~Tetrahydrofuran	109-99-9	1.80E+04	n	9.40E+04	n	3.40E+03	n	
3.80E+00	H			3.00E-03	I	5.00E-02	H	V	1	0.1		Furazolidone	67-45-8	1.40E+01	c	6.00E+01	c	2.00E+00	c	
1.50E+00	C	4.30E-04	C						1	0.1	1.01E+04	Furfural	98-01-1	2.10E+02	n	2.60E+03	n	3.80E+01	n	
									1	0.1		Furium	531-82-8	3.60E+01	c	1.50E+02	c	5.10E+00	c	
3.00E-02	I	8.60E-06	C						1	0.1		Furmecyclo	60568-05-0	1.80E+03	c	7.70E+03	c	1.10E+02	c	
				4.00E-04	I				1	0.1		Glufosinate, Ammonium	77182-82-2	2.50E+01	n	3.30E+02	n	8.00E+00	n	
						8.00E-05	C		1	0.1		Glutaraldehyde	111-30-8	1.10E+05	nm	4.80E+05	nm			
				4.00E-04	I	1.00E-03	H	V	1		1.06E+05	Glycidyl	765-34-4	2.30E+01	n	2.10E+02	n	1.70E+00	n	
				1.00E-01	I				1	0.1		Glyphosate	1071-83-6	6.30E+03	n	8.20E+04	n	2.00E+03	n	7.0E+02
				1.00E-02	X			V	1			Guanidine	113-00-8	7.80E+02	n	1.20E+04	n	2.00E+02	n	
				2.00E-02	P				1	0.1		Guanidine Chloride	50-01-1	1.30E+03	n	1.60E+04	n	4.00E+02	n	
4.50E+00	I	1.30E-03	I	5.00E-05	I				1	0.1		Haloxyfop, Methyl	69806-40-2	3.20E+00	n	4.10E+01	n	7.60E-01	n	
				5.00E-04	I			V	1			Heptachlor	76-44-8	1.30E+01	c**	6.30E+01	c**	1.40E-01	c**	4.0E-01
9.10E+00	I	2.60E-03	I	1.30E-05	I			V	1			Heptachlor Epoxide	1024-57-3	1.00E+00	n	1.50E+01	n	1.20E-01	n	2.0E-01
				2.00E-03	I			V	1			Hexabromobenzene	87-82-1	1.60E+02	n	2.30E+03	n	4.00E+01	n	
				2.00E-04	I				1	0.1		Hexabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-153)	68631-49-2	1.30E+01	n	1.60E+02	n	4.00E+00	n	
1.60E+00	I	4.60E-04	I	8.00E-04	I			V	1		1.68E+01	Hexachlorobenzene	118-74-1	2.10E+01	c**	9.60E+01	c**	9.80E-01	c*	1.0E+00
7.80E-02	I	2.20E-05	I	1.00E-03	P			V	1			Hexachlorobutadiene	87-68-3	7.80E+01	ns	5.30E+02	c**s	6.50E+00	n	
6.30E+00	I	1.80E-03	I	8.00E-03	A				1	0.1		Hexachlorocyclohexane, Alpha-	319-84-6	8.60E+00	c*	3.60E+01	c	7.20E-01	c	
1.80E+00	I	5.30E-04	I						1	0.1		Hexachlorocyclohexane, Beta-	319-85-7	3.00E+01	c	1.30E+02	c	2.50E+00	c	
1.10E+00	C	3.10E-04	C	3.00E-04	I				1	0.04		Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	2.10E+01	n	2.50E+02	c**	3.60E+00	n	2.0E-01
1.80E+00	I	5.10E-04	I						1	0.1		Hexachlorocyclohexane, Technical	608-73-1	3.00E+01	c	1.30E+02	c	2.50E+00	c	
4.00E-02	I	1.10E-05	C	6.00E-03	I	2.00E-04	I	V	1		1.57E+01	Hexachlorocyclopentadiene	77-47-4	1.80E+00	n	7.50E+00	n	4.10E-01	n	5.0E+01
				7.00E-04	I	3.00E-02	I	V	1			Hexachloroethane	67-72-1	4.50E+01	n	4.60E+02	n	6.20E+00	n	
				3.00E-04	I				1	0.1		Hexachlorophene	70-30-4	1.90E+01	n	2.50E+02	n	6.00E+00	n	
1.10E-01	I			3.00E-03	I				1	0.015		Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	2.30E+02	n	2.80E+03	c**	6.00E+01	n	
						1.00E-05	I	V	1		3.39E+03	Hexamethylene Diisocyanate, 1,6-	822-06-0	3.10E+00	n	1.30E+01	n	2.10E-02	n	
				4.00E-04	P				1	0.1		Hexamethylphosphoramide	680-31-9	2.50E+01	n	3.30E+02	n	8.00E+00	n	
						7.00E-01	I	V	1		1.41E+02	Hexane, N-	110-54-3	6.10E+02	ns	2.50E+03	ns	1.50E+03	n	
				2.00E+00	P				1	0.1		Hexanedioic Acid	124-04-9	1.30E+05	nm	1.60E+06	nm	4.00E+04	n	
				5.00E-03	I	3.00E-02	I	V	1		3.28E+03	Hexanone, 2-	591-78-6	2.00E+02	n	1.30E+03	n	3.80E+01	n	
				3.30E-02	I				1	0.1		Hexazinone	51235-04-2	2.10E+03	n	2.70E+04	n	6.40E+02	n	
				2.50E-02	I				1	0.1		Hexythiazox	78587-05-0	1.60E+03	n	2.10E+04	n	1.10E+02	n	
				3.00E-04	I				1	0.1		Hydramethylnon	67485-29-4	1.90E+01	n	2.50E+02	n	5.90E+00	n	
3.00E+00	I	4.90E-03	I			3.00E-05	P	V	1			Hydrazine	302-01-2	2.30E+01	c	1.10E+02	c	6.30E-02	n	
3.00E+00	I	4.90E-03	I						1			Hydrazine Sulfate	10034-93-2	2.30E+01	c	1.10E+02	c	2.60E+00	c	
						2.00E-02	I	V	1			Hydrogen Chloride	7647-01-0	2.80E+07	nm	1.20E+08	nm	4.20E+01	n	
				4.00E-02	C	1.40E-02	C	V	1			Hydrogen Fluoride	7664-39-3	3.10E+03	n	4.70E+04	n	2.80E+01	n	
6.00E-02	P			4.00E-02	P	2.00E-03	I	V	1	0.1		Hydrogen Sulfide	7783-06-4	2.80E+06	nm	1.20E+07	nm	4.20E+00	n	
									1	0.1		Hydroquinone	123-31-9	9.00E+02	c**	3.80E+03	c**	1.30E+02	c**	
				1.30E-02	I				1	0.1		Imazalil	35554-44-0	8.20E+02	n	1.10E+04	n	1.90E+02	n	
				2.50E-01	I				1	0.1		Imazaquin	81335-37-7	1.60E+04	n	2.10E+05	nm	4.90E+03	n	
				2.50E-01	I				1	0.1		Imazethapyr	81335-77-5	1.60E+04	n	2.10E+05	nm	4.70E+03	n	
				1.00E-02	A				1			Iodine	7553-56-2	7.80E+02	n	1.20E+04	n	2.00E+02	n	
				4.00E-02	I				1	0.1		Iprodione	36734-19-7	2.50E+03	n	3.30E+04	n	7.40E+02	n	

Regional Removal Management Level (RML) Summary Table (TR=1E-04, HQ=1) May 2016

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Toxicity and Chemical-specific Information											Contaminant		Removal Levels								
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)	
				7.00E-01	P				1			Iron	7439-89-6	5.50E+04	n	8.20E+05	nm	1.40E+04	n		
9.50E-04	I			3.00E-01	I		V		1		1.00E+04	Isobutyl Alcohol	78-83-1	2.30E+04	ns	3.50E+05	nms	5.90E+03	n		
				2.00E-01	I	2.00E+00	C		1	0.1		Isophorone	78-59-1	1.30E+04	c	1.60E+05	nm	3.80E+03	n		
				1.50E-02	I		V		1			Isopropalin	33820-53-0	1.20E+03	n	1.80E+04	n	4.00E+01	n		
				2.00E+00	P	2.00E-01	P	V	1		1.09E+05	Isopropanol	67-63-0	5.60E+03	n	2.40E+04	n	4.10E+02	n		
				1.00E-01	I				1	0.1		Isopropyl Methyl Phosphonic Acid	1832-54-8	6.30E+03	n	8.20E+04	n	2.00E+03	n		
				5.00E-02	I				1	0.1		Isoxaben	82558-50-7	3.20E+03	n	4.10E+04	n	7.30E+02	n		
						3.00E-01	A	V	1			JP-7	NA	4.30E+08	nm	1.80E+09	nm	6.30E+02	n		
				2.00E-03	I				1	0.1		Lactofen	77501-63-4	1.30E+02	n	1.60E+03	n	2.50E+01	n		
												Lead Compounds									
5.00E-01	C	1.50E-01	C	2.00E-02	C	2.00E-04	C	M	0.025			~Lead Chromate	7758-97-6	3.00E+01	c*	6.20E+02	c*	4.10E+00	c*		
8.50E-03	C	1.20E-05	C						1			~Lead Phosphate	7446-27-7	8.20E+03	c	3.80E+04	c	9.10E+02	c		
8.50E-03	C	1.20E-05	C						1	0.1		~Lead acetate	301-04-2	6.40E+03	c	2.70E+04	c	9.20E+02	c		
									1			~Lead and Compounds	7439-92-1	4.00E+02	L	8.00E+02	L	1.50E+01	L	1.5E+01	
8.50E-03	C	1.20E-05	C						1	0.1		~Lead subacetate	1335-32-6	6.40E+03	c	2.70E+04	c	9.20E+02	c		
				1.00E-07	I		V		1		2.43E+00	~Tetraethyl Lead	78-00-2	7.80E-03	n	1.20E-01	n	1.30E-03	n		
				5.00E-06	P		V		1		3.83E+02	Lewisite	541-25-3	3.90E-01	n	5.80E+00	n	9.00E-02	n		
				2.00E-03	I				1	0.1		Linuron	330-55-2	1.30E+02	n	1.60E+03	n	3.30E+01	n		
				2.00E-03	P				1			Lithium	7439-93-2	1.60E+02	n	2.30E+03	n	4.00E+01	n		
				5.00E-04	I				1	0.1		MCPA	94-74-6	3.20E+01	n	4.10E+02	n	7.50E+00	n		
				1.00E-02	I				1	0.1		MCPB	94-81-5	6.30E+02	n	8.20E+03	n	1.50E+02	n		
				1.00E-03	I				1	0.1		MCPB	93-65-2	6.30E+01	n	8.20E+02	n	1.60E+01	n		
				2.00E-02	I				1	0.1		Malathion	121-75-5	1.30E+03	n	1.60E+04	n	3.90E+02	n		
				1.00E-01	I	7.00E-04	C		1	0.1		Maleic Anhydride	108-31-6	6.30E+03	n	8.00E+04	n	1.90E+03	n		
				5.00E-01	I				1	0.1		Maleic Hydrazide	123-33-1	3.20E+04	n	4.10E+05	nm	1.00E+04	n		
				1.00E-04	P				1	0.1		Malononitrile	109-77-3	6.30E+00	n	8.20E+01	n	2.00E+00	n		
				3.00E-02	H				1	0.1		Mancozeb	8018-01-7	1.90E+03	n	2.50E+04	n	5.40E+02	n		
				5.00E-03	I				1	0.1		Maneb	12427-38-2	3.20E+02	n	4.10E+03	n	9.80E+01	n		
				1.40E-01	I	5.00E-05	I		1			Manganese (Diet)	7439-96-5								
				2.40E-02	S	5.00E-05	I		0.04			Manganese (Non-diet)	7439-96-5	1.80E+03	n	2.60E+04	n	4.30E+02	n		
				9.00E-05	H				1	0.1		Mephosolan	950-10-7	5.70E+00	n	7.40E+01	n	1.80E+00	n		
				3.00E-02	I				1	0.1		Mepiquat Chloride	24307-26-4	1.90E+03	n	2.50E+04	n	6.00E+02	n		
												Mercury Compounds									
				3.00E-04	I	3.00E-04	S		0.07			~Mercuric Chloride (and other Mercury salts)	7487-94-7	2.30E+01	n	3.50E+02	n	5.70E+00	n	2.0E+00	
						3.00E-04	I	V	1		3.13E+00	~Mercury (elemental)	7439-97-6	1.10E+01	ns	4.60E+01	ns	6.30E-01	n	2.0E+00	
				1.00E-04	I				1			~Methyl Mercury	22967-92-6	7.80E+00	n	1.20E+02	n	2.00E+00	n		
				8.00E-05	I				1	0.1		~Phenylmercuric Acetate	62-38-4	5.10E+00	n	6.60E+01	n	1.60E+00	n		
				3.00E-05	I		V		1			Merphos	150-50-5	2.30E+00	n	3.50E+01	n	6.00E-01	n		
				3.00E-05	I				1	0.1		Merphos Oxide	78-48-8	1.90E+00	n	2.50E+01	n	8.50E-02	n		
				6.00E-02	I				1	0.1		Metaxyl	57837-19-1	3.80E+03	n	4.90E+04	n	1.20E+03	n		
				1.00E-04	I	3.00E-02	P	V	1		4.58E+03	Methacrylonitrile	126-98-7	7.50E+00	n	1.00E+02	n	1.90E+00	n		
				5.00E-05	I				1	0.1		Methamidophos	10265-92-6	3.20E+00	n	4.10E+01	n	1.00E+00	n		
				2.00E+00	I	2.00E+01	I	V	1		1.06E+05	Methanol	67-56-1	1.20E+05	nms	1.20E+06	nms	2.00E+04	n		
				1.00E-03	I				1	0.1		Methidathion	950-37-8	6.30E+01	n	8.20E+02	n	1.90E+01	n		
4.90E-02	C	1.40E-05	C	2.50E-02	I				1	0.1		Methomyl	16752-77-5	1.60E+03	n	2.10E+04	n	5.00E+02	n		
									1	0.1		Methoxy-5-nitroaniline, 2-	99-59-2	1.10E+03	c	4.70E+03	c	1.50E+02	c		
				5.00E-03	I				1	0.1		Methoxychlor	72-43-5	3.20E+02	n	4.10E+03	n	3.70E+01	n	4.0E+01	
				8.00E-03	P	1.00E-03	P	V	1		1.15E+05	Methoxyethanol Acetate, 2-	110-49-6	1.10E+02	n	5.10E+02	n	2.10E+00	n		
				5.00E-03	P	2.00E-02	I	V	1		1.06E+05	Methoxyethanol, 2-	109-86-4	3.30E+02	n	3.50E+03	n	2.90E+01	n		
				1.00E+00	X		V		1		2.90E+04	Methyl Acetate	79-20-9	7.80E+04	ns	1.20E+06	nms	2.00E+04	n		
						2.00E-02	P	V	1		6.75E+03	Methyl Acrylate	96-33-3	1.50E+02	n	6.10E+02	n	4.20E+01	n		
				6.00E-01	I	5.00E+00	I	V	1		2.84E+04	Methyl Ethyl Ketone (2-Butanone)	78-93-3	2.70E+04	n	1.90E+05	nms	5.60E+03	n		
				1.00E-03	X	1.00E-03	P	2.00E-05	X	V	1	1.80E+05	Methyl Hydrazine	60-34-4	1.00E+00	n	4.40E+00	n	4.20E-02	n	
						3.00E+00	I	V	1		3.36E+03	Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1	3.30E+04	ns	1.40E+05	nms	6.30E+03	n		

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Toxicity and Chemical-specific Information											Contaminant		Removal Levels								
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 _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)	
				1.00E-03	C V				1		1.01E+04	Methyl Isocyanate	624-83-9	4.60E+00	n	1.90E+01	n	2.10E+00	n		
				1.40E+00	I	7.00E-01	I	V	1		2.36E+03	Methyl Methacrylate	80-62-6	4.40E+03	ns	1.90E+04	ns	1.40E+03	n		
				2.50E-04	I				1	0.1		Methyl Parathion	298-00-0	1.60E+01	n	2.10E+02	n	4.50E+00	n		
				6.00E-02	X				1	0.1		Methyl Phosphonic Acid	993-13-5	3.80E+03	n	4.90E+04	n	1.20E+03	n		
9.90E-02	C	2.80E-05	C	6.00E-03	H	4.00E-02	H	V	1		3.93E+02	Methyl Styrene (Mixed Isomers)	25013-15-4	3.20E+02	n	2.60E+03	ns	2.30E+01	n		
1.80E-03	C	2.60E-07	C						1	0.1		Methyl methanesulfonate	66-27-3	5.50E+02	c	2.30E+03	c	7.90E+01	c		
				3.00E+00	I	V			1		8.87E+03	Methyl tert-Butyl Ether (MTBE)	1634-04-4	4.70E+03	c**	2.10E+04	c**s	1.40E+03	c**		
9.00E-03	P			3.00E-04	X				1	0.1		Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2	1.90E+01	n	2.50E+02	n	6.00E+00	n		
8.30E+00	C	2.40E-03	C	2.00E-02	X				1	0.1		Methyl-5-Nitroaniline, 2-	99-55-8	1.30E+03	n	1.60E+04	n	3.80E+02	n		
									1	0.1		Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	6.50E+00	c	2.80E+01	c	9.40E-01	c		
1.30E-01	C	3.70E-05	C	1.00E-02	A				1	0.1		Methylaniline Hydrochloride, 2-	636-21-5	4.20E+02	c	1.80E+03	c	6.00E+01	c		
				2.00E-04	X				1	0.1		Methylarsonic acid	124-58-3	6.30E+02	n	8.20E+03	n	2.00E+02	n		
									1	0.1		Methylbenzene,1,4-diamine monohydrochloride, 2-	74612-12-7	1.30E+01	n	1.60E+02	n	4.00E+00	n		
1.00E-01	X			3.00E-04	X				1	0.1		Methylbenzene-1,4-diamine sulfate, 2-	615-50-9	1.90E+01	n	2.50E+02	n	6.00E+00	n		
2.20E+01	C	6.30E-03	C					M	1	0.1		Methylcholanthrene, 3-	56-49-5	5.50E-01	c	1.00E+01	c	1.10E-01	c		
2.00E-03	I	1.00E-08	I	6.00E-03	I	6.00E-01	I	V	M	1	3.32E+03	Methylene Chloride	75-09-2	3.50E+02	n	3.20E+03	n	1.10E+02	n	5.0E+00	
1.00E-01	P	4.30E-04	C	2.00E-03	P				M	1	0.1	Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	1.20E+02	c**	1.60E+03	n	1.60E+01	c**		
4.60E-02	I	1.30E-05	C						1	0.1		Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	1.20E+03	c	5.00E+03	c	4.80E+01	c		
1.60E+00	C	4.60E-04	C			2.00E-02	C		1	0.1		Methylenebisbenzenamine, 4,4'-	101-77-9	3.40E+01	c	1.40E+02	c	4.70E+00	c		
				6.00E-04	I				1	0.1		Methylenediphenyl Diisocyanate	101-68-8	8.50E+05	nm	3.60E+06	nm				
				7.00E-02	H			V	1		5.00E+02	Methylstyrene, Alpha-	98-83-9	5.50E+03	ns	8.20E+04	ns	7.80E+02	n		
				1.50E-01	I				1	0.1		Metolachlor	51218-45-2	9.50E+03	n	1.20E+05	nm	2.70E+03	n		
				2.50E-02	I				1	0.1		Metribuzin	21087-64-9	1.60E+03	n	2.10E+04	n	4.90E+02	n		
				2.50E-01	I				1	0.1		Metsulfuron-methyl	74223-64-6	1.60E+04	n	2.10E+05	nm	4.90E+03	n		
				3.00E+00	P			V	1		3.42E-01	Mineral oils	8012-95-1	2.30E+05	nms	3.50E+06	nms	6.00E+04	n		
1.80E+01	C	5.10E-03	C	2.00E-04	I			V	1			Mirex	2385-85-5	3.60E+00	c**	1.70E+01	c*	8.80E-02	c*		
				2.00E-03	I				1	0.1		Molinat	2212-67-1	1.30E+02	n	1.60E+03	n	3.00E+01	n		
				5.00E-03	I				1			Molybdenum	7439-98-7	3.90E+02	n	5.80E+03	n	1.00E+02	n		
				1.00E-01	I				1			Monochloramine	10599-90-3	7.80E+03	n	1.20E+05	nm	2.00E+03	n	4.0E+03	
				2.00E-03	P				1	0.1		Monomethylaniline	100-61-8	1.30E+02	n	1.60E+03	n	3.80E+01	n		
				2.50E-02	I				1	0.1		Myclobutanil	88671-89-0	1.60E+03	n	2.10E+04	n	4.50E+02	n		
				3.00E-04	X				1	0.1		N,N'-Diphenyl-1,4-benzenediamine	74-31-7	1.90E+01	n	2.50E+02	n	3.60E+00	n		
				2.00E-03	I			V	1			Naled	300-76-5	1.60E+02	n	2.30E+03	n	4.00E+01	n		
				3.00E-02	X	1.00E-01	P	V	1			Naphtha, High Flash Aromatic (HFAN)	64742-95-6	2.30E+03	n	3.50E+04	n	1.50E+02	n		
1.80E+00	C	0.00E+00	C						1	0.1		Naphthylamine, 2-	91-59-8	3.00E+01	c	1.30E+02	c	3.90E+00	c		
				1.00E-01	I				1	0.1		Napropamide	15299-99-7	6.30E+03	n	8.20E+04	n	1.60E+03	n		
				2.60E-04	C	1.10E-02	C	1.40E-05	C	1	0.1	Nickel Acetate	373-02-4	6.70E+02	n	8.10E+03	n	2.20E+02	n		
				2.60E-04	C	1.10E-02	C	1.40E-05	C	1	0.1	Nickel Carbonate	3333-67-3	6.70E+02	n	8.10E+03	n	2.20E+02	n		
				2.60E-04	C	1.10E-02	C	1.40E-05	C V	1		Nickel Carbonyl	13463-39-3	8.20E+02	n	1.10E+04	n	2.90E-02	n		
				2.60E-04	C	1.10E-02	C	1.40E-05	C	0.04		Nickel Hydroxide	12054-48-7	8.20E+02	n	1.10E+04	n	2.00E+02	n		
				2.60E-04	C	1.10E-02	C	2.00E-05	C	0.04		Nickel Oxide	1313-99-1	8.40E+02	n	1.20E+04	n	2.00E+02	n		
				2.40E-04	I	1.10E-02	C	1.40E-05	C	0.04		Nickel Refinery Dust	NA	8.20E+02	n	1.10E+04	n	2.20E+02	n		
				2.60E-04	C	2.00E-02	I	9.00E-05	A	0.04		Nickel Soluble Salts	7440-02-0	1.50E+03	n	2.20E+04	n	3.90E+02	n		
1.70E+00	C	4.80E-04	I	1.10E-02	C	1.40E-05	C		0.04			Nickel Subulfide	12035-72-2	4.10E+01	c*	1.90E+02	c*	4.50E+00	c*		
				2.60E-04	C	1.10E-02	C	1.40E-05	C	1	0.1	Nickelocene	1271-28-9	6.70E+02	n	8.10E+03	n	2.20E+02	n		
				1.60E+00	I				1			Nitrate	14797-55-8	1.30E+05	nm	1.90E+06	nm	3.20E+04	n	1.0E+04	
									1			Nitrate + Nitrite (as N)	NA							1.0E+04	
				1.00E-01	I				1			Nitrite	14797-65-0	7.80E+03	n	1.20E+05	nm	2.00E+03	n	1.0E+03	
				1.00E-02	X	5.00E-05	X		1	0.1		Nitroaniline, 2-	88-74-4	6.30E+02	n	8.00E+03	n	1.90E+02	n		
2.00E-02	P			4.00E-03	P	6.00E-03	P		1	0.1		Nitroaniline, 4-	100-01-6	2.50E+02	n	3.30E+03	n	7.80E+01	n		
				4.00E-05	I	2.00E-03	I	9.00E-03	I	V	1	3.05E+03	Nitrobenzene	98-95-3	1.30E+02	n	1.30E+03	n	1.30E+01	n	
				3.00E+03	P				1	0.1		Nitrocellulose	9004-70-0	1.90E+08	nm	2.50E+09	nm	6.00E+07	n		
				7.00E-02	H				1	0.1		Nitrofurantoin	67-20-9	4.40E+03	n	5.70E+04	n	1.40E+03	n		
1.30E+00	C	3.70E-04	C						1	0.1		Nitrofurazone	59-87-0	4.20E+01	c	1.80E+02	c	6.00E+00	c		

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Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
1.70E-02	P			1.00E-04	P				1	0.1		Nitroglycerin	55-63-0	6.30E+00	n	8.20E+01	n	2.00E+00	n	
		8.80E-06	P	1.00E-01	I	5.00E-03	P	V	1		1.80E+04	Nitroguanidine	556-88-7	6.30E+03	n	8.20E+04	n	2.00E+03	n	
		2.70E-03	H			2.00E-02	I	V	1		4.86E+03	Nitromethane	75-52-5	8.80E+01	n	3.70E+02	n	1.00E+01	n	
												Nitropropane, 2-	79-46-9	1.40E+00	c	6.00E+00	c	2.10E-01	c	
2.70E+01	C	7.70E-03	C					M	1	0.1		Nitroso-N-ethylurea, N-	759-73-9	4.50E-01	c	8.50E+00	c	9.20E-02	c	
1.20E+02	C	3.40E-02	C					M	1	0.1		Nitroso-N-methylurea, N-	684-93-5	1.00E-01	c	1.90E+00	c	2.10E-02	c	
5.40E+00	I	1.60E-03	I					V	1			Nitroso-di-N-butylamine, N-	924-16-3	9.90E+00	c	4.60E+01	c	2.70E-01	c	
7.00E+00	I	2.00E-03	C						1	0.1		Nitroso-di-N-propylamine, N-	621-64-7	7.80E+00	c	3.30E+01	c	1.10E+00	c	
2.80E+00	I	8.00E-04	C						1	0.1		Nitrosodiethanolamine, N-	1116-54-7	1.90E+01	c	8.20E+01	c	2.80E+00	c	
1.50E+02	I	4.30E-02	I					M	1	0.1		Nitrosodiethylamine, N-	55-18-5	8.10E-02	c	1.50E+00	c	1.70E-02	c	
5.10E+01	I	1.40E-02	I	8.00E-06	P	4.00E-05	X	V	M	1	2.37E+05	Nitrosodimethylamine, N-	62-75-9	2.00E-01	c**	3.40E+00	c**	1.10E-02	c**	
4.90E-03	I	2.60E-06	C						1	0.1		Nitrosodiphenylamine, N-	86-30-6	1.10E+04	c	4.70E+04	c	1.20E+03	c	
2.20E+01	I	6.30E-03	C					V	1		1.08E+05	Nitrosomethylethylamine, N-	10595-95-6	2.00E+00	c	9.10E+00	c	7.10E-02	c	
6.70E+00	C	1.90E-03	C						1	0.1		Nitrosomorpholine [N-]	59-89-2	8.10E+00	c	3.40E+01	c	1.20E+00	c	
9.40E+00	C	2.70E-03	C						1	0.1		Nitrosopiperidine [N-]	100-75-4	5.80E+00	c	2.40E+01	c	8.20E-01	c	
2.10E+00	I	6.10E-04	I						1	0.1		Nitrosopyrrolidine, N-	930-55-2	2.60E+01	c	1.10E+02	c	3.70E+00	c	
2.20E-01	P			1.00E-04	X				1	0.1		Nitrotoluene, m-	99-08-1	6.30E+00	n	8.20E+01	n	1.70E+00	n	
1.60E-02	P			9.00E-04	P			V	1		1.51E+03	Nitrotoluene, o-	88-72-2	7.00E+01	n	1.10E+03	n	1.60E+01	n	
				4.00E-03	P				1	0.1		Nitrotoluene, p-	99-99-0	2.50E+02	n	3.30E+03	n	7.10E+01	n	
				3.00E-04	X	2.00E-02	P	V	1		6.86E+00	Nonane, n-	111-84-2	1.10E+01	ns	7.20E+01	ns	5.30E+00	n	
				4.00E-02	I				1	0.1		Norflurazon	27314-13-2	2.50E+03	n	3.30E+04	n	7.70E+02	n	
				3.00E-03	I				1	0.1		Octabromodiphenyl Ether	32536-52-0	1.90E+02	n	2.50E+03	n	6.00E+01	n	
				5.00E-02	I				1	0.006		Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0	3.90E+03	n	5.70E+04	n	1.00E+03	n	
				2.00E-03	H				1	0.1		Octamethylpyrophosphoramide	152-16-9	1.30E+02	n	1.60E+03	n	4.00E+01	n	
				5.00E-02	I				1	0.1		Oryzalin	19044-88-3	3.20E+03	n	4.10E+04	n	8.10E+02	n	
				5.00E-03	I				1	0.1		Oxadiazon	19666-30-9	3.20E+02	n	4.10E+03	n	4.70E+01	n	
				2.50E-02	I				1	0.1		Oxamyl	23135-22-0	1.60E+03	n	2.10E+04	n	5.00E+02	n	2.0E+02
				3.00E-03	I				1	0.1		Oxyfluorfen	42874-03-3	1.90E+02	n	2.50E+03	n	3.20E+01	n	
				1.30E-02	I				1	0.1		Paclitaxel	76738-62-0	8.20E+02	n	1.10E+04	n	2.30E+02	n	
				4.50E-03	I				1	0.1		Paraquat Dichloride	1910-42-5	2.80E+02	n	3.70E+03	n	9.00E+01	n	
				6.00E-03	H				1	0.1		Parathion	56-38-2	3.80E+02	n	4.90E+03	n	8.60E+01	n	
				5.00E-02	H			V	1			Pebulate	1114-71-2	3.90E+03	n	5.80E+04	n	5.60E+02	n	
				4.00E-02	I				1	0.1		Pendimethalin	40487-42-1	2.50E+03	n	3.30E+04	n	1.80E+02	n	
				2.00E-03	I			V	1		3.12E-01	Pentabromodiphenyl Ether	32534-81-9	1.60E+02	ns	2.30E+03	ns	4.00E+01	n	
9.00E-02	P			1.00E-04	I				1	0.1		Pentabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-99)	60348-60-9	6.30E+00	n	8.20E+01	n	2.00E+00	n	
				8.00E-04	I			V	1		4.57E+02	Pentachlorobenzene	608-93-5	6.30E+01	n	9.30E+02	n	3.20E+00	n	
								V	1			Pentachloroethane	76-01-7	7.70E+02	cs	3.60E+03	cs	6.50E+01	c	
2.60E-01	H			3.00E-03	I			V	1			Pentachloronitrobenzene	82-68-8	2.30E+02	n	1.30E+03	c**	1.20E+01	c**	
4.00E-01	I	5.10E-06	C	5.00E-03	I				1	0.25		Pentachlorophenol	87-86-5	1.00E+02	c**	4.00E+02	c**	4.10E+00	c**	1.0E+00
4.00E-03	X			2.00E-03	P				1	0.1		Pentaerythritol tetranitrate (PETN)	78-11-5	1.30E+02	n	1.60E+03	n	3.90E+01	n	
				1.00E+00	P	V			1		3.88E+02	Pentane, n-	109-66-0	8.10E+02	ns	3.40E+03	ns	2.10E+03	n	
				7.00E-04	I				1			Perchlorates								
				7.00E-04	I				1			~Ammonium Perchlorate	7790-98-9	5.50E+01	n	8.20E+02	n	1.40E+01	n	
				7.00E-04	I				1			~Lithium Perchlorate	7791-03-9	5.50E+01	n	8.20E+02	n	1.40E+01	n	
				7.00E-04	I				1			~Perchlorate and Perchlorate Salts	14797-73-0	5.50E+01	n	8.20E+02	n	1.40E+01	n	1.5E+01(F)
				7.00E-04	I				1			~Potassium Perchlorate	7778-74-7	5.50E+01	n	8.20E+02	n	1.40E+01	n	
				7.00E-04	I				1			~Sodium Perchlorate	7601-89-0	5.50E+01	n	8.20E+02	n	1.40E+01	n	
				2.00E-02	P			V	1			Perfluorobutane Sulfonate	375-73-5	1.60E+03	n	2.30E+04	n	3.80E+02	n	
				5.00E-02	I				1	0.1		Permethrin	52645-53-1	3.20E+03	n	4.10E+04	n	1.00E+03	n	
2.20E-03	C	6.30E-07	C						1	0.1		Phenacetin	62-44-2	2.50E+04	c	1.00E+05	cm	3.40E+03	c	
				2.50E-01	I				1	0.1		Phenmedipham	13684-63-4	1.60E+04	n	2.10E+05	nm	4.00E+03	n	
				3.00E-01	I	2.00E-01	C		1	0.1		Phenol	108-95-2	1.90E+04	n	2.50E+05	nm	5.80E+03	n	
				4.00E-03	I				1	0.1		Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1	2.50E+02	n	3.30E+03	n	7.80E+01	n	
				5.00E-04	X				1	0.1		Phenothiazine	92-84-2	3.20E+01	n	4.10E+02	n	4.30E+00	n	

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Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
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 _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
				6.00E-03	I				1	0.1		Phenylenediamine, m-	108-45-2	3.80E+02	n	4.90E+03	n	1.20E+02	n	
4.70E-02	H			1.90E-01	H				1	0.1		Phenylenediamine, o-	95-54-5	1.20E+03	c	4.90E+03	c	1.60E+02	c	
1.90E-03	H								1	0.1		Phenylenediamine, p-	106-50-3	1.20E+04	n	1.60E+05	nm	3.80E+03	n	
									1	0.1		Phenylphenol, 2-	90-43-7	2.80E+04	c	1.20E+05	cm	3.00E+03	c	
				2.00E-04	H				1	0.1		Phorate	298-02-2	1.30E+01	n	1.60E+02	n	3.00E+00	n	
						3.00E-04	I	V	1		1.61E+03	Phosgene	75-44-5	3.10E-01	n	1.30E+00	n			
				2.00E-02	I				1	0.1		Phosmet	732-11-6	1.30E+03	n	1.60E+04	n	3.70E+02	n	
				4.90E+01	P				1			Phosphates, Inorganic								
				4.90E+01	P				1			~Aluminum metaphosphate	13776-88-0	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Ammonium polyphosphate	68333-79-9	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Calcium pyrophosphate	7790-76-3	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Diammonium phosphate	7783-28-0	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Dicalcium phosphate	7757-93-9	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Dimagnesium phosphate	7782-75-4	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Dipotassium phosphate	7758-11-4	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Disodium phosphate	7558-79-4	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Monoaluminum phosphate	13530-50-2	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Monoammonium phosphate	7722-76-1	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Monocalcium phosphate	7758-23-8	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Monomagnesium phosphate	7757-86-0	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Monopotassium phosphate	7778-77-0	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Monosodium phosphate	7558-80-7	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Polyphosphoric acid	8017-16-1	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Potassium triphosphate	13845-36-8	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Sodium acid pyrophosphate	7758-16-9	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Sodium aluminum phosphate (acidic)	7785-88-8	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Sodium aluminum phosphate (anhydrous)	10279-59-1	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Sodium aluminum phosphate (tetrahydrate)	10305-76-7	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Sodium hexametaphosphate	10124-56-8	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Sodium polyphosphate	68915-31-1	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Sodium trimetaphosphate	7785-84-4	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Sodium tripolyphosphate	7758-29-4	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Tetrapotassium phosphate	7320-34-5	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Tetrasodium pyrophosphate	7722-88-5	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Trialuminum sodium tetra decahydrogenoctaorthophosphate (dihydrate)	15136-87-5	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Tricalcium phosphate	7758-87-4	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Trimagnesium phosphate	7757-87-1	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Tripotassium phosphate	7778-53-2	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				4.90E+01	P				1			~Trisodium phosphate	7601-54-9	3.80E+06	nm	5.70E+07	nm	9.70E+05	n	
				3.00E-04	I	3.00E-04	I	V	1			Phosphine	7803-51-2	2.30E+01	n	3.50E+02	n	5.70E-01	n	
				4.90E+01	P	1.00E-02	I		1			Phosphoric Acid	7664-38-2	3.00E+06	nm	2.90E+07	nm	9.70E+05	n	
				2.00E-05	I			V	1			Phosphorus, White	7723-14-0	1.60E+00	n	2.30E+01	n	4.00E-01	n	
												Phthalates								
1.40E-02	I	2.40E-06	C	2.00E-02	I				1	0.1		~Bis(2-ethylhexyl)phthalate	117-81-7	1.30E+03	n	1.60E+04	c**	4.00E+02	n	6.0E+00
1.90E-03	P			2.00E-01	I				1	0.1		~Butyl Benzyl Phthalate	85-68-7	1.30E+04	n	1.20E+05	c**m	1.60E+03	c**	
				1.00E+00	I				1	0.1		~Butylphthalyl Butylglycolate	85-70-1	6.30E+04	n	8.20E+05	nm	1.30E+04	n	
				1.00E-01	I				1	0.1		~Dibutyl Phthalate	84-74-2	6.30E+03	n	8.20E+04	n	9.00E+02	n	
				8.00E-01	I				1	0.1		~Diethyl Phthalate	84-66-2	5.10E+04	n	6.60E+05	nm	1.50E+04	n	
				1.00E-01	I			V	1			~Dimethylterephthalate	120-61-6	7.80E+03	n	1.20E+05	nm	1.90E+03	n	
				1.00E-02	P				1	0.1		~Octyl Phthalate, di-N-	117-84-0	6.30E+02	n	8.20E+03	n	2.00E+02	n	
				1.00E+00	H				1	0.1		~Phthalic Acid, P-	100-21-0	6.30E+04	n	8.20E+05	nm	1.90E+04	n	
				2.00E+00	I	2.00E-02	C		1	0.1		~Phthalic Anhydride	85-44-9	1.30E+05	nm	1.60E+06	nm	3.90E+04	n	
				7.00E-02	I				1	0.1		Picloram	1918-02-1	4.40E+03	n	5.70E+04	n	1.40E+03	n	5.0E+02
				1.00E-04	X				1	0.1		Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3	6.30E+00	n	8.20E+01	n	2.00E+00	n	

Regional Removal Management Level (RML) Summary Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
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 _i (mg/m ³)	k e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
				9.00E-04	X				1	0.1		Picric Acid (2,4,6-Trinitrophenol)	88-89-1	5.70E+01	n	7.40E+02	n	1.80E+01	n	
3.00E+01	C	8.60E-03	C	1.00E-02	I				1	0.1		Pirimiphos, Methyl	29232-93-7	6.30E+02	n	8.20E+03	n	1.20E+02	n	
				7.00E-06	H				1	0.1		Polybrominated Biphenyls	59536-65-1	4.40E-01	n	5.70E+00	n	1.40E-01	n	
												Polychlorinated Biphenyls (PCBs)								
7.00E-02	S	2.00E-05	S	7.00E-05	I			V	1	0.14		~Aroclor 1016	12674-11-2	4.10E+00	n	5.10E+01	n	1.40E+00	n	
2.00E+00	S	5.70E-04	S					V	1	0.14		~Aroclor 1221	11104-28-2	2.00E+01	c	8.30E+01	c	4.70E-01	c	
2.00E+00	S	5.70E-04	S					V	1	0.14		~Aroclor 1232	11141-16-5	1.70E+01	c	7.20E+01	c	4.70E-01	c	
2.00E+00	S	5.70E-04	S					V	1	0.14		~Aroclor 1242	53469-21-9	2.30E+01	c	9.50E+01	c	7.80E-01	c	
2.00E+00	S	5.70E-04	S					V	1	0.14		~Aroclor 1248	12672-29-6	2.30E+01	c	9.50E+01	c	7.80E-01	c	
2.00E+00	S	5.70E-04	S	2.00E-05	I			V	1	0.14		~Aroclor 1254	11097-69-1	1.20E+00	n	1.50E+01	n	4.00E-01	n	
2.00E+00	S	5.70E-04	S					V	1	0.14		~Aroclor 1260	11096-82-5	2.40E+01	c	9.90E+01	c	7.80E-01	c	
3.90E+00	E	1.10E-03	E	6.00E-04	X			V	1	0.14		~Aroclor 5460	11126-42-4	3.50E+01	n	4.40E+02	n	1.20E+01	n	
				2.30E-05	E	1.30E-03	E	V	1	0.14		~Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	1.40E+00	n	1.70E+01	n	4.00E-01	c**	
3.90E+00	E	1.10E-03	E	2.30E-05	E	1.30E-03	E	V	1	0.14		~Hexachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 167)	52663-72-6	1.40E+00	n	1.70E+01	n	4.00E-01	c**	
3.90E+00	E	1.10E-03	E	2.30E-05	E	1.30E-03	E	V	1	0.14		~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	1.40E+00	n	1.70E+01	n	4.00E-01	c**	
3.90E+00	E	1.10E-03	E	2.30E-05	E	1.30E-03	E	V	1	0.14		~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 156)	38380-08-4	1.40E+00	n	1.70E+01	n	4.00E-01	c**	
3.90E+03	E	1.10E+00	E	2.30E-08	E	1.30E-06	E	V	1	0.14		~Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	1.40E-03	n	1.70E-02	n	4.00E-04	c**	
3.90E+00	E	1.10E-03	E	2.30E-05	E	1.30E-03	E	V	1	0.14		~Pentachlorobiphenyl, 2',3,4,4',5'- (PCB 123)	65510-44-3	1.40E+00	n	1.70E+01	n	4.00E-01	c**	
3.90E+00	E	1.10E-03	E	2.30E-05	E	1.30E-03	E	V	1	0.14		~Pentachlorobiphenyl, 2,3',4,4',5'- (PCB 118)	31508-00-6	1.40E+00	n	1.70E+01	n	4.00E-01	c**	
3.90E+00	E	1.10E-03	E	2.30E-05	E	1.30E-03	E	V	1	0.14		~Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	1.40E+00	n	1.70E+01	n	4.00E-01	c**	
3.90E+00	E	1.10E-03	E	2.30E-05	E	1.30E-03	E	V	1	0.14		~Pentachlorobiphenyl, 2,3,4,4',5'- (PCB 114)	74472-37-0	1.40E+00	n	1.70E+01	n	4.00E-01	c**	
1.30E+04	E	3.80E+00	E	7.00E-09	E	4.00E-07	E	V	1	0.14		~Pentachlorobiphenyl, 3,3',4,4',5'- (PCB 126)	57465-28-8	4.10E-04	n	5.10E-03	n	1.20E-04	c**	
2.00E+00	I	5.70E-04	I					V	1	0.14		~Polychlorinated Biphenyls (high risk)	1336-36-3	2.30E+01	c	9.40E+01	c			
4.00E-01	I	1.00E-04	I					V	1	0.14		~Polychlorinated Biphenyls (low risk)	1336-36-3					4.40E+00	c	
7.00E-02	I	2.00E-05	I					V	1	0.14		~Polychlorinated Biphenyls (lowest risk)	1336-36-3							5.0E-01
1.30E+01	E	3.80E-03	E	7.00E-06	E	4.00E-04	E	V	1	0.14		~Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	4.10E-01	n	5.10E+00	n	1.40E-01	n	
3.90E+01	E	1.10E-02	E	2.30E-06	E	1.30E-04	E	V	1	0.14		~Tetrachlorobiphenyl, 3,4,4',5'- (PCB 81)	70362-50-4	1.40E-01	n	1.70E+00	n	4.00E-02	c**	
				6.00E-04	I			V	1	0.1		Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9	8.50E+05	nm	3.60E+06	nm			
												Polynuclear Aromatic Hydrocarbons (PAHs)								
				6.00E-02	I			V	1	0.13		~Acenaphthene	83-32-9	3.60E+03	n	4.50E+04	n	5.30E+02	n	
				3.00E-01	I			V	1	0.13		~Anthracene	120-12-7	1.80E+04	n	2.30E+05	nm	1.80E+03	n	
7.30E-01	E	1.10E-04	C					V	M	1	0.13	~Benzo[a]anthracene	56-55-3	1.60E+01	c	2.90E+02	c	1.20E+00	c	
1.20E+00	C	1.10E-04	C							1	0.13	~Benzo[j]fluoranthene	205-82-3	4.20E+01	c	1.80E+02	c	6.50E+00	c	
7.30E+00	I	1.10E-03	C						M	1	0.13	~Benzo[a]pyrene	50-32-8	1.60E+00	c	2.90E+01	c	3.40E-01	c	2.0E-01
7.30E-01	E	1.10E-04	C						M	1	0.13	~Benzo[b]fluoranthene	205-99-2	1.60E+01	c	2.90E+02	c	3.40E+00	c	
7.30E-02	E	1.10E-04	C						M	1	0.13	~Benzo[k]fluoranthene	207-08-9	1.60E+02	c	2.90E+03	c	3.40E+01	c	
				8.00E-02	I			V	1	0.13		~Chloronaphthalene, Beta-	91-58-7	4.80E+03	n	6.00E+04	n	7.50E+02	n	
7.30E-03	E	1.10E-05	C						M	1	0.13	~Chrysene	218-01-9	1.60E+03	c	2.90E+04	c	3.40E+02	c	
7.30E+00	E	1.20E-03	C						M	1	0.13	~Dibenz[a,h]anthracene	53-70-3	1.60E+00	c	2.90E+01	c	3.40E-01	c	
1.20E+01	C	1.10E-03	C							1	0.13	~Dibenzo[a,e]pyrene	192-65-4	4.20E+00	c	1.80E+01	c	6.50E-01	c	
2.50E+02	C	7.10E-02	C						M	1	0.13	~Dimethylbenz[a]anthracene, 7,12-	57-97-6	4.60E-02	c	8.40E-01	c	1.00E-02	c	
				4.00E-02	I					1	0.13	~Fluoranthene	206-44-0	2.40E+03	n	3.00E+04	n	8.00E+02	n	
				4.00E-02	I			V	1	0.13		~Fluorene	86-73-7	2.40E+03	n	3.00E+04	n	2.90E+02	n	
7.30E-01	E	1.10E-04	C						M	1	0.13	~Indeno[1,2,3-cd]pyrene	193-39-5	1.60E+01	c	2.90E+02	c	3.40E+00	c	
2.90E-02	P			7.00E-02	A			V	1	0.13	3.94E+02	~Methylnaphthalene, 1-	90-12-0	1.80E+03	c**s	7.30E+03	c**s	1.10E+02	c**	
				4.00E-03	I			V	1	0.13		~Methylnaphthalene, 2-	91-57-6	2.40E+02	n	3.00E+03	n	3.60E+01	n	
1.20E+00	C	3.40E-05	C	2.00E-02	I	3.00E-03	I	V	1	0.13		~Naphthalene	91-20-3	1.30E+02	n	5.90E+02	n	6.10E+00	n	
				3.00E-02	I			V	1	0.13		~Nitropyrene, 4-	57835-92-4	4.20E+01	c	1.80E+02	c	1.90E+00	c	
										1	0.1	~Pyrene	129-00-0	1.80E+03	n	2.30E+04	n	1.20E+02	n	
1.50E-01	I			2.00E-02	P				1	0.1		Potassium Perfluorobutane Sulfonate	29420-49-3	1.30E+03	n	1.60E+04	n	4.00E+02	n	
				9.00E-03	I				1	0.1		Prochloraz	67747-09-5	3.60E+02	c**	1.50E+03	c**	3.80E+01	c**	
				6.00E-03	H			V	1			Profluralin	26399-36-0	4.70E+02	n	7.00E+03	n	2.60E+01	n	
				1.50E-02	I				1	0.1		Prometon	1610-18-0	9.50E+02	n	1.20E+04	n	2.50E+02	n	
				4.00E-03	I				1	0.1		Prometryn	7287-19-6	2.50E+02	n	3.30E+03	n	6.00E+01	n	

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Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
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 _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
				1.30E-02	I				1	0.1		Propachlor	1918-16-7	8.20E+02	n	1.10E+04	n	2.50E+02	n	
				5.00E-03	I				1	0.1		Propanil	709-98-8	3.20E+02	n	4.10E+03	n	8.20E+01	n	
				2.00E-02	I				1	0.1		Propargite	2312-35-8	1.30E+03	n	1.60E+04	n	1.60E+02	n	
				2.00E-03	I		V		1		1.11E+05	Propargyl Alcohol	107-19-7	1.60E+02	n	2.30E+03	n	4.00E+01	n	
				2.00E-02	I				1	0.1		Propazine	139-40-2	1.30E+03	n	1.60E+04	n	3.40E+02	n	
				2.00E-02	I				1	0.1		Propham	122-42-9	1.30E+03	n	1.60E+04	n	3.50E+02	n	
				1.30E-02	I				1	0.1		Propiconazole	60207-90-1	8.20E+02	n	1.10E+04	n	2.10E+02	n	
						8.00E-03	I	V	1		3.26E+04	Propionaldehyde	123-38-6	7.50E+01	n	3.10E+02	n	1.70E+01	n	
				1.00E-01	X	1.00E+00	X	V	1		2.64E+02	Propyl benzene	103-65-1	3.80E+03	ns	2.40E+04	ns	6.60E+02	n	
						3.00E+00	C	V	1		3.49E+02	Propylene	115-07-1	2.20E+03	ns	9.30E+03	ns	6.30E+03	n	
				2.00E+01	P				1	0.1		Propylene Glycol	57-55-6	1.30E+06	nm	1.60E+07	nm	4.00E+05	n	
						2.70E-04	A		1	0.1		Propylene Glycol Dinitrate	6423-43-4	3.90E+05	nm	1.60E+06	nm	4.10E+02	n	
				7.00E-01	H	2.00E+00	I	V	1		1.06E+05	Propylene Glycol Monomethyl Ether	107-98-2	4.10E+04	n	3.70E+05	nms	3.20E+03	n	
2.40E-01	I	3.70E-06	I			3.00E-02	I	V	1		7.77E+04	Propylene Oxide	75-56-9	2.10E+02	c**	9.70E+02	c**	2.70E+01	c**	
				7.50E-02	I				1	0.1		Propyzamide	23950-58-5	4.70E+03	n	6.20E+04	n	1.20E+03	n	
				1.00E-03	I			V	1		5.30E+05	Pyridine	110-86-1	7.80E+01	n	1.20E+03	n	2.00E+01	n	
				5.00E-04	I				1	0.1		Quinalphos	13593-03-8	3.20E+01	n	4.10E+02	n	5.10E+00	n	
3.00E+00	I								1	0.1		Quinoline	91-22-5	1.80E+01	c	7.70E+01	c	2.40E+00	c	
				9.00E-03	I				1	0.1		Quizalofop-ethyl	76578-14-8	5.70E+02	n	7.40E+03	n	1.20E+02	n	
						3.00E-02	A		1			Refractory Ceramic Fibers	NA	4.30E+07	nm	1.80E+08	nm			
				3.00E-02	I				1	0.1		Resmethrin	10453-86-8	1.90E+03	n	2.50E+04	n	6.70E+01	n	
				5.00E-02	H			V	1			Ronnel	299-84-3	3.90E+03	n	5.80E+04	n	4.10E+02	n	
2.20E-01	C	6.30E-05	C						1	0.1		Rotenone	83-79-4	2.50E+02	n	3.30E+03	n	6.10E+01	n	
				5.00E-03	I				1	0.1		Safrole	94-59-7	5.50E+01	c	1.00E+03	c	9.60E+00	c	
				5.00E-03	I				1			Selenious Acid	7783-00-8	3.90E+02	n	5.80E+03	n	1.00E+02	n	
						2.00E-02	C		1			Selenium	7782-49-2	3.90E+02	n	5.80E+03	n	1.00E+02	n	5.0E+01
				5.00E-03	C	2.00E-02	C		1			Selenium Sulfide	7446-34-6	3.90E+02	n	5.80E+03	n	1.00E+02	n	
				9.00E-02	I				1	0.1		Sethoxydim	74051-80-2	5.70E+03	n	7.40E+04	n	1.00E+03	n	
						3.00E-03	C		1			Silica (crystalline, respirable)	7631-86-9	4.30E+06	nm	1.80E+07	nm			
1.20E-01	H			5.00E-03	I				0.04			Silver	7440-22-4	3.90E+02	n	5.80E+03	n	9.40E+01	n	
				5.00E-03	I				1	0.1		Simazine	122-34-9	3.20E+02	n	1.90E+03	c**	6.10E+01	c**	4.0E+00
				1.30E-02	I				1	0.1		Sodium Acifluorfen	62476-59-9	8.20E+02	n	1.10E+04	n	2.60E+02	n	
5.00E-01	C	1.50E-01	C			2.00E-02	C	2.00E-04	C	M	0.025	Sodium Azide	26628-22-8	3.10E+02	n	4.70E+03	n	8.00E+01	n	
				2.00E-02	C				1			Sodium Dichromate	10588-01-9	3.00E+01	c*	6.20E+02	c*	4.10E+00	c*	
2.70E-01	H			3.00E-02	I				1	0.1		Sodium Diethyldithiocarbamate	148-18-5	2.00E+02	c**	8.50E+02	c*	2.90E+01	c*	
				5.00E-02	A	1.30E-02	C		1			Sodium Fluoride	7681-49-4	3.90E+03	n	5.80E+04	n	1.00E+03	n	
				2.00E-05	I				1	0.1		Sodium Fluoroacetate	62-74-8	1.30E+00	n	1.60E+01	n	4.00E-01	n	
				1.00E-03	H				1			Sodium Metavanadate	13718-26-8	7.80E+01	n	1.20E+03	n	2.00E+01	n	
				8.00E-04	P				1			Sodium Tungstate	13472-45-2	6.30E+01	n	9.30E+02	n	1.60E+01	n	
				8.00E-04	P				1			Sodium Tungstate Dihydrate	10213-10-2	6.30E+01	n	9.30E+02	n	1.60E+01	n	
2.40E-02	H			3.00E-02	I				1	0.1		Stirofos (Tetrachlorovinphos)	961-11-5	1.90E+03	n	9.60E+03	c**	2.80E+02	c**	
5.00E-01	C	1.50E-01	C			2.00E-02	C	2.00E-04	C	M	0.025	Strontium Chromate	7789-06-2	3.00E+01	c*	6.20E+02	c*	4.10E+00	c*	
				6.00E-01	I				1			Strontium, Stable	7440-24-6	4.70E+04	n	7.00E+05	nm	1.20E+04	n	
				3.00E-04	I				1	0.1		Strychnine	57-24-9	1.90E+01	n	2.50E+02	n	5.90E+00	n	
				2.00E-01	I	1.00E+00	I	V	1		8.67E+02	Styrene	100-42-5	6.00E+03	ns	3.50E+04	ns	1.20E+03	n	1.0E+02
				3.00E-03	P				1	0.1		Styrene-Acrylonitrile (SAN) Trimer	NA	1.90E+02	n	2.50E+03	n	4.80E+01	n	
				1.00E-03	P	2.00E-03	X		1	0.1		Sulfolane	126-33-0	6.30E+01	n	8.20E+02	n	2.00E+01	n	
				8.00E-04	P				1	0.1		Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9	5.10E+01	n	6.60E+02	n	1.10E+01	n	
						1.00E-03	C	V	1			Sulfur Trioxide	7446-11-9	1.40E+06	nm	6.00E+06	nm	2.10E+00	n	
2.50E-02	I	7.10E-06	I			1.00E-03	C		1			Sulfuric Acid	7664-93-9	1.40E+06	nm	6.00E+06	nm			
				5.00E-02	H				1	0.1		Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl ester	140-57-8	2.20E+03	c**	9.20E+03	c**	1.30E+02	c**	
				3.00E-02	H				1	0.1		TCMTB	21564-17-0	1.90E+03	n	2.50E+04	n	4.80E+02	n	
				7.00E-02	I				1	0.1		Tebuthiuron	34014-18-1	4.40E+03	n	5.70E+04	n	1.40E+03	n	
				2.00E-02	H				1	0.1		Temephos	3383-96-8	1.30E+03	n	1.60E+04	n	4.00E+02	n	

Regional Removal Management Level (RML) Summary Table (TR=1E-04, HQ=1) May 2016

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Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
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 _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
				1.30E-02	I				1	0.1		Terbacil	5902-51-2	8.20E+02	n	1.10E+04	n	2.50E+02	n	
				2.50E-05	H			V	1		3.09E+01	Terbufos	13071-79-9	2.00E+00	n	2.90E+01	n	2.40E-01	n	
				1.00E-03	I				1	0.1		Terbutryn	886-50-0	6.30E+01	n	8.20E+02	n	1.30E+01	n	
				1.00E-04	I				1	0.1		Tetrabromodiphenyl ether, 2,2',4,4'-(BDE-47)	5436-43-1	6.30E+00	n	8.20E+01	n	2.00E+00	n	
				3.00E-04	I			V	1			Tetrachlorobenzene, 1,2,4,5-	95-94-3	2.30E+01	n	3.50E+02	n	1.70E+00	n	
2.60E-02	I	7.40E-06	I	3.00E-02	I			V	1		6.80E+02	Tetrachloroethane, 1,1,1,2-	630-20-6	2.00E+02	c*	8.80E+02	c*s	5.70E+01	c**	
2.00E-01	I	5.80E-05	C	2.00E-02	I			V	1		1.90E+03	Tetrachloroethane, 1,1,2,2-	79-34-5	6.00E+01	c*	2.70E+02	c*	7.60E+00	c*	
2.10E-03	I	2.60E-07	I	6.00E-03	I	4.00E-02	I	V	1		1.66E+02	Tetrachloroethylene	127-18-4	8.10E+01	n	3.90E+02	ns	4.10E+01	n	5.0E+00
				3.00E-02	I				1	0.1		Tetrachlorophenol, 2,3,4,6-	58-90-2	1.90E+03	n	2.50E+04	n	2.40E+02	n	
2.00E+01	H							V	1			Tetrachlorotoluene, p- alpha, alpha, alpha-	5216-25-1	3.50E+00	c	1.60E+01	c	1.30E-01	c	
				5.00E-04	I				1	0.1		Tetraethyl Dithiopyrophosphate	3689-24-5	3.20E+01	n	4.10E+02	n	7.10E+00	n	
				2.00E-03	P	8.00E+01	I	V	1		2.05E+03	Tetrafluoroethane, 1,1,1,2-	811-97-2	1.00E+05	nms	4.30E+05	nms	7.10E+05	n	
									1	0.0007		Tetryl (Trinitrophenylmethylnitramine)	479-45-8	1.60E+02	n	2.30E+03	n	3.90E+01	n	
				2.00E-05	S				1			Thallic Oxide	1314-32-5	1.60E+00	n	2.30E+01	n	4.00E-01	n	
				1.00E-05	X				1			Thallium (I) Nitrate	10102-45-1	7.80E-01	n	1.20E+01	n	2.00E-01	n	
				1.00E-05	X				1			Thallium (Soluble Salts)	7440-28-0	7.80E-01	n	1.20E+01	n	2.00E-01	n	2.0E+00
				1.00E-05	X			V	1			Thallium Acetate	563-68-8	7.80E-01	n	1.20E+01	n	2.00E-01	n	
				2.00E-05	X			V	1			Thallium Carbonate	6533-73-9	1.60E+00	n	2.30E+01	n	4.00E-01	n	
				1.00E-05	X				1			Thallium Chloride	7791-12-0	7.80E-01	n	1.20E+01	n	2.00E-01	n	
				1.00E-05	S				1			Thallium Selenite	12039-52-0	7.80E-01	n	1.20E+01	n	2.00E-01	n	
				2.00E-05	X				1			Thallium Sulfate	7446-18-6	1.60E+00	n	2.30E+01	n	4.00E-01	n	
				1.30E-02	I				1	0.1		Thiensusulfuron-methyl	79277-27-3	8.20E+02	n	1.10E+04	n	2.60E+02	n	
				1.00E-02	I				1	0.1		Thiobencarb	28249-77-6	6.30E+02	n	8.20E+03	n	1.60E+02	n	
				7.00E-02	X				1	0.0075		Thiodiglycol	111-48-8	5.40E+03	n	7.90E+04	n	1.40E+03	n	
				3.00E-04	H				1	0.1		Thiofanox	39196-18-4	1.90E+01	n	2.50E+02	n	5.30E+00	n	
				8.00E-02	I				1	0.1		Thiophanate, Methyl	23564-05-8	5.10E+03	n	6.60E+04	n	1.60E+03	n	
				5.00E-03	I				1	0.1		Thiram	137-26-8	3.20E+02	n	4.10E+03	n	9.80E+01	n	
				6.00E-01	H				1			Tin	7440-31-5	4.70E+04	n	7.00E+05	nm	1.20E+04	n	
				1.00E-04	A	V			1			Titanium Tetrachloride	7550-45-0	1.40E+05	nm	6.00E+05	nm	2.10E-01	n	
				8.00E-02	I	5.00E+00	I	V	1		8.18E+02	Toluene	108-88-3	4.90E+03	ns	4.70E+04	ns	1.10E+03	n	1.0E+03
				1.10E-05	C	8.00E-06	C	V	1			Toluene-2,4-diisocyanate	584-84-9	6.40E+00	n	2.70E+01	n	1.70E-02	n	
1.80E-01	X			2.00E-04	X				1	0.1		Toluene-2,5-diamine	95-70-5	1.30E+01	n	1.60E+02	n	4.00E+00	n	
				1.10E-05	C	8.00E-06	C	V	1		1.71E+03	Toluene-2,6-diisocyanate	91-08-7	5.30E+00	n	2.20E+01	n	1.70E-02	n	
1.60E-02	P	5.10E-05	C						1	0.1		Toluidine, o- (Methylaniline, 2-)	95-53-4	3.40E+03	c	1.40E+04	c	4.70E+02	c	
3.00E-02	P			4.00E-03	X				1	0.1		Toluidine, p-	106-49-0	2.50E+02	n	3.30E+03	n	7.70E+01	n	
				3.00E+00	P			V	1		3.42E-01	Total Petroleum Hydrocarbons (Aliphatic High)	NA	2.30E+05	nms	3.50E+06	nms	6.00E+04	n	
				6.00E-01	P	V			1		1.41E+02	Total Petroleum Hydrocarbons (Aliphatic Low)	NA	5.20E+02	ns	2.20E+03	ns	1.30E+03	n	
				1.00E-02	X	1.00E-01	P	V	1		6.86E+00	Total Petroleum Hydrocarbons (Aliphatic Medium)	NA	9.60E+01	ns	4.40E+02	ns	1.00E+02	n	
				4.00E-02	P				1	0.1		Total Petroleum Hydrocarbons (Aromatic High)	NA	2.50E+03	n	3.30E+04	n	8.00E+02	n	
				4.00E-03	P	3.00E-02	P	V	1		1.82E+03	Total Petroleum Hydrocarbons (Aromatic Low)	NA	8.20E+01	n	4.20E+02	n	3.30E+01	n	
1.10E+00	I	3.20E-04	I	4.00E-03	P	3.00E-03	P	V	1			Total Petroleum Hydrocarbons (Aromatic Medium)	NA	1.10E+02	n	6.00E+02	n	5.50E+00	n	
				7.50E-03	I				1	0.1		Toxaphene	8001-35-2	4.90E+01	c	2.10E+02	c	7.10E+00	c	3.0E+00
				3.00E-04	A			V	1			Tralometrin	66841-25-6	4.70E+02	n	6.20E+03	n	1.50E+02	n	
				8.00E+01	X				1	0.1		Tri-n-butyltin	688-73-3	2.30E+01	n	3.50E+02	n	3.70E+00	n	
				3.00E-02	I				1	0.1		Triacetin	102-76-1	5.10E+06	nm	6.60E+07	nm	1.60E+06	n	
									1	0.1		Triadimefon	43121-43-3	1.90E+03	n	2.50E+04	n	5.50E+02	n	
				1.30E-02	I			V	1			Triallate	2303-17-5	1.00E+03	n	1.50E+04	n	1.20E+02	n	
				1.00E-02	I				1	0.1		Triasulfuron	82097-50-5	6.30E+02	n	8.20E+03	n	2.00E+02	n	
				8.00E-03	I				1	0.1		Tribenuron-methyl	101200-48-0	5.10E+02	n	6.60E+03	n	1.60E+02	n	
9.00E-03	P			5.00E-03	I			V	1			Tribromobenzene, 1,2,4-	615-54-3	3.90E+02	n	5.80E+03	n	4.50E+01	n	
				1.00E-02	P				1	0.1		Tributyl Phosphate	126-73-8	6.30E+02	n	8.20E+03	n	1.20E+02	n	
				3.00E-04	P				1	0.1		Tributyltin Compounds	NA	1.90E+01	n	2.50E+02	n	6.00E+00	n	
				3.00E-04	I				1	0.1		Tributyltin Oxide	56-35-9	1.90E+01	n	2.50E+02	n	5.70E+00	n	
				3.00E+01	I	3.00E+01	H	V	1		9.10E+02	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1	4.00E+04	ns	1.70E+05	nms	5.50E+04	n	

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Toxicity and Chemical-specific Information											Contaminant		Removal Levels							
SFO (mg/kg-day) ⁻¹	k _e y	IUR (ug/m ³) ⁻¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _i (mg/m ³)	k _e y	muta- gen	GIABS	ABS	C _{sat} (mg/kg)	Analyte	CAS No.	Resident Soil (mg/kg)	key	Industrial Soil (mg/kg)	key	Tapwater (ug/L)	key	MCL (ug/L)
7.00E-02	I			2.00E-02	I					1	0.1	Trichloroacetic Acid	76-03-9	7.80E+02	c**	3.30E+03	c**	1.10E+02	c**	6.0E+01
2.90E-02	H									1	0.1	Trichloroaniline HCl, 2,4,6-	33663-50-2	1.90E+03	c	7.90E+03	c	2.70E+02	c	
7.00E-03	X			3.00E-05	X					1	0.1	Trichloroaniline, 2,4,6-	634-93-5	1.90E+00	n	2.50E+01	n	4.00E-01	n	
				8.00E-04	X			V		1		Trichlorobenzene, 1,2,3-	87-61-6	6.30E+01	n	9.30E+02	n	7.00E+00	n	
2.90E-02	P			1.00E-02	I	2.00E-03	P	V		1	4.04E+02	Trichlorobenzene, 1,2,4-	120-82-1	5.80E+01	n	2.60E+02	n	4.00E+00	n	7.0E+01
				2.00E+00	I	5.00E+00	I	V		1	6.40E+02	Trichloroethane, 1,1,1-	71-55-6	8.10E+03	ns	3.60E+04	ns	8.00E+03	n	2.0E+02
5.70E-02	I	1.60E-05	I	4.00E-03	I	2.00E-04	X	V		1	2.16E+03	Trichloroethane, 1,1,2-	79-00-5	1.50E+00	n	6.30E+00	n	4.10E-01	n	5.0E+00
4.60E-02	I	4.10E-06	I	5.00E-04	I	2.00E-03	I	V	M	1	6.92E+02	Trichloroethylene	79-01-6	4.10E+00	n	1.90E+01	n	2.80E+00	n	5.0E+00
				3.00E-01	I			V		1	1.23E+03	Trichlorofluoromethane	75-69-4	2.30E+04	ns	3.50E+05	nms	5.20E+03	n	
				1.00E-01	I					1	0.1	Trichlorophenol, 2,4,5-	95-95-4	6.30E+03	n	8.20E+04	n	1.20E+03	n	
1.10E-02	I	3.10E-06	I	1.00E-03	P					1	0.1	Trichlorophenol, 2,4,6-	88-06-2	6.30E+01	n	8.20E+02	n	1.20E+01	n	
				1.00E-02	I					1	0.1	Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5	6.30E+02	n	8.20E+03	n	1.60E+02	n	
				8.00E-03	I					1	0.1	Trichlorophenoxypropionic acid, -2,4,5	93-72-1	5.10E+02	n	6.60E+03	n	1.10E+02	n	5.0E+01
3.00E+01	I			5.00E-03	I			V		1	1.28E+03	Trichloropropane, 1,1,2-	598-77-6	3.90E+02	n	5.80E+03	ns	8.80E+01	n	
				4.00E-03	I	3.00E-04	I	V	M	1	1.40E+03	Trichloropropane, 1,2,3-	96-18-4	5.10E-01	c**	1.10E+01	c**	7.50E-02	c**	
				3.00E-03	X	3.00E-04	P	V		1	3.11E+02	Trichloropropene, 1,2,3-	96-19-5	7.30E-01	n	3.10E+00	n	6.20E-01	n	
										1	0.1	Tricresyl Phosphate (TCP)	1330-78-5	1.30E+03	n	1.60E+04	n	1.60E+02	n	
										1	0.1	Triphane	58138-08-2	1.90E+02	n	2.50E+03	n	1.80E+01	n	
						7.00E-03	I	V		1	2.79E+04	Triethylamine	121-44-8	1.20E+02	n	4.80E+02	n	1.50E+01	n	
				2.00E+00	P					1	0.1	Triethylene Glycol	112-27-6	1.30E+05	nm	1.60E+06	nm	4.00E+04	n	
7.70E-03	I			7.50E-03	I			V		1	4.81E+03	Trifluoroethane, 1,1,1-	420-46-2	1.50E+04	ns	6.20E+04	ns	4.20E+04	n	
										1		Trifuralin	1582-09-8	5.90E+02	n	8.80E+03	n	4.00E+01	n	
2.00E-02	P			1.00E-02	P					1	0.1	Trimethyl Phosphate	512-56-1	6.30E+02	n	8.20E+03	n	2.00E+02	n	
						5.00E-03	P	V		1	2.93E+02	Trimethylbenzene, 1,2,3-	526-73-8	4.90E+01	n	2.10E+02	n	1.00E+01	n	
						7.00E-03	P	V		1	2.19E+02	Trimethylbenzene, 1,2,4-	95-63-6	5.80E+01	n	2.40E+02	ns	1.50E+01	n	
				1.00E-02	X			V		1	1.82E+02	Trimethylbenzene, 1,3,5-	108-67-8	7.80E+02	ns	1.20E+04	ns	1.20E+02	n	
				1.00E-02	X			V		1	2.96E+01	Trimethylpentene, 2,4,4-	25167-70-8	7.80E+02	ns	1.20E+04	ns	6.50E+01	n	
				3.00E-02	I					1	0.019	Trinitrobenzene, 1,3,5-	99-35-4	2.20E+03	n	3.20E+04	n	5.90E+02	n	
3.00E-02	I			5.00E-04	I					1	0.032	Trinitrotoluene, 2,4,6-	118-96-7	3.60E+01	n	5.10E+02	n	9.80E+00	n	
				2.00E-02	P					1	0.1	Triphenylphosphine Oxide	791-28-6	1.30E+03	n	1.60E+04	n	3.60E+02	n	
				2.00E-02	A					1	0.1	Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8	1.30E+03	n	1.60E+04	n	3.60E+02	n	
2.30E+00	C	6.60E-04	C	1.00E-02	X					1	0.1	Tris(1-chloro-2-propyl)phosphate	13674-84-5	6.30E+02	n	8.20E+03	n	1.90E+02	n	
2.00E-02	P			7.00E-03	P			V		1	4.67E+02	Tris(2,3-dibromopropyl)phosphate	126-72-7	2.80E+01	c	1.30E+02	c	6.80E-01	c	
										1	0.1	Tris(2-chloroethyl)phosphate	115-96-8	4.40E+02	n	5.70E+03	n	1.40E+02	n	
3.20E-03	P			1.00E-01	P					1	0.1	Tris(2-ethylhexyl)phosphate	78-42-2	6.30E+03	n	7.20E+04	c**	2.00E+03	n	
				8.00E-04	P					1		Tungsten	7440-33-7	6.30E+01	n	9.30E+02	n	1.60E+01	n	
				3.00E-03	I	4.00E-05	A			1		Uranium (Soluble Salts)	NA	2.30E+02	n	3.50E+03	n	6.00E+01	n	3.0E+01
1.00E+00	C	2.90E-04	C	8.30E-03	P				M	1	0.1	Urethane	51-79-6	1.20E+01	c	2.30E+02	c	2.50E+00	c	
				9.00E-03	I	7.00E-06	P			0.026		Vanadium Pentoxide	1314-62-1	6.60E+02	n	8.40E+03	n	1.50E+02	n	
				5.00E-03	S	1.00E-04	A			0.026		Vanadium and Compounds	7440-62-2	3.90E+02	n	5.80E+03	n	8.60E+01	n	
				1.00E-03	I			V		1		Vernolate	1929-77-7	7.80E+01	n	1.20E+03	n	1.10E+01	n	
				2.50E-02	I					1	0.1	Vinclozolin	50471-44-8	1.60E+03	n	2.10E+04	n	4.40E+02	n	
				1.00E+00	H	2.00E-01	I	V		1	2.75E+03	Vinyl Acetate	108-05-4	9.10E+02	n	3.80E+03	ns	4.10E+02	n	
3.20E-05	H					3.00E-03	I	V		1	2.47E+03	Vinyl Bromide	593-60-2	4.30E+00	n	1.80E+01	n	6.30E+00	n	
7.20E-01	I	4.40E-06	I	3.00E-03	I	1.00E-01	I	V	M	1	3.92E+03	Vinyl Chloride	75-01-4	5.90E+00	c*	1.70E+02	c**	1.90E+00	c*	2.0E+00
				3.00E-04	I					1	0.1	Warfarin	81-81-2	1.90E+01	n	2.50E+02	n	5.60E+00	n	
				2.00E-01	S	1.00E-01	S	V		1	3.90E+02	Xylene, p-	106-42-3	5.60E+02	ns	2.40E+03	ns	1.90E+02	n	
				2.00E-01	S	1.00E-01	S	V		1	3.88E+02	Xylene, m-	108-38-3	5.50E+02	ns	2.40E+03	ns	1.90E+02	n	
				2.00E-01	S	1.00E-01	S	V		1	4.34E+02	Xylene, o-	95-47-6	6.50E+02	ns	2.80E+03	ns	1.90E+02	n	
				2.00E-01	I	1.00E-01	I	V		1	2.60E+02	Xylenes	1330-20-7	5.80E+02	ns	2.50E+03	ns	1.90E+02	n	1.0E+04
				3.00E-04	I					1		Zinc Phosphide	1314-84-7	2.30E+01	n	3.50E+02	n	6.00E+00	n	
				3.00E-01	I					1		Zinc and Compounds	7440-66-6	2.30E+04	n	3.50E+05	nm	6.00E+03	n	
				5.00E-02	I					1	0.1	Zineb	12122-67-7	3.20E+03	n	4.10E+04	n	9.90E+02	n	
				8.00E-05	X					1		Zirconium	7440-67-7	6.30E+00	n	9.30E+01	n	1.60E+00	n	

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k e IUR (ug/m ³) ⁻¹	k e RfD _o (mg/kg- day)	k e RfC (mg/m ³) ⁻¹	k e 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-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)				
8.7E-03	I	4.0E-03	I	I	V	1.1E+05	1.4E+09	8.7E+03	1	0.1	Acephate	30560-19-1	8.0E+03	2.8E+04	1.1E+03	6.2E+03	3.1E+02	1.3E+03	8.2E+01	2.5E+02			
		2.2E-06	I			1.4E+09	1.4E+09	1			75-07-0				1.1E+03				8.2E+01				
		2.0E-02	I			1.4E+09	1.4E+09	1			34256-82-1				1.1E+03	1.6E+03	6.6E+03	8.2E+01	1.3E+03				
		9.0E-01	I	3.1E+01	A	V	1.1E+05	1.4E+09	1.4E+04	1	0.1	Acetone	67-64-1				7.0E+04		4.4E+05	6.1E+04			
		2.0E-03	X			1.4E+09	1.4E+09	1			75-86-5							2.8E+06	2.8E+06				
		6.0E-02	I	V		1.3E+05	1.4E+09	1.3E+04	1		75-05-8							8.1E+02	8.1E+02				
3.8E+00	C	1.3E-03	C			1.0E-01	I	V	2.5E+03	1.4E+09	6.0E+04	1			98-86-2				7.8E+03	7.8E+03			
		5.0E-04	I	2.0E-05	I	V	2.3E+04	1.4E+09	6.9E+03	1	0.1	Acetylaminofluorene, 2-Acrolein	53-96-3	1.8E+01	6.5E+01	2.9E+05	1.4E+01						
		2.0E-03	I	6.0E-03	I	M	1.4E+09	1			107-02-8					3.9E+01		1.4E-01	1.4E-01				
5.0E-01	I	1.0E-04	I			2.0E-03	I	6.0E-03	I	M	1.4E+09	1			79-06-1	1.6E+02	6.6E+02	8.5E+06	1.3E+02				
		5.0E-01	I	1.0E-03	I	V	1.1E+05	1.4E+09	9.5E+04	1	0.1	Acrylamide	79-10-7	3.1E+01	1.2E+02	1.4E+06	2.4E+01						
5.4E-01	I	6.8E-05	I	4.0E-02	A	2.0E-03	I	V	1.1E+04	1.4E+09	7.7E+03	1			107-13-1	3.1E+03		1.6E+01	1.6E+01				
		6.0E-03	P			1.4E+09	1				111-69-3							8.5E+06	8.5E+06				
5.6E-02	C	1.0E-02	I			1.4E+09	1				15972-60-8	1.2E+03	4.4E+03		9.7E+02	7.8E+02	3.3E+03		6.3E+02				
		1.0E-03	I			1.4E+09	1				116-06-3					7.8E+01	3.3E+02		6.3E+01				
		1.0E-03	I			1.4E+09	1				1646-88-4					7.8E+01	3.3E+02		6.3E+01				
1.7E+01	I	4.9E-03	I	3.0E-05	I	V	1.4E+09	1.7E+06	1		309-00-2	4.1E+00		9.8E+01	3.9E+00	2.3E+00			2.3E+00				
		5.0E-03	I	1.0E-04	X	V	1.1E+05	1.4E+09	3.4E+04	1		107-18-6					3.9E+02		3.6E+00	3.5E+00			
2.1E-02	C	6.0E-06	C			1.0E+00	P	5.0E-03	P		1.4E+09	1			107-05-1	3.3E+03		7.4E+01	7.2E+01				
		1.0E+00	P	5.0E-03	P		1.4E+09	1			7429-90-5					7.8E+04		7.1E+06	7.7E+04				
		4.0E-04	I			1.4E+09	1				20859-73-8					3.1E+01			3.1E+01				
2.1E+01	C	6.0E-03	C			9.0E-03	I				834-12-8					7.0E+02	3.0E+03		5.7E+02				
		2.5E-03	I			1.4E+09	1				92-67-1	3.3E+00	1.2E+01	6.4E+04	2.6E+00								
		8.0E-02	P			1.4E+09	1				591-27-5					6.3E+03	2.6E+04		5.1E+03				
		2.0E-02	P			1.4E+09	1				123-30-8					1.6E+03	6.6E+03		1.3E+03				
		2.5E-03	I			1.4E+09	1				33089-61-1					2.0E+02	8.2E+02		1.6E+02				
		1.0E-01	I	V		1.4E+09	1				7664-41-7					1.6E+04		8.2E+01	1.6E+04				
		2.0E-01	I	3.0E-03	X	V	1.4E+04	1.4E+09	2.6E+04	1		7773-06-0							8.2E+01				
		1.4E+09	1			1.4E+09	1				15-85-4								8.2E+01				
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I				1.4E+09	1			62-53-3	1.2E+04	4.3E+04	2.4E+08	9.5E+03				
4.0E-02	P	2.0E-03	X			1.4E+09	1				1.4E+09	1			44-65-1	1.7E+03	6.2E+03	1.4E+03	5.5E+02				
		4.0E-04	I			1.4E+09	0.15				1440-36-0					1.6E+02	6.6E+02	1.4E+06	4.4E+02				
		5.0E-04	H			1.4E+09	0.15				1314-60-9					3.1E+01			3.1E+01				
		4.0E-04	H			1.4E+09	0.15				1332-81-6					3.1E+01			3.1E+01				
		2.0E-04	I			1.4E+09	0.15				1309-64-4							2.8E+05	2.8E+05				
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C				1.4E+09	1	0.03		7440-38-2	7.7E+01	5.5E+02	8.9E+04	6.8E+01				
		3.5E-06	C	5.0E-05	I		1.4E+09	1			7784-42-1					3.9E+01	3.3E+02	2.1E+04	3.5E+01				
		5.0E-02	I			1.4E+09	1				3337-71-1					2.7E-01		7.1E+04	2.7E-01				
2.3E-01	C	3.5E-02	I			1.4E+09	1				1912-24-9	3.0E+02	1.1E+03		2.4E+02	3.9E+03	1.6E+04		3.2E+03				
8.8E-01	C	2.5E-04	C			1.4E+09	1				492-80-8	7.9E+01	2.8E+02	1.5E+06	6.2E+01	2.7E+03	1.2E+04		2.2E+03				
		4.0E-04	I			1.4E+09	1				65195-55-3					3.1E+01	1.3E+02		2.5E+01				
1.1E-01	I	3.1E-05	I	3.0E-03	A	1.0E-02	A				1.4E+09	1			86-50-0	6.3E+02		4.7E+03	5.6E+02				
		1.0E+00	P	7.0E-06	P		1.4E+09	1			103-33-3					2.3E+02	9.9E+02	1.4E+07	1.9E+02				
		2.0E-01	I	5.0E-04	H		1.4E+09	0.07			7440-39-3					1.6E+04		7.1E+05	1.5E+04				
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M			1.4E+09	0.025			10294-40-3	3.1E+01		9.2E+02	3.0E+01				
		3.0E-01	I	V		1.4E+09	3.1E+05	1			1861-40-1					1.6E+03		2.8E+05	1.6E+03				
		5.0E-02	I			1.4E+09	1				17804-35-2					2.3E+04	3.3E+05	9.9E+03	8.6E+03				
		2.0E-01	I			1.4E+09	1				83055-99-6					7.8E+04	3.3E+05	9.9E+03	8.6E+03				
		3.0E-02	I			1.4E+09	1				25057-89-0					3.9E+03	1.6E+04		3.2E+03				
4.0E-03	P	1.0E-01	I	V		1.2E+03	1.4E+09	2.3E+04	1		100-52-7	1.7E+04			1.7E+04	1.6E+04	6.6E+04		1.3E+04				
5.5E-02	I	7.8E-06	I	4.0E-03	I	3.0E-02	I	V	1.8E+03	1.4E+09	3.5E+03	1			3.1E+02			1.1E+02	8.2E+01				
1.0E-01	X	3.0E-04	X			1.4E+09	1				6369-59-1	7.0E+02	2.5E+03	1.3E+02	5.4E+02	2.3E+01	9.9E+01		1.9E+01				
		1.0E-03	P			1.3E+03	1.4E+09	1.9E+04	1		108-98-5					7.8E+01			7.8E+01				
2.3E+02	I	6.7E-02	I	3.0E-03	I		1.4E+09	1			92-87-5	6.7E-02	2.6E-01	2.1E+03	5.3E-02	2.3E+02	9.9E+02		1.9E+02				
		4.0E+00	I			1.4E+09	1				65-85-0					3.1E+05	1.3E+06		2.5E+05				
1.3E+01	I			V		3.2E+02	1.4E+09	6.8E+04	1		98-07-7	5.3E+00			5.3E+00	7.8E+03	3.3E+04		6.3E+03				
		1.0E-01	P			1.4E+09	1				100-51-6								6.3E+03				

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRVT; A = ATSDR; C = Cal EPA; X = APPENDIX PPRVT SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1				
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg- day)	k e y	RfC _o (mg/m ³) ⁻¹	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-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Child THQ=1 (mg/kg)	Child THQ=1 (mg/kg)	Child THQ=1 (mg/kg)	Child THI=1 (mg/kg)	
1.7E-01	I	4.9E-05	C	2.0E-03	P	1.0E-03	P	V	1.5E+03	1.4E+09	2.6E+04	1		Benzyl Chloride	100-44-7	4.1E+02		1.5E+02	1.1E+02	1.6E+02		2.7E+01	2.3E+01	
		2.4E-03	I	2.0E-03	I	2.0E-05	I			1.4E+09		0.007		Beryllium and compounds	7440-41-7			1.6E+05	1.6E+05	1.6E+02		2.8E+04	1.6E+02	
				9.0E-03	P					1.4E+09			0.1	Bifenox	42576-02-3					7.0E+02	3.0E+03		5.7E+02	
				1.5E-02	I					1.4E+09			0.1	Biphenthrin	82657-04-3					1.2E+03	4.9E+03		9.5E+02	
8.0E-03	I			4.0E-04	X	V				1.4E+09	1.1E+05	1		Biphenyl, 1,1'-	92-52-4	8.7E+03			8.7E+03	3.9E+04		4.8E+01	4.7E+01	
				4.0E-02	I	V			1.0E+03	1.4E+09	3.5E+04	1		Bis(2-chloro-1-methylethyl) ether	108-60-1					3.1E+03			3.1E+03	
				3.0E-03	P					1.4E+09			0.1	Bis(2-chloroethoxy)methane	111-91-1					2.3E+02	9.9E+02		1.9E+02	
1.1E+00	I	3.3E-04	I					V	5.1E+03	1.4E+09	4.3E+04	1		Bis(2-chloroethoxy)ether	111-44-4	6.3E+01		3.6E+01	2.3E+01					
2.2E+02	I	6.2E-02	I					V	4.2E+03	1.4E+09	1.9E+03	1		Bis(chloromethyl)ether	542-88-1	3.2E-01		8.5E-03	8.3E-03					
				5.0E-02	I					1.4E+09			0.1	Bisphenol A	80-05-7					3.9E+03	1.6E+04		3.2E+03	
				2.0E-01	I	2.0E-02	H			1.4E+09			1	Boron And Borates Only	7440-42-8					1.6E+04		2.8E+07	1.6E+04	
				2.0E+00	P	2.0E-02	P	V		1.4E+09			1	Boron Trichloride	10294-34-5					1.6E+05		2.8E+07	1.6E+05	
				4.0E-02	C	1.3E-02	C	V		1.4E+09			1	Boron Trifluoride	7637-07-2					3.1E+03		1.8E+07	3.1E+03	
7.0E-01	I			4.0E-03	I					1.4E+09			1	Bromate	15541-45-4	9.9E+01			9.9E+01				3.1E+02	3.1E+02
2.0E+00	X	6.0E-04	X					V	2.4E+03	1.4E+09	5.9E+03	1		Bromo-2-chloroethane, 1-	107-04-0	3.5E+01		2.8E+00	2.6E+00					
				8.0E-03	I	6.0E-02	I	V	6.8E+02	1.4E+09	8.4E+03	1		Bromobenzene	108-86-1					6.3E+02		5.2E+02	2.9E+02	
				4.0E-02	X	V			4.0E+03	1.4E+09	3.6E+03	1		Bromochloromethane	74-97-5							1.5E+02	1.5E+02	
6.2E-02	I	3.7E-05	C	2.0E-02	I			V	9.3E+02	1.4E+09	4.0E+03	1		Bromodichloromethane	75-27-4	1.1E+03		3.0E+01	2.9E+01	1.6E+03			1.6E+03	
7.9E-03	I	1.1E-06	I	2.0E-02	I			V	9.2E+02	1.4E+09	9.7E+03	1		Bromoform	75-25-2	8.8E+03		2.5E+03	1.9E+03	1.6E+03			1.6E+03	
				1.4E-03	I	5.0E-03	I	V	3.6E+03	1.4E+09	1.4E+03	1		Bromomethane	74-83-9					1.1E+02		7.3E+00	6.8E+00	
				5.0E-03	H			V	1.4E+09	1.2E+05		1		Bromophos	2104-96-3					3.9E+02			3.9E+02	
				2.0E-02	I			V	1.4E+09			1	0.1	Bromoxynil	1689-84-5					1.6E+03	6.6E+03		1.3E+03	
				2.0E-02	I			V	1.4E+09	4.7E+05		1		Bromoxynil Octanoate	1689-99-2					1.6E+03			1.6E+03	
3.4E+00	C	3.0E-05	I			2.0E-03	I	V	6.7E+02	1.4E+09	8.7E+02	1		Butadiene, 1,3-	106-99-0	2.0E+01		8.1E+00	5.8E+00			1.8E+00	1.8E+00	
				1.0E-01	I			V	7.6E+03	1.4E+09	3.0E+04	1		Butanol, N-	71-36-3					7.8E+03			7.8E+03	
				2.0E+00	P	3.0E+01	P	V	2.1E+04	1.4E+09	2.9E+04	1		Butyl alcohol, sec-	78-92-2					1.6E+05		9.1E+05	1.3E+05	
				5.0E-02	I			V	1.4E+09	8.6E+04		1		Butylate	2008-41-5					3.9E+03			3.9E+03	
2.0E-04	C	5.7E-08	C							1.4E+09			0.1	Butylated hydroxyanisole	25013-16-5	3.5E+05	1.2E+06	6.7E+09	2.7E+05					
3.6E-03	P			3.0E-01	P					1.4E+09			0.1	Butylated hydroxytoluene	128-37-0	1.9E+04	6.9E+04		1.5E+04	2.3E+04	9.9E+04		1.9E+04	
				5.0E-02	P			V	1.1E+02	1.4E+09	8.1E+03	1		Butylbenzene, n-	104-51-8					3.9E+03			3.9E+03	
				1.0E-01	X	V			1.5E+02	1.4E+09	7.4E+03	1		Butylbenzene, sec-	135-98-8					7.8E+03			7.8E+03	
				1.0E-01	X	V			1.8E+02	1.4E+09	7.4E+03	1		Butylbenzene, tert-	98-06-6					7.8E+03			7.8E+03	
				2.0E-02	A					1.4E+09			0.1	Cacodylic Acid	15-60-5					1.6E+03	6.6E+03		1.3E+03	
1.8E-03	I	1.0E-03	I	1.0E-05	A					1.4E+09		0.025	0.001	Cadmium (Diet)	1440-43-9			2.1E+05	2.1E+05	7.8E+01	8.2E+02	1.4E+04	7.1E+01	
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M	1.4E+09		0.025			Cadmium (Water)	1440-43-9	3.1E+01		9.2E+02	3.0E+01	1.6E+03		2.8E+05	1.6E+03	
				5.0E-01	I	2.2E-03	C			1.4E+09			0.1	Calcium Chromate	3765-19-0					3.9E+04	1.6E+05	3.1E+06	3.1E+04	
				1.5E-01	C	4.3E-05	C	2.0E-03	I	1.4E+09			0.1	Captafol	2425-06-1	4.6E+02	1.6E+03	8.9E+06	3.6E+02	1.6E+02	6.6E+02		1.3E+02	
2.3E-03	C	6.6E-07	C	1.3E-01	I					1.4E+09			0.1	Captan	133-06-2	3.0E+04	1.1E+05	5.8E+08	2.4E+04	1.0E+04	4.3E+04		8.2E+03	
				1.0E-01	I					1.4E+09			0.1	Carbaryl	63-25-2					7.8E+03	3.3E+04		6.3E+03	
				5.0E-03	I					1.4E+09			0.1	Carbofuran	1563-66-2					3.9E+02	1.6E+03		3.2E+02	
				1.0E-01	I	7.0E-01	I	V	7.4E+02	1.4E+09	1.2E+03	1		Carbon Disulfide	75-15-0					7.8E+03		8.5E+02	7.7E+02	
7.0E-02	I	6.0E-06	I	4.0E-03	I	1.0E-01	I	V	4.6E+02	1.4E+09	1.5E+03	1		Carbon Tetrachloride	56-23-5	9.9E+02		7.0E+01	6.5E+01	3.1E+02		1.6E+02	1.0E+02	
				1.0E-01	P	V			5.9E+03	1.4E+09	6.5E+02	1		Carbonyl Sulfide	463-58-1							6.7E+01	6.7E+01	
				1.0E-02	I					1.4E+09			0.1	Carbosulfan	55285-14-8					7.8E+02	3.3E+03		6.3E+02	
				1.0E-01	I					1.4E+09			0.1	Carboxin	5234-68-4					7.8E+03	3.3E+04		6.3E+03	
				9.0E-04	I					1.4E+09			1	Ceric oxide	1306-38-3							1.3E+06	1.3E+06	
				1.0E-01	I			V	1.4E+09	1.5E+05		1		Chloral Hydrate	302-17-0					7.8E+03			7.8E+03	
				1.5E-02	I					1.4E+09			0.1	Chloramben	133-90-4					1.2E+03	4.9E+03		9.5E+02	
4.0E-01	H									1.4E+09			0.1	Chloranil	118-75-2	1.7E+02	6.1E+02		1.3E+02					
3.5E-01	I	1.0E-04	I	5.0E-04	I	7.0E-04	I	V	1.4E+09	1.5E+06		0.04		Chlordane	12789-03-6	2.0E+02	1.8E+03	4.3E+03	1.7E+02	3.9E+01	4.1E+02	1.1E+03	3.5E+01	
1.0E+01	I	4.6E-03	C	3.0E-04	I					1.4E+09			0.1	Chlordecone (Kepone)	143-50-0	7.0E+00	2.5E+01	8.3E+04	5.4E+00	2.3E+01	9.9E+01		1.9E+01	
				7.0E-04	A					1.4E+09			0.1	Chlorfenvinphos	470-90-6					5.5E+01	2.3E+02		4.4E+01	
				2.0E-02	I					1.4E+09			0.1	Chlorimuron, Ethyl-	90982-32-4					1.6E+03	6.6E+03		1.3E+03	
				1.0E-01	I	1.5E-04	A	V	2.8E+03	1.4E+09	1.2E+03	1		Chlorine	7782-50-5					7.8E+03		1.8E-01	1.8E-01	
				3.0E-02	I	2.0E-04	I	V		1.4E+09			1	Chlorine Dioxide	10049-04-4					2.3E+03		2.8E+05	2.3E+03	
				3.0E-02	I					1.4E+09			1	Chlorite (Sodium Salt)	7758-19-2					2.3E+03			2.3E+03	
				5.0E+01	I																			

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1				
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg- day)	k e y	RfC _o (mg/m ³) ⁻¹	k e y	o v e r l o a d	muta- gen	C _{mt} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
4.6E-01	H									1.4E+09					Chloro-2-methylaniline HCl, 4-	3165-93-3	1.5E+02	5.4E+02		1.2E+02				
1.0E-01	P	7.7E-05	C	3.0E-03	X					1.4E+09					Chloro-2-methylaniline, 4-	95-69-2	7.0E+02	2.5E+03	5.0E+06	5.4E+02	2.3E+02	9.9E+02		1.9E+02
2.7E-01	X									1.2E+04	1.4E+09	1.6E+04	1		Chloroacetaldehyde, 2-	107-20-0	2.6E+02			2.6E+02				
										1.4E+09					Chloroacetic Acid	79-11-8								
						3.0E-05	I			1.4E+09					Chloroacetophenone, 2-	532-27-4						4.3E+04	4.3E+04	
2.0E-01	P			4.0E-03	I					1.4E+09					Chloroaniline, p-	106-47-8	3.5E+02	1.2E+03		2.7E+02	3.1E+02	1.3E+03		2.5E+02
				2.0E-02	I	5.0E-02	P	V		7.6E+02	1.4E+09	6.5E+03	1		Chlorobenzene	108-90-7					1.6E+03		3.4E+02	2.8E+02
1.1E-01	C	3.1E-05	C	2.0E-02	I					1.4E+09					Chlorobenzilate	510-15-6	6.3E+02	2.2E+03	1.2E+07	4.9E+02	1.6E+03	6.6E+03		1.3E+03
				3.0E-02	X					1.4E+09					Chlorobenzoic Acid, p-	74-11-3					2.3E+03	9.9E+03		1.9E+03
				3.0E-03	P	3.0E-01	P	V		2.9E+02	1.4E+09	6.8E+03	1		Chlorobenzoic acid, 4-	98-56-6					2.3E+02		2.1E+03	2.1E+02
				4.0E-02	P			V		7.3E+02	1.4E+09	1.8E+03	1		Chlorobutane, 1-	109-69-3					3.1E+03			3.1E+03
				5.0E+01	I	V		V		1.7E+03	1.4E+09	9.4E+02	1		Chlorodifluoromethane	75-45-6							4.9E+04	4.9E+04
				2.0E-02	P			V		1.1E+05	1.4E+09	7.8E+04	1		Chloroethanol, 2-	107-07-3					1.6E+03			1.6E+03
3.1E-02	C	2.3E-05	I	1.0E-02	I	9.8E-02	A	V		2.5E+03	1.4E+09	2.6E+03	1		Chloroform	67-66-3	2.2E+03		3.2E+01	3.2E+01	7.8E+02		2.7E+02	2.0E+02
				9.0E-02	I	V		V		1.3E+03	1.4E+09	1.2E+03	1		Chloromethane	74-87-3							1.1E+02	1.1E+02
2.4E+00	C	6.9E-04	C					V		9.3E+03	1.4E+09	5.3E+03	1		Chloromethyl Methyl Ether	107-30-2	2.9E+01		2.2E+00	2.0E+00				
3.0E-01	P			3.0E-03	P	1.0E-05	X			1.4E+09					Chloronitrobenzene, o-	88-73-3	2.3E+02	8.2E+02		1.8E+02	2.3E+02	9.9E+02	1.4E+04	1.9E+02
6.0E-02	P			7.0E-04	P	2.0E-03	P			1.4E+09					Chloronitrobenzene, p-	100-00-5	1.2E+03	4.1E+03		9.0E+02	5.5E+01	2.3E+02	2.8E+06	4.4E+01
				5.0E-03	I	V		V		2.7E+04	1.4E+09	1.4E+05	1		Chlorophenol, 2-	95-57-8					3.9E+02			3.9E+02
				4.0E-04	C	V		V		6.2E+02	1.4E+09	4.7E+03	1		Chloropicrin	76-06-2							2.0E+00	2.0E+00
3.1E-03	C	8.9E-07	C	1.5E-02	I			V		1.4E+09					Chloroethanol	1897-45-6	2.2E+04	8.0E+04	4.3E+08	1.8E+04	1.2E+03	4.9E+03		9.5E+02
				2.0E-02	I			V		9.1E+02	1.4E+09	8.1E+03	1		Chlorotoluene, o-	95-49-8					1.6E+03			1.6E+03
				2.0E-02	X	V		V		2.5E+02	1.4E+09	7.3E+03	1		Chlorotoluene, p-	106-43-4					1.6E+03			1.6E+03
2.4E+02	C	6.9E-02	C					V		1.4E+09					Chlorzotocin	54749-90-5	2.9E-01	1.0E+00	5.5E+03	2.3E-01				
				2.0E-01	I			V		1.4E+09					Chlorpropopham	101-21-3					1.6E+04	6.6E+04		1.3E+04
				1.0E-03	A			V		1.4E+09					Chlorpyrifos	2921-88-2					7.8E+01	3.3E+02		6.3E+01
				1.0E-02	H			V		1.4E+09					Chlorpyrifos Methyl	5598-13-0					7.8E+02	3.3E+03		6.3E+02
				5.0E-02	I			V		1.4E+09					Chlorsulfuron	64902-72-3					3.9E+03	1.6E+04		3.2E+03
				1.0E-02	I			V		1.4E+09					Chlorthal-dimethyl	1861-32-1					7.8E+02	3.3E+03		6.3E+02
				8.0E-04	H			V		1.4E+09					Chlorthiophos	60238-56-4					6.3E+01	2.6E+02		5.1E+01
				1.5E+00	I			V		1.4E+09	0.013				Chromium(III), insoluble Salts	16065-83-1					1.2E+05			1.2E+05
5.0E-01	J	8.4E-02	S	3.0E-03	I	1.0E-04	I	M		1.4E+09	0.025				Chromium(VI)	18540-29-9	3.1E+01	1.6E+03	3.0E+01		2.3E+02		1.4E+05	2.3E+02
								V		1.4E+09	0.013				Chromium, Total	7440-47-3								
				1.3E-02	I			V		1.4E+09					Clofentazine	74115-24-5					1.0E+03	4.3E+03		8.2E+02
				9.0E-03	P	3.0E-04	P	6.0E-06	P	1.4E+09					Cobalt	7440-48-4			4.2E+04	4.2E+04	2.3E+01		8.5E+03	2.3E+01
				6.2E-04	I			V	M	1.4E+09					Coke Oven Emissions	8007-45-2								
				4.0E-02	H			V		1.4E+09					Copper	7440-50-8					3.1E+03			3.1E+03
				5.0E-02	I	6.0E-01	C			1.4E+09					Cresol, m-	108-39-4					3.9E+03	1.6E+04	8.5E+08	3.2E+03
				5.0E-02	I	6.0E-01	C			1.4E+09					Cresol, o-	95-48-7					3.9E+03	1.6E+04	8.5E+08	3.2E+03
				1.0E-01	A	6.0E-01	C			1.4E+09					Cresol, p-	106-44-5					7.8E+03	3.3E+04	8.5E+08	6.3E+03
				1.0E-01	A			V		1.4E+09					Cresol, p-chloro-m-	59-50-7					7.8E+03	3.3E+04		6.3E+03
				1.0E-01	A	6.0E-01	C			1.4E+09					Cresols	1319-77-3					7.8E+03	3.3E+04	8.5E+08	6.3E+03
1.9E+00	H			1.0E-03	P			V		1.7E+04	1.4E+09	1.9E+04	1		Crotonaldehyde, trans-	123-73-9	3.7E+01			3.7E+01	7.8E+01			7.8E+01
				1.0E-01	I	4.0E-01	I	V		2.7E+02	1.4E+09	6.2E+03	1		Cumene	98-82-8					7.8E+03		2.6E+03	1.9E+03
2.2E-01	C	6.3E-05	C					V		1.4E+09					Cupferron	135-20-6	3.2E+02	1.1E+03	6.1E+06	2.5E+02	1.6E+02	6.6E+02		1.3E+02
8.4E-01	H			2.0E-03	H			V		1.4E+09					Cyanazine	21725-46-2	8.3E+01	2.9E+02		6.5E+01				
				1.0E-03	I			V		1.4E+09					Cyanides									
				5.0E-03	I			V		1.4E+09					~Calcium Cyanide	592-01-8					7.8E+01			7.8E+01
				6.0E-04	I	8.0E-04	S	V		9.5E+05	1.4E+09	5.3E+04	1		~Copper Cyanide	544-92-3					3.9E+02			3.9E+02
				1.0E-03	I			V		1.4E+09					~Cyanide (CN-)	57-12-5					4.7E+01		4.4E+01	2.3E+01
				1.0E-03	I			V		1.4E+09					~Cyanogen	460-19-5					7.8E+01			7.8E+01
				9.0E-02	I			V		1.4E+09					~Cyanogen Bromide	506-68-3					7.0E+03			7.0E+03
				5.0E-02	I			V		1.4E+09					~Cyanogen Chloride	506-77-4					3.9E+03			3.9E+03
				6.0E-04	I	8.0E-04	I	V		1.0E+07	1.4E+09	5.2E+04	1		~Hydrogen Cyanide	74-90-8					4.7E+01		4.4E+01	2.3E+01
				2.0E-03	I			V		1.4E+09					~Potassium Cyanide	151-50-8					1.6E+02			1.6E+02
				5.0E-03	I			V		1.4E+09	0.04				~Potassium Silver Cyanide	506-61-6					3.9E+02			3.9E+02
				1.0E-01	I			V		1.4E+09	0.04				~Silver Cyanide	506-64-9					7.8E+03			7.8E+03
				1.0E-03	I			V		1.4E+09					~Sodium Cyanide	143-33-9					7.8E+01			7.8E+01
				2.0E-04	P																			

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg- day)	k e y	RfC _o (mg/m ³) ⁻¹	k e y	muta- gen	C _{mt} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)			
				2.0E-04	X			V		1.4E+09			1	*Thiocyanic Acid	463-56-9							1.6E+01		1.6E+01		
				5.0E-02	I					1.4E+09			1	*Zinc Cyanide	557-21-1							3.9E+03		3.9E+03		
				6.0E+00	I	V		V	1.2E+02	1.4E+09	1.0E+03		1	Cyclohexane	110-82-7							6.5E+03		6.5E+03		
2.3E-02		H		5.0E+00	I	7.0E-01	P	V	5.1E+03	1.4E+09	4.2E+04		1	Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	3.0E+03	1.1E+04		2.4E+03			3.9E+05		3.0E+04	2.8E+04	
				5.0E-03	P	1.0E+00	X	V	2.8E+02	1.4E+09	1.5E+03		1	Cyclohexanone	108-94-1							3.9E+02		1.5E+03	3.1E+02	
				2.0E-01	I			V	2.9E+05	1.4E+09	7.5E+04		1	Cyclohexylamine	108-91-8							1.6E+04			1.6E+04	
				2.5E-02	I					1.4E+09			1	0.1	Cyfluthrin	68359-37-5						2.0E+03	8.2E+03		1.6E+03	
				5.0E-03	I					1.4E+09			1	0.1	Cyhalothrin	68085-85-8						3.9E+02	1.6E+03		3.2E+02	
				1.0E-02	I					1.4E+09			1	0.1	Cypermethrin	52315-07-8						7.8E+02	3.3E+03		6.3E+02	
				7.5E-03	I					1.4E+09			1	0.1	Cyromazine	66215-27-8						5.9E+02	2.5E+03		4.7E+02	
2.4E-01	I	6.9E-05	C							1.4E+09			1	0.1	DDD	72-54-8	2.9E+02	1.0E+03	5.5E+06	2.3E+02						
3.4E-01	I	9.7E-05	C					V		1.4E+09	2.1E+06		1		DDE, p,p'-	72-55-9	2.0E+02		6.1E+03	2.0E+02						
3.4E-01	I	9.7E-05	C	5.0E-04	I					1.4E+09			1	0.03	DDT	50-29-3	2.0E+02	2.4E+03	3.9E+06	1.9E+02			3.9E+01	5.5E+02	3.7E+01	
				3.0E-02	I					1.4E+09			1	0.1	Dalapon	75-99-0						2.3E+03	9.9E+03		1.9E+03	
1.8E-02	C	5.1E-06	C	1.5E-01	I					1.4E+09			1	0.1	Daminozide	1596-84-5	3.9E+03	1.4E+04	7.5E+07	3.0E+03			1.2E+04	4.9E+04	9.5E+03	
7.0E-04	I			7.0E-03	I					1.4E+09			1	0.1	Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209)	1163-19-5	9.9E+04	3.5E+05		7.8E+04			5.5E+02	2.3E+03	4.4E+02	
				4.0E-05	I					1.4E+09			1	0.1	Demeton	8065-48-3						3.1E+00	1.3E+01		2.5E+00	
1.2E-03	I			6.0E-01	I					1.4E+09			1	0.1	Di(2-ethylhexyl)adipate	103-23-1	5.8E+04	2.1E+05		4.5E+04			4.7E+04	2.0E+05		3.8E+04
6.1E-02	H			7.0E-04	A					1.4E+09			1	0.1	Diallate	2303-16-4	1.1E+03	4.1E+03		8.9E+02			5.5E+01	2.3E+02		4.4E+01
				1.0E-02	X			V		1.4E+09	5.2E+05		1		Dibenzothiophene	132-65-0							7.8E+02			7.8E+02
8.0E-01	P	6.0E-03	P	2.0E-04	I	2.0E-04	I	V	M	9.8E+02	1.4E+09	3.2E+04	1		Dibromo-3-chloropropane, 1,2-	96-12-8	1.9E+01		5.4E-01	5.3E-01			1.6E+01		6.7E+00	4.7E+00
				4.0E-04	X			V		1.6E+02	1.4E+09	1.9E+04	1		Dibromobenzene, 1,3-	108-36-1							3.1E+01			3.1E+01
				1.0E-02	I			V		1.4E+09	2.7E+04		1		Dibromobenzene, 1,4-	106-37-6							7.8E+02			7.8E+02
8.4E-02	I			2.0E-02	I			V		8.0E+02	1.4E+09	8.0E+03	1		Dibromochloromethane	124-48-1	8.3E+02			8.3E+02			1.6E+03			1.6E+03
2.0E+00	I	6.0E-04	I	9.0E-03	I	9.0E-03	I	V		1.3E+03	1.4E+09	8.6E+03	1		Dibromoethane, 1,2-	106-93-4	3.5E+01		4.0E+00	3.6E+00			7.0E+02		8.1E+01	7.3E+01
				4.0E-03	X	V		V		2.8E+03	1.4E+09	5.6E+03	1		Dibromomethane (Methylene Bromide)	74-95-3							2.4E+01			2.4E+01
				3.0E-04	P					1.4E+09			1	0.1	Dibutyltin Compounds	NA							2.3E+01	9.9E+01		1.9E+01
				3.0E-02	I					1.4E+09			1	0.1	Dicamba	1918-00-9							2.3E+03	9.9E+03		1.9E+03
4.2E-03	P							V		5.5E+02	1.4E+09	3.2E+03	1		Dichloro-2-butene, 1,4-	764-41-0			2.1E-01	2.1E-01						
4.2E-03	P							V		5.2E+02	1.4E+09	1.1E+04	1		Dichloro-2-butene, cis-1,4-	1476-11-5			7.4E-01	7.4E-01						
4.2E-03	P							V		7.6E+02	1.4E+09	1.1E+04	1		Dichloro-2-butene, trans-1,4-	110-57-6			7.4E-01	7.4E-01						
5.0E-02	I			4.0E-03	I					1.4E+09			1	0.1	Dichloroacetic Acid	79-43-6	1.4E+03	4.9E+03		1.1E+03			3.1E+02	1.3E+03		2.5E+02
				9.0E-02	I	2.0E-01	H	V		3.8E+02	1.4E+09	1.2E+04	1		Dichlorobenzene, 1,2-	95-50-1							7.0E+03		2.4E+03	1.8E+03
5.4E-03	C	1.1E-05	C	7.0E-02	A	8.0E-01	I	V		1.4E+09	1.0E+04		1		Dichlorobenzene, 1,4-	106-46-7	1.3E+04		2.7E+02	2.6E+02			5.5E+03		8.7E+03	3.4E+03
4.5E-01	I	3.4E-04	C							1.4E+09			1	0.1	Dichlorobenzidine, 3,3'-	91-94-1	1.5E+02	5.5E+02	1.1E+06	1.2E+02			7.0E+02	3.0E+03		5.7E+02
				9.0E-03	X					1.4E+09			1	0.1	Dichlorobenzophenone, 4,4'-	90-98-2							1.6E+04			8.7E+01
				2.0E-01	I	1.0E-01	X	V		8.5E+02	1.4E+09	8.4E+02	1		Dichlorodifluoromethane	75-71-8							8.8E+01			8.7E+01
5.7E-03	C	1.6E-06	C	2.0E-01	P					1.7E+03	1.4E+09	2.1E+03	1		Dichloroethane, 1,1-	75-34-3	1.2E+04		3.7E+02	3.6E+02			1.6E+04			1.6E+04
9.1E-02	I	2.6E-05	I	6.0E-03	X	7.0E-03	P	V		3.0E+03	1.4E+09	4.6E+03	1		Dichloroethane, 1,2-	107-06-2	7.6E+02		4.9E+01	4.6E+01			4.7E+02		3.3E+01	3.1E+01
				5.0E-02	I	2.0E-01	I	V		1.2E+03	1.4E+09	1.2E+03	1		Dichloroethylene, 1,1-	75-35-4							3.9E+03		2.4E+02	2.3E+02
				2.0E-03	I			V		2.4E+03	1.4E+09	2.5E+03	1		Dichloroethylene, 1,2-cis-	156-59-2							1.6E+02			1.6E+02
				2.0E-02	I			V		1.9E+03	1.4E+09	1.8E+03	1		Dichloroethylene, 1,2-trans-	156-60-5							1.6E+03			1.6E+03
				3.0E-03	I					1.4E+09			1	0.1	Dichlorophenol, 2,4-	120-83-2							2.3E+02	9.9E+02		1.9E+02
				1.0E-02	I					1.4E+09			1	0.05	Dichlorophenoxy Acetic Acid, 2,4-	94-75-7							7.8E+02	6.6E+03		7.0E+02
				8.0E-03	I					1.4E+09			1	0.1	Dichlorophenoxybutyric Acid, 4-(2,4-	94-82-6							6.3E+02	2.6E+03		5.1E+02
3.6E-02	C	1.0E-05	C	9.0E-02	A	4.0E-03	I	V		1.4E+03	1.4E+09	3.8E+03	1		Dichloropropane, 1,2-	78-87-5	1.9E+03		1.1E+02	1.0E+02			7.0E+03		1.6E+01	1.6E+01
				2.0E-02	P			V		1.5E+03	1.4E+09	6.8E+03	1		Dichloropropane, 1,3-	142-28-9							1.6E+03			1.6E+03
				3.0E-03	I					1.4E+09			1	0.1	Dichloropropanol, 2,3-	616-23-9							2.3E+02	9.9E+02		1.9E+02
1.0E-01	I	4.0E-06	I	3.0E-02	I	2.0E-02	I	V		1.6E+03	1.4E+09	3.6E+03	1		Dichloropropene, 1,3-	542-75-6	7.0E+02		2.5E+02	1.8E+02			2.3E+03		7.4E+01	7.2E+01
2.9E-01	I	8.3E-05	C	5.0E-04	I	5.0E-04	I			1.4E+09			1	0.1	Dichlorvos	62-73-7	2.4E+02	8.5E+02	4.6E+06	1.9E+02			3.9E+01	1.6E+02	7.1E+05	3.2E+01
				1.0E-04	I					1.4E+09			1	0.1	Dicrotophos	141-66-2							7.8E+00	3.3E+01		6.3E+00
				8.0E-02	P	3.0E-04	X	V		2.6E+02	1.4E+09	4.1E+03	1		Dicyclopentadiene	77-73-6							6.3E+03		1.3E+00	1.3E+00
1.6E+01	I	4.6E-03	I	5.0E-05	I					1.4E+09																

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRVT; A = ATSDR; C = Cal EPA; X = APPENDIX PPRVT SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)																							
Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³) ⁻¹	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³)	k _e (y)	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Child THQ=1 (mg/kg)	Child THQ=1 (mg/kg)	Child THQ=1 (mg/kg)	Child THI=1 (mg/kg)	
				1.0E-03	P		V		1.1E+05	1.4E+09	1.4E+05	1	0.1	Diethylformamide	617-84-5					7.8E+01			7.8E+01
3.5E+02	C	1.0E-01	C							1.4E+09			0.1	Diethylstilbestrol	56-53-1	2.0E-01	7.1E-01	3.8E+03	1.6E-01	6.3E+03	2.6E+04		5.1E+03
				8.0E-02	I					1.4E+09			0.1	Difenzoquat	43222-48-6					1.6E+03	6.6E+03		1.3E+03
				2.0E-02	I					1.4E+09			0.1	Diffubenzuron	35367-38-5								
4.4E-02	C	1.3E-05	C							1.4E+03	1.4E+09	1.2E+03	1	Difluoroethane, 1,1-	75-37-6							4.8E+04	4.8E+04
										1.4E+09	1.2E+05	1	0.1	Dihydrosafrole	94-58-6	1.6E+03	2.7E+03	9.9E+02					
										1.4E+09	3.1E+03	1	0.1	Diisopropyl Ether	108-20-3							2.2E+03	2.2E+03
				8.0E-02	I		V		5.3E+02	1.4E+09	3.8E+04	1	0.1	Diisopropyl Methylphosphonate	1445-75-6					6.3E+03			6.3E+03
				2.0E-02	I					1.4E+09			0.1	Dimethipin	55290-64-7					1.6E+03	6.6E+03		1.3E+03
				2.0E-04	I					1.4E+09			0.1	Dimethoate	60-51-5					1.6E+01	6.6E+01		1.3E+01
1.6E+00	P									1.4E+09			0.1	Dimethoxybenzidine, 3,3'-	119-90-4	4.3E+01	1.5E+02		3.4E+01				
1.7E-03	P			6.0E-02	P					1.4E+09			0.1	Dimethyl methylphosphonate	756-79-6	4.1E+04	1.5E+05		3.2E+04	4.7E+03	2.0E+04		3.8E+03
4.6E+00	C	1.3E-03	C							1.4E+09			0.1	Dimethylamino azobenzene [p-]	60-11-7	1.5E+01	5.4E+01	2.9E+05	1.2E+01				
5.8E-01	H									1.4E+09			0.1	Dimethylaniline HCl, 2,4-	21436-96-4	1.2E+02	4.3E+02		9.4E+01				
2.0E-01	P			2.0E-03	X					1.4E+09			0.1	Dimethylaniline, 2,4-	95-68-1	3.5E+02	1.2E+03		2.7E+02	1.6E+02	6.6E+02		1.3E+02
				2.0E-03	I		V		8.3E+02	1.4E+09	3.1E+04	1	0.1	Dimethylaniline, N,N-	121-69-7					1.6E+02			1.6E+02
1.1E+01	P									1.4E+09			0.1	Dimethylbenzidine, 3,3'-	119-93-7	6.3E+00	2.2E+01		4.9E+00				
				1.0E-01	P	3.0E-02	I	V		1.1E+05	1.4E+09	1.3E+05	1	Dimethylformamide	68-12-2					7.8E+03		4.0E+03	2.6E+03
				1.0E-04	X	2.0E-06	X	V		1.7E+05	1.4E+09	2.8E+04	1	Dimethylhydrazine, 1,1-	57-14-7					7.8E+00		5.8E-02	5.7E-02
5.5E+02	C	1.6E-01	C							1.9E+05	1.4E+09	1.7E+05	1	Dimethylhydrazine, 1,2-	540-73-8	1.3E-01	2.9E-01	8.8E-02					
				2.0E-02	I					1.4E+09			0.1	Dimethylphenol, 2,4-	105-67-9					1.6E+03	6.6E+03		1.3E+03
				6.0E-04	I					1.4E+09			0.1	Dimethylphenol, 2,6-	576-26-1					4.7E+01	2.0E+02		3.8E+01
4.5E-02	C	1.3E-05	C							1.4E+09			0.1	Dimethylphenol, 3,4-	95-65-8					7.8E+01	3.3E+02		6.3E+01
				8.0E-05	X					1.4E+09			0.1	Dimethylvinylchloride	513-37-1	1.5E+03	1.2E+02	1.1E+02					
										1.4E+09			0.1	Dinitro-o-cresol, 4,6-	534-52-1					6.3E+00	2.6E+01		5.1E+00
				2.0E-03	I					1.4E+09			0.1	Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5					1.6E+02	6.6E+02		1.3E+02
				1.0E-04	P					1.4E+09			0.1	Dinitrobenzene, 1,2-	528-29-0					7.8E+00	3.3E+01		6.3E+00
				1.0E-04	I					1.4E+09			0.1	Dinitrobenzene, 1,3-	99-65-0					7.8E+00	3.3E+01		6.3E+00
				1.0E-04	P					1.4E+09			0.1	Dinitrobenzene, 1,4-	100-25-4					7.8E+00	3.3E+01		6.3E+00
6.8E-01	I			2.0E-03	I					1.4E+09			0.1	Dinitrophenol, 2,4-	51-28-5					1.6E+02	6.6E+02		1.3E+02
										1.4E+09			0.1	Dinitrotoluene Mixture, 2,4/2,6-	NA	1.0E+02	3.6E+02	8.0E+01					
3.1E-01	C	8.9E-05	C	2.0E-03	I					1.4E+09		0.102		Dinitrotoluene, 2,4-	121-14-2	2.2E+02	7.8E+02	4.3E+06	1.7E+02	1.6E+02	6.5E+02		1.3E+02
1.5E+00	P			3.0E-04	X					1.4E+09		0.099		Dinitrotoluene, 2,6-	606-20-2	4.6E+01	1.7E+02		3.6E+01	2.3E+01	1.0E+02		1.9E+01
				2.0E-03	S					1.4E+09		0.006		Dinitrotoluene, 2-Amino-4,6-	35572-78-2					1.6E+02	1.1E+04		1.5E+02
4.5E-01	X			2.0E-03	S					1.4E+09		0.009		Dinitrotoluene, 4-Amino-2,6-	19406-51-0					1.6E+02	7.3E+03		1.5E+02
				9.0E-04	X					1.4E+09		0.1		Dinitrotoluene, Technical grade	15321-14-6	1.5E+02	5.5E+02	1.2E+02		7.0E+01	3.0E+02		5.7E+01
				1.0E-03	I					1.4E+09		0.1		Dinoseb	88-85-7					7.8E+01	3.3E+02		6.3E+01
1.0E-01	I	5.0E-06	I	3.0E-02	I	3.0E-02	I	V		1.2E+05	1.4E+09	4.0E+04	1	Dioxane, 1,4-	123-91-1	7.0E+02	2.2E+03	5.3E+02		2.3E+03		1.2E+03	8.1E+02
6.2E+03	I	1.3E+00	I							1.4E+09			0.03	Dioxins	NA								
										1.4E+09			0.03	*Hexachlorodibenzo-p-dioxin, Mixture	NA	1.1E-02	1.3E-01	2.9E+02	1.0E-02				
1.3E+05	C	3.8E+01	C	7.0E-10	I	4.0E-08	C	V		1.4E+09	2.0E+06	1	0.03	*TCDD, 2,3,7,8-	1746-01-6	5.3E-04	6.3E-03	1.4E-02	4.8E-04	5.5E-05	7.7E-04	8.2E-02	5.1E-05
				3.0E-02	I					1.4E+09			0.1	Diphenamid	957-51-7					2.3E+03	9.9E+03		1.9E+03
				8.0E-04	X					1.4E+09			0.1	Diphenyl Sulfone	127-63-9					6.3E+01	2.6E+02		5.1E+01
				2.5E-02	I					1.4E+09			0.1	Diphenylamine	122-39-4					2.0E+03	8.2E+03		1.6E+03
8.0E-01	I	2.2E-04	I							1.4E+09			0.1	Diphenylhydrazine, 1,2-	122-66-7	8.7E+01	3.1E+02	1.7E+06	6.8E+01				
				2.2E-03	I					1.4E+09			0.1	Diquat	85-00-7					1.7E+02	7.3E+02		1.4E+02
7.1E+00	C	1.4E-01	C							1.4E+09			0.1	Direct Black 38	1937-37-7	9.8E+00	3.5E+01	2.7E+03	7.6E+00				
7.4E+00	C	1.4E-01	C							1.4E+09			0.1	Direct Blue 6	2602-46-2	9.4E+00	3.3E+01	2.7E+03	7.3E+00				
6.7E+00	C	1.4E-01	C							1.4E+09			0.1	Direct Brown 95	16071-86-6	1.0E+01	3.7E+01	2.7E+03	8.1E+00				
				4.0E-05	I					1.4E+09			0.1	Disulfoton	298-04-4					3.1E+00	1.3E+01		2.5E+00
				1.0E-02	I		V			1.4E+09	4.5E+04	1	0.1	Dithiane, 1,4-	505-29-3					7.8E+02			7.8E+02
				2.0E-03	I					1.4E+09			0.1	Diuron	330-54-1					1.6E+02	6.6E+02		1.3E+02
				4.0E-03	I					1.4E+09			0.1	Dodine	2439-10-3					3.1E+02	1.3E+03		2.5E+02
				2.5E-02	I		V			1.4E+09	1.2E+05	1	0.1	EPTC	759-94-4					2.0E+03			2.0E+03
				6.0E-03	I		V			1.4E+09	4.1E+05	1	0.1	Endosulfan	115-29-7					4.7E+02			4.7E+02
				2.0E-02	I					1.4E+09			0.1	Endothall	145-73-3					1.6E+03	6.6E+03		1.3E+03
				3.0E-04	I					1.4E+09			0.1	Endrin	72-20-8					2.3E+01	9.9E+01		1.9E+01
9.9E-03	I	1.2E-06	I	6.0E-03	P	1.0E-03	I	V		1.1E+04	1.4E+09	1.9E+04	1	Epichlorohydrin	106-89-8	7.0E+03	4.4E+03	2.7E+03		4.7E+02		2.0E+01	1.9E+01

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRVT; A = ATSDR; C = Cal EPA; X = APPENDIX PPRVT SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)																									
Toxicity and Chemical-specific Information															Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1				
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg- day)	k e y	RfC _o (mg/m ³)	k e y	o m u t a g e n	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)		
				4.0E-02 5.0E-03	P I				1.4E+09 1.4E+09			1	0.1	Ethanol, 2-(2-methoxyethoxy)- Ethephon	111-77-3 16672-87-0					3.1E+03 3.9E+02	1.3E+04 1.6E+03		2.5E+03 3.2E+02		
				5.0E-04	I				1.4E+09			1	0.1	Ethion	563-12-2					3.9E+01	1.6E+02		3.2E+01		
				1.0E-01 9.0E-02	P P	6.0E-02 2.0E-01	V I		2.4E+04 1.1E+05	1.4E+09 1.4E+09	6.2E+04 9.8E+04	1	1	Ethoxyethanol Acetate, 2- Ethoxyethanol, 2-	111-15-9 110-80-5					7.8E+03 7.0E+03	3.8E+03 2.1E+04		2.6E+03 5.2E+03		
				9.0E-01 5.0E-03	I P	7.0E-02 8.0E-03	P I	V V	1.1E+04 2.5E+03	1.4E+09 1.4E+09	8.6E+03 6.3E+03	1	1	Ethyl Acetate Ethyl Acrylate	141-78-6 140-88-5					7.0E+04 3.9E+02	6.3E+02 5.3E+01		6.2E+02 4.7E+01		
				1.0E-01	I	1.0E+01	I	V	2.1E+03	1.4E+09	1.3E+03	1		Ethyl Chloride (Chloroethane)	75-00-3						1.4E+04		1.4E+04		
				2.0E-01 1.0E-05	I I			V P	1.0E+04 3.0E-01	1.4E+09 1.4E+09	3.1E+03 5.8E+03	1	0.1	Ethyl Ether Ethyl Methacrylate Ethyl-p-nitrophenyl Phosphonate	60-29-7 97-63-2 2104-64-5					1.6E+04 7.8E-01		1.8E+03 3.3E+00	1.6E+04 6.3E-01		
1.1E-02	C	2.5E-06	C	1.0E-01 7.0E-02 9.0E-02	I P P	1.0E+00 1.4E+09 1.9E+05	I V V		4.8E+02 1.4E+09 1.4E+09	1.4E+09 1.4E+09 1.4E+09	5.7E+03 1.8E+05	1	0.1	Ethylbenzene Ethylene Cyanohydrin Ethylene Diamine	100-41-4 109-78-4 107-15-3	6.3E+03		6.4E+02	5.8E+02	7.8E+03 5.5E+03 7.0E+03	2.3E+04	5.9E+03	3.4E+03 4.4E+03 7.0E+03		
3.1E-01	C	8.8E-05	C	2.0E+00 1.0E-01	I I	4.0E-01 1.6E+00	C I		1.4E+09 1.4E+09	1.4E+09	1.4E+09	1	0.1	Ethylene Glycol Ethylene Glycol Monobutyl Ether Ethylene Oxide	107-21-1 111-76-2 75-21-8	2.2E+02		1.9E+01	1.8E+01	1.6E+05 7.8E+03	6.6E+05 3.3E+04	5.7E+08 2.3E+09	1.3E+05 6.3E+03 1.9E+02		
4.5E-02 6.5E+01	C C	1.3E-05 1.9E-02	C C	8.0E-05 3.0E+00	I I	4.0E-01 1.6E+00	C V		1.4E+09 1.4E+09	1.4E+09	2.4E+04	1	0.1	Ethylene Thiourea Ethyleneimine Ethylphthalyl Ethyl Glycolate	96-45-7 151-56-4 84-72-0	1.5E+03 1.1E+00	5.5E+03	2.9E+07 3.5E-01	1.2E+03 2.7E-01	6.3E+00 2.3E+04	2.6E+01	5.7E+08 2.3E+09	5.1E+00 6.3E+03 1.9E+02		
				2.5E-04 2.5E-02 2.5E-02	I I I				1.4E+09 1.4E+09 1.4E+09	1.4E+09	1.4E+04	1	0.1	Fenamiphos Fenproprathrin Fenvalerate	22224-92-6 39515-41-8 51630-58-1					2.0E+01 2.0E+03 2.0E+03	8.2E+01 8.2E+03 8.2E+03		1.6E+01 1.6E+03 1.6E+03		
				1.3E-02 4.0E-02 6.0E-02	I C I	1.4E+09 1.3E-02 1.3E-02	C C		1.4E+09 1.4E+09	1.4E+09	1.4E+09	1	0.1	Fluometuron Fluoride Fluorine (Soluble Fluoride)	2164-17-2 16984-48-8 7782-41-4					1.0E+03 3.1E+03 4.7E+03	4.3E+03	1.8E+07 1.8E+07	8.2E+02 3.1E+03 4.7E+03		
				8.0E-02 2.0E-02 7.0E-04	I I I				1.4E+09 1.4E+09 1.4E+09	1.4E+09	1.4E+09	1	0.1	Fluridone Flurprimidol Flusilazole	59756-60-4 56425-91-3 85509-19-9					6.3E+03 1.6E+03 5.5E+01	2.6E+04 6.6E+03 2.3E+02		5.1E+03 1.3E+03 4.4E+01		
3.5E-03	I			6.0E-02 1.0E-02 1.0E-01	I I I				1.4E+09 1.4E+09 1.4E+09	1.4E+09	1.4E+09	1	0.1	Flutolanil Fluvalinate Folpet	66332-96-5 9409-94-5 133-07-3				1.6E+04	2.0E+04 7.1E+04			4.7E+03 7.8E+02 7.8E+03	2.0E+04 3.3E+03 3.3E+04	3.8E+03 6.3E+02 6.3E+03
1.9E-01	I			2.0E-03 1.3E-05	I I				1.4E+09 1.4E+09	1.4E+09	7.8E+04	1	0.1	Fomesafen Fonofos Formaldehyde	72178-02-0 44-22-9 50-00-0	3.7E+02	1.3E+03		2.9E+02	1.6E+02 1.6E+04	6.6E+02	8.0E+02	1.3E+02 7.6E+02		
				9.0E-01 3.0E+00	P I	3.0E-04 1.4E+09	X I	V V	1.1E+05 1.4E+09	1.4E+09	9.3E+04	1	0.1	Formic Acid Fosetyl-AL Furans	64-18-6 39148-24-8					7.0E+04 2.3E+05	2.9E+01 9.9E+05		2.9E+01 1.9E+05		
				1.0E-03 1.0E-03 9.0E-01	X I I			V V V	1.4E+09 6.2E+03 1.7E+05	1.4E+09 1.4E+09 1.4E+09	1.6E+05 2.6E+03 1.2E+04	1	0.03	*Dibenzofuran *Furan *Tetrahydrofuran	132-64-9 110-00-9 109-99-9					7.8E+01 7.8E+01 7.0E+04	1.1E+03 1.1E+03 9.9E+05		7.3E+01 7.3E+01 1.8E+04		
3.8E+00 1.5E+00	H C			3.0E-03	I	5.0E-02	H V		1.0E+04	1.4E+09	4.9E+04	1	0.1	Furazolidone Furfural Furium	67-45-8 98-01-1 531-82-8	1.8E+01	6.5E+01		1.4E+01	2.3E+02		2.5E+03	2.1E+02		
3.0E-02	I	8.6E-06	C	4.0E-04	I				1.4E+09	1.4E+09	1.4E+09	1	0.1	Furmecyclox Glufosinate, Ammonium Glutaraldehyde	60568-05-0 77182-82-2 111-30-8	2.3E+03	8.2E+03	4.4E+07	1.8E+03	3.1E+01	1.3E+02		1.1E+05	2.5E+01 1.1E+05	
				4.0E-04 1.0E-01 1.0E-02	I I X	1.0E-03	H V		1.1E+05	1.4E+09	8.4E+04	1	0.1	Glycidyl Glyphosate Guanidine	765-34-4 1071-83-6 113-00-8					3.1E+01 7.8E+03 7.8E+02	3.3E+04	8.8E+01	2.3E+01 6.3E+03 7.8E+02		
4.5E+00 9.1E+00	I I	1.3E-03 2.6E-03	I I	2.0E-02 5.0E-04	P I			V V	1.4E+09 1.4E+09	1.4E+09	4.8E+05	1	0.1	Guanidine Chloride Haloxypol, Methyl Heptachlor	50-01-1 69806-40-2 76-44-8	1.5E+01		1.0E+02	1.3E+01	1.6E+03 3.9E+00 3.9E+01	6.6E+03 1.6E+01		1.3E+03 3.2E+00 3.9E+01		
				2.0E-02 5.0E-05 2.0E-03 2.0E-04	P I I I			V V V	1.4E+09 1.4E+09	1.4E+09	8.4E+05	1	0.1	Heptachlor Epoxide Hexabromobenzene Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BDE-153)	1024-57-3 87-82-1 68631-49-2	7.6E+00		9.1E+01	7.0E+00	1.0E+00 1.6E+02 1.6E+01	1.3E+02 6.6E+01		1.0E+00 1.6E+02 1.3E+01		
1.6E+00 7.8E-02 6.3E+00	I I I	4.6E-04 2.2E-05 1.8E-03	I I A	8.0E-04 1.0E-03 8.0E-03	I P A			V V	1.4E+09 1.7E+01 1.4E+09	1.4E+09	6.8E+04 1.1E+04	1	0.1	Hexachlorobenzene Hexachlorobutadiene Hexachlorocyclohexane, Alpha-	118-74-1 87-68-3 319-84-6	4.3E+01 8.9E+02 1.1E+01	4.1E+01 1.4E+02	2.1E+01 1.2E+02		6.3E+01 7.8E+01 6.3E+02	2.6E+03		6.3E+01 7.8E+01 2.6E+03		

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRVT; A = ATSDR; C = Cal EPA; X = APPENDIX PPRVT SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg- day)	k e y	RfC _o (mg/m ³) ⁻¹	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-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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	I			3.0E-04	S				1.4E+09				0.07	**Mercuric Chloride (and other Mercury salts)	7487-94-7					2.3E+01		4.3E+05	2.3E+01
				3.0E-04	I	V			3.1E+00	1.4E+09	3.5E+04		1	**Mercury (elemental)	7439-97-6							1.1E+01	1.1E+01
1.0E-04	I			8.0E-05	I				1.4E+09	1.4E+09			1	**Methyl Mercury	22967-92-6					7.8E+00			7.8E+00
				8.0E-05	I				1.4E+09	1.4E+09			0.1	**Phenylmercuric Acetate	62-38-4					6.3E+00	2.6E+01		5.1E+00
3.0E-05	I			3.0E-05	I			V	1.4E+09	1.9E+06			1	Merphos	150-50-5					2.3E+00			2.3E+00
3.0E-05	I			6.0E-02	I				1.4E+09	1.4E+09			1	Merphos Oxide	78-48-8					2.3E+00	9.9E+00		1.9E+00
6.0E-02	I			6.0E-02	I				1.4E+09	1.4E+09			0.1	Metalaxyl	57837-19-1					4.7E+03	2.0E+04		3.8E+03
1.0E-04	I			5.0E-05	I			P	4.6E+03	1.4E+09	6.8E+03		1	Methacrylonitrile	126-98-7					7.8E+00		2.1E+02	7.5E+00
2.0E+00	I			2.0E+00	I	V			1.1E+05	1.4E+09	2.9E+04		1	Methamidophos	10265-92-6					3.9E+00	1.6E+01		3.2E+00
1.0E-03	I			2.5E-02	I				1.4E+09	1.4E+09			1	Methanol	67-56-1					1.6E+05		6.1E+05	1.2E+05
4.9E-02	C	1.4E-05	C	2.5E-02	I				1.4E+09	1.4E+09			1	Methidathion	950-37-8					7.8E+01	3.3E+02		6.3E+01
				2.5E-02	I				1.4E+09	1.4E+09			1	Methomyl	16752-77-5					2.0E+03	8.2E+03		1.6E+03
				5.0E-03	I				1.4E+09	1.4E+09			1	Methoxy-5-nitroaniline, 2-	99-59-2	1.4E+03	5.0E+03	2.7E+07	1.1E+03				
				8.0E-03	P	1.0E-03	P	V	1.2E+05	1.4E+09	1.2E+05		1	Methoxychlor	72-43-5					3.9E+02	1.6E+03		3.2E+02
				5.0E-03	P	2.0E-02	I	V	1.1E+05	1.4E+09	1.0E+05		1	Methoxyethanol Acetate, 2-	110-49-6					6.3E+02		1.3E+02	1.1E+02
				1.0E+00	X			V	2.9E+04	1.4E+09	8.1E+03		1	Methoxyethanol, 2-	109-86-4					3.9E+02		2.1E+03	3.3E+02
				2.0E-02	P			V	6.8E+03	1.4E+09	7.0E+03		1	Methyl Acetate	79-20-9					7.8E+04			7.8E+04
				6.0E-01	I	5.0E+00	I	V	2.8E+04	1.4E+09	1.2E+04		1	Methyl Acrylate	96-33-3					4.7E+04		1.5E+02	1.5E+02
				1.0E-03	X			P	1.8E+05	1.4E+09	5.0E+04		1	Methyl Ethyl Ketone (2-Butanone)	78-93-3					4.7E+04		6.4E+04	2.7E+04
				3.0E+00	I			V	3.4E+03	1.4E+09	1.1E+04		1	Methyl Hydrazine	60-34-4			1.4E+01	1.4E+01	7.8E+01		1.1E+00	1.0E+00
				1.0E-03	C			V	1.0E+04	1.4E+09	4.4E+03		1	Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1							3.3E+04	3.3E+04
				1.4E+00	I	7.0E-01	I	V	2.4E+03	1.4E+09	6.3E+03		1	Methyl Isocyanate	624-83-9							4.6E+00	4.6E+00
				2.5E-04	I				1.4E+09	1.4E+09			0.1	Methyl Methacrylate	80-62-6					1.1E+05		4.6E+03	4.4E+03
				6.0E-02	X				1.4E+09	1.4E+09			1	Methyl Parathion	298-00-0					2.0E+01	8.2E+01		1.6E+01
				6.0E-02	X				1.4E+09	1.4E+09			1	Methyl Phosphonic Acid	993-13-5					4.7E+03	2.0E+04		3.8E+03
9.9E-02	C	2.8E-05	C	6.0E-03	H	4.0E-02	H	V	3.9E+02	1.4E+09	2.4E+04		1	Methyl Styrene (Mixed Isomers)	25013-15-4					4.7E+02		1.0E+03	3.2E+02
1.8E-03	C	2.6E-07	C	3.0E+00	I	V			8.9E+03	1.4E+09	4.9E+03		1	Methyl methanesulfonate	66-27-3	7.0E+02	2.5E+03	1.4E+07	5.5E+02				
				3.0E+00	I	V			8.9E+03	1.4E+09	4.9E+03		1	Methyl tert-Butyl Ether (MTBE)	1634-04-4	3.9E+04	5.3E+03	4.7E+03				1.5E+04	1.5E+04
9.0E-03	P			2.0E-02	X				1.4E+09	1.4E+09			1	Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2					2.3E+01	9.9E+01		1.9E+01
8.3E+00	C	2.4E-03	C	3.0E+00	I	V			1.4E+09	1.4E+09			1	Methyl-5-Nitroaniline, 2-	99-55-8	7.7E+03	2.7E+04		6.0E+03	1.6E+03	6.6E+03		1.3E+03
				3.0E+00	I	V			1.4E+09	1.4E+09			1	Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	8.4E+00	3.0E+01	1.6E+05	6.5E+00				
1.3E-01	C	3.7E-05	C	1.0E-02	A				1.4E+09	1.4E+09			1	Methylaliline Hydrochloride, 2-	616-21-5	5.3E+02	1.9E+03	1.0E+07	4.2E+02				
				2.0E-04	X				1.4E+09	1.4E+09			1	Methylarsonic acid	114-58-3					7.8E+02	3.3E+03		6.3E+02
1.0E-01	X			2.0E-04	X				1.4E+09	1.4E+09			1	Methylbenzene, 1,4-diamine monohydrochloride, 2-	74612-12-7					1.6E+01	6.6E+01		1.3E+01
2.2E+01	C	6.3E-03	C	3.0E-04	X				1.4E+09	1.4E+09			1	Methylbenzene, 1,4-diaminesulfate, 2-	615-50-9	7.0E+02	2.5E+03		5.4E+02	2.3E+01	9.9E+01		1.9E+01
2.0E-03	I	1.0E-08	I	6.0E-03	I	6.0E-01	I	V	3.3E+03	1.4E+09	2.2E+03		1	Methylcholanthrene, 3-	56-49-5	7.0E-01	2.7E+00	2.2E+04	5.5E-01				
1.0E-01	P	4.3E-04	C	2.0E-03	P			M	1.4E+09	1.4E+09			1	Methylene Chloride	75-09-2	7.7E+03	2.2E+04	5.7E+03		4.7E+02		1.4E+03	3.5E+02
4.6E-02	I	1.3E-05	C	1.6E+00	C	4.6E-04	C		1.4E+09	1.4E+09			1	Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	1.5E+02	6.0E+02	3.2E+05	1.2E+02	1.6E+02	6.6E+02		1.3E+02
1.6E+00	C	4.6E-04	C	2.0E-02	C				1.4E+09	1.4E+09			1	Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	1.5E+03	5.4E+03	2.9E+07	1.2E+03				
				6.0E-04	I				1.4E+09	1.4E+09			1	Methylenebisbenzenamine, 4,4'-	101-77-9	4.3E+01	1.5E+02	8.3E+05	3.4E+01			2.8E+07	2.8E+07
				7.0E-02	H			V	5.0E+02	1.4E+09	1.3E+04		1	Methylenediphenyl Diisocyanate	101-68-8							8.5E+05	8.5E+05
				1.5E-01	I				1.4E+09	1.4E+09			1	Methylstyrene, Alpha-	98-83-9					5.5E+03			5.5E+03
				2.5E-02	I				1.4E+09	1.4E+09			1	Metolachlor	51218-45-2					1.2E+04	4.9E+04		9.5E+03
				2.5E-01	I				1.4E+09	1.4E+09			1	Metribuzin	21087-64-9					2.0E+03	8.2E+03		1.6E+03
				3.0E+00	P			V	3.4E-01	1.4E+09	1.4E+03		1	Metsulfuron-methyl	74223-64-6					2.0E+04	8.2E+04		1.6E+04
1.8E+01	C	5.1E-03	C	2.0E-04	I			V	1.4E+09	8.6E+05			1	Mineral oils	8012-95-1	3.9E+00		4.7E+01	3.6E+00	2.3E+05			2.3E+05
				2.0E-03	I				1.4E+09	1.4E+09			1	Mirex	2385-85-5					1.6E+01			1.6E+01
				5.0E-03	I				1.4E+09	1.4E+09			1	Molinate	2212-67-1					1.6E+02	6.6E+02		1.3E+02
				1.0E-01	I				1.4E+09	1.4E+09			1	Molybdenum	7439-98-7					3.9E+02			3.9E+02
				2.0E-03	P				1.4E+09	1.4E+09			1	Monochloramine	10599-90-3					7.8E+03			7.8E+03
				2.5E-02	I				1.4E+09	1.4E+09			1	Monomethylaniline	100-61-8					1.6E+02	6.6E+02		1.3E+02
				3.0E-04	X				1.4E+09	1.4E+09			1	Myclobutanil	88671-89-0					2.0E+03	8.2E+03		1.6E+03
				2.0E-03	I			V	1.4E+09	5.7E+04			1	N,N'-Diphenyl-1,4-benzenediamine	74-31-7					2.3E+01	9.9E+01		1.9E+01
				3.0E-02	X	1.0E-01	P	V	1.4E+09	1.4E+09			1	Naled	300-76-5					1.6E+02			1.6E+02
				2.6E-04	C	1.1E-02	C	1.4E-05	C	1.4E+09	1.4E+09		1	Naphtha, High Flash Aromatic (HFAN)	64742-95-6					2.3E+03		1.4E+08	2.3E+03
1.8E+00	C	0.0E+00	C	1.0E-01	I				1.4E+09	1.4E+09			1	Naphthylamine, 2-	91-59								

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg- day)	k e y	RfC _o (mg/m ³) ⁻¹	k e y	muta- gen	C _{mt} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)
2.6E-04	C	1.1E-02	C	1.4E-05	C	V				1.4E+09			1	Nickel Carbonyl	13463-39-3			1.5E+06	1.5E+06	8.6E+02		2.0E+04	8.2E+02
2.6E-04	C	1.1E-02	C	1.4E-05	C					1.4E+09			0.04	Nickel Hydroxide	12054-48-7			1.5E+06	1.5E+06	8.6E+02		2.0E+04	8.2E+02
2.6E-04	C	1.1E-02	C	2.0E-05	C					1.4E+09			0.04	Nickel Oxide	1313-99-1			1.5E+06	1.5E+06	8.6E+02		2.8E+04	8.4E+02
2.4E-04	I	1.1E-02	C	1.4E-05	C					1.4E+09			0.04	Nickel Refinery Dust	NA			1.6E+06	1.6E+06	8.6E+02		2.0E+04	8.2E+02
2.6E-04	C	2.0E-02	I	9.0E-05	A					1.4E+09			0.04	Nickel Soluble Salts	7440-02-0			1.5E+06	1.5E+06	1.6E+03		1.3E+05	1.5E+03
1.7E+00	C	4.8E-04	I	1.1E-02	C	1.4E-05	C			1.4E+09			0.04	Nickel Subsulfide	12035-72-2	4.1E+01		8.0E+05	4.1E+01	8.6E+02		2.0E+04	8.2E+02
2.6E-04	C	1.1E-02	C	1.4E-05	C					1.4E+09			1	Nickelocene	1271-28-9			1.5E+06	1.5E+06	8.6E+02	3.6E+03	2.0E+04	6.7E+02
1.6E+00										1.4E+09			1	Nitrate	14797-55-8					1.3E+05		2.0E+04	1.3E+05
				1.0E-01	I					1.4E+09			1	Nitrate + Nitrite (as N)	NA								7.8E+03
				1.0E-02	X	5.0E-05	X			1.4E+09			1	Nitrite	14797-65-0								7.8E+03
				1.0E-02	X	5.0E-05	X			1.4E+09			0.1	Nitroaniline, 2-	88-74-4								6.3E+02
2.0E-02	P	4.0E-03	P	6.0E-03	P					1.4E+09			1	Nitroaniline, 4-	100-01-6	3.5E+03	1.2E+04	5.1E+02	2.7E+03	3.1E+02	1.3E+03	8.5E+06	2.5E+02
		2.0E-03	I	9.0E-03	I	V		3.1E+03	1.4E+09	7.3E+04			1	Nitrobenzene	98-95-3				1.6E+02	1.6E+02	6.9E+02		1.3E+02
		3.0E+03	P						1.4E+09				0.1	Nitrocellulose	9004-70-0				5.1E+02	2.3E+08	9.9E+08		1.9E+08
1.3E+00	C	3.7E-04	C	7.0E-02	H					1.4E+09			1	Nitrofurantoin	67-20-9	5.3E+01	1.9E+02	1.0E+06	4.2E+01	5.5E+03	2.3E+04		4.4E+03
1.7E-02	P	1.0E-04	P							1.4E+09			1	Nitrofurazone	59-87-0	4.1E+03	1.5E+04		3.2E+03	7.8E+00	3.3E+01		6.3E+00
		1.0E-01	I							1.4E+09			1	Nitroglycerin	55-63-0					7.8E+00	3.3E+01		6.3E+00
8.8E-06	P	2.7E-03	H	5.0E-03	P	V		1.8E+04	1.4E+09	1.7E+04			1	Nitroguanidine	556-88-7					7.8E+03	3.3E+04		6.3E+03
2.7E+01	C	7.7E-03	C	2.0E-02	I	V		4.9E+03	1.4E+09	1.3E+04			1	Nitromethane	75-52-5			5.4E+02	5.4E+02			8.8E+01	2.7E+02
1.2E+02	C	3.4E-02	C							1.4E+09			1	Nitropropane, 2-	79-46-9			1.4E+00	1.4E+00			2.7E+02	2.7E+02
5.4E+00	I	1.6E-03	I							1.4E+09	2.4E+05		1	Nitroso-N-ethylurea, N-	759-73-9	5.7E-01	2.2E+00	1.8E+04	4.5E-01	1.3E-01	5.0E-01	4.1E+03	1.0E-01
7.0E+00	I	2.0E-03	C							1.4E+09			1	Nitroso-N-methylurea, N-	684-93-5	1.3E-01	5.0E-01	4.1E+03	1.0E-01	1.3E+01	4.3E+01	9.9E+00	
2.8E+00	I	8.0E-04	C							1.4E+09			1	Nitroso-di-N-butylamine, N-	924-16-3	1.3E+01		4.3E+01	9.9E+00				
1.5E+02	I	4.3E-02	I							1.4E+09			1	Nitroso-di-N-propylamine, N-	621-64-7	9.9E+00	3.5E+01	1.9E+05	7.8E+00				
5.1E+01	I	1.4E-02	I	8.0E-06	P	4.0E-05	X	V	M	2.4E+05	1.4E+09	8.2E+04	1	Nitrosodimethylamine, N-	62-75-9	3.0E-01		6.0E-01	2.0E-01	6.3E-01		3.4E+00	5.3E-01
4.9E-03	I	2.6E-06	C							1.4E+09			1	Nitrosodiethanolamine, N-	1116-54-7	2.5E+01	8.8E+01	4.8E+05	1.9E+01				
2.2E+01	I	6.3E-03	C							1.4E+09			1	Nitrosodithylamine, N-	55-18-5	1.0E-01	4.0E-01	3.2E+03	8.1E-02				
6.7E+00	C	1.9E-03	C							1.4E+09			1	Nitrosodiphenylamine, N-	86-30-6	1.4E+04	5.0E+04	1.5E+08	1.1E+04				
9.4E+00	C	2.7E-03	C							1.4E+09			1	Nitrosomethylethylamine, N-	10595-95-6	3.2E+00		5.4E+00	2.0E+00				
2.1E+00	I	6.1E-04	I							1.4E+09			1	Nitrosomorpholine [N-]	59-89-2	1.0E+01	3.7E+01	2.0E+05	8.1E+00				
				1.0E-04	X					1.4E+09			1	Nitrosopiperidine [N-]	180-75-4	7.4E+00	2.6E+01	1.4E+05	5.8E+00				
2.2E-01	P	9.0E-04	P					1.5E+03	1.4E+09	1.4E+05			1	Nitrosopyrrolidine, N-	910-55-2	3.3E+01	1.2E+02	6.3E+05	2.6E+01	3.2E+02	4.3E+03	1.5E+04	3.2E+02
1.6E-02	P	4.0E-03	P							1.4E+09			1	Nitrotoluene, m-	99-08-1					7.8E+00	3.3E+01		6.3E+00
				3.0E-04	X	2.0E-02	P	V		6.9E+00	1.4E+09	1.0E+03	1	Nitrotoluene, o-	88-72-2	3.2E+02			3.2E+02	7.0E+01		7.0E+01	
				3.0E-03	I					1.4E+09			1	Nitrotoluene, p-	99-99-0	4.3E+03	1.5E+04		3.4E+03	3.1E+02	1.3E+03		2.5E+02
				3.0E-04	X	2.0E-02	P	V		6.9E+00	1.4E+09	1.0E+03	1	Nonane, n-	111-84-2					2.3E+01		2.2E+01	1.1E+01
				4.0E-02	I					1.4E+09			1	Norflurazon	27314-13-2					3.1E+03	1.3E+04		2.5E+03
				3.0E-03	I					1.4E+09			1	Octabromodiphenyl Ether	32536-52-0					2.3E+02	9.9E+02		1.9E+02
				5.0E-02	I					1.4E+09			1	Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0					3.9E+03	2.7E+05		3.9E+03
				2.0E-03	H					1.4E+09			1	Octamethylpyrophosphoramide	152-16-9					1.6E+02	6.6E+02		1.3E+02
				5.0E-02	I					1.4E+09			1	Oryzalin	19044-88-3					3.9E+03	1.6E+04		3.2E+03
				5.0E-03	I					1.4E+09			1	Oxadiazon	19666-30-9					3.9E+02	1.6E+03		3.2E+02
				2.5E-02	I					1.4E+09			1	Oxamyl	23135-22-0					2.0E+03	8.2E+03		1.6E+03
				3.0E-03	I					1.4E+09			1	Oxyfluorfen	42874-03-3					2.3E+02	9.9E+02		1.9E+02
				1.3E-02	I					1.4E+09			1	Paclitaxel	76738-62-0					1.0E+03	4.3E+03		8.2E+02
				4.5E-03	I					1.4E+09			1	Paraquat Dichloride	1910-42-5					3.5E+02	1.5E+03		2.8E+02
				6.0E-03	H					1.4E+09			1	Parathion	56-38-2					4.7E+02	2.0E+03		3.8E+02
				5.0E-02	H					1.4E+09	4.5E+04		1	Pebulate	1114-71-2					3.9E+03			3.9E+03
				4.0E-02	I					1.4E+09			1	Pendimethalin	40487-42-1					3.1E+03	1.3E+04		2.5E+03
				2.0E-03	I			3.1E-01	1.4E+09	5.1E+05			1	Pentabromodiphenyl Ether	32534-81-9					1.6E+02			1.6E+02
				1.0E-04	I					1.4E+09			1	Pentabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-99)	60348-60-9					7.8E+00	3.3E+01		6.3E+00
9.0E-02	P	8.0E-04	I							1.4E+09	8.1E+04		1	Pentachlorobenzene	608-93-5					6.3E+01			6.3E+01
								4.6E+02	1.4E+09	9.7E+03			1	Pentachloroethane	76-01-7	7.7E+02			7.7E+02				
2.6E-01	H	3.0E-03	I							1.4E+09	4.3E+05		1	Pentachloronitrobenzene	82-68-8	2.7E+02			2.7E+02	2.3E+02			2.3E+02
4.0E-01	I	5.1E-06	C	5.0E-03	I					1.4E+09			0.25	Pentachlorophenol	87-86-5	1.7E+02	2.5E+02	7.5E+07	1.0E+02	3.9E+02	6.6E+02		2.5E+02
4.0E-03	X	2.0E-03	P							1.4E+09			0.1	Pentaerythritol tetranitrate (PETN)	78-11-5	1.7E+04	6.2E+04		1.4E+04	1.6E+02	6.6E+02		1.3E+02
				1.0E+00	P	V		3.9E+02	1.4E+09	7.8E+02			1	Pentane, n-	109-66-0							8.1E+02	8.1E+02

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg- day)	k e y	RfC _o (mg/m ³) ⁻¹	k e y	muta- gen	C _{soil} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)		
				7.0E-04	I					1.4E+09			1	~Lithium Perchlorate	7791-03-9							5.5E+01	5.5E+01		
				7.0E-04	I					1.4E+09			1	~Perchlorate and Perchlorate Salts	14797-73-0							5.5E+01	5.5E+01		
				7.0E-04	I					1.4E+09			1	~Potassium Perchlorate	7778-74-7							5.5E+01	5.5E+01		
				7.0E-04	I					1.4E+09			1	~Sodium Perchlorate	7601-89-0							5.5E+01	5.5E+01		
				2.0E-02	P			V		1.4E+09	1.3E+05		1	Perfluorobutane Sulfonate	375-73-5							1.6E+03	1.6E+03		
				5.0E-02	I					1.4E+09			1	0.1	Permethrin	52645-53-1						3.9E+03	1.6E+04	3.2E+03	
2.2E-03	C	6.3E-07	C							1.4E+09			1	0.1	Phenacetin	62-44-2	3.2E+04	1.1E+05	6.1E+08	2.5E+04					
				2.5E-01	I					1.4E+09			1	0.1	Phenmedipham	13684-63-4						2.0E+04	8.2E+04	1.6E+04	
				3.0E-01	I	2.0E-01	C			1.4E+09			1	0.1	Phenol	108-95-2						2.3E+04	9.9E+04	2.8E+08	1.9E+04
				4.0E-03	I					1.4E+09			1	0.1	Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1						3.1E+02	1.3E+03	2.5E+02	
				5.0E-04	X					1.4E+09			1	0.1	Phenothiazine	92-84-2						3.9E+01	1.6E+02	3.2E+01	
				6.0E-03	I					1.4E+09			1	0.1	Phenylenediamine, m-	108-45-2						4.7E+02	2.0E+03	3.8E+02	
4.7E-02	H									1.4E+09			1	0.1	Phenylenediamine, o-	95-54-5	1.5E+03	5.3E+03		1.2E+03					
				1.9E-01	H					1.4E+09			1	0.1	Phenylenediamine, p-	106-50-3						1.5E+04	6.3E+04	1.2E+04	
1.9E-03	H									1.4E+09			1	0.1	Phenylphenol, 2-	90-43-7	3.6E+04	1.3E+05		2.8E+04					
				2.0E-04	H					1.4E+09			1	0.1	Phorate	298-02-2						1.6E+01	6.6E+01	1.3E+01	
						3.0E-04	I	V		1.6E+03	1.4E+09	9.8E+02	1		Phosgene	75-44-5							3.1E-01	3.1E-01	
				2.0E-02	I					1.4E+09			1	0.1	Phosmet	732-11-6						1.6E+03	6.6E+03	1.3E+03	
															Phosphates, Inorganic										
				4.9E+01	P					1.4E+09			1		~Aluminum metaphosphate	13776-88-0						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Ammonium polyphosphate	68333-79-9						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Calcium pyrophosphate	7790-76-3						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Diammonium phosphate	7783-28-0						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Dicalcium phosphate	7757-93-9						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Dimagnesium phosphate	7782-75-4						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Dipotassium phosphate	7758-11-4						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Disodium phosphate	7558-79-4						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Monoaluminum phosphate	13530-50-2						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Monoammonium phosphate	7722-76-1						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Monocalcium phosphate	7758-23-8						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Monomagnesium phosphate	7757-86-0						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Monopotassium phosphate	7778-77-0						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Monosodium phosphate	7758-80-7						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Polyphosphoric acid	8017-16-1						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Potassium triphosphate	11845-36-8						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Sodium acid pyrophosphate	7758-16-9						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Sodium aluminum phosphate (acidic)	7785-88-8						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Sodium aluminum phosphate (anhydrous)	10279-59-1						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Sodium aluminum phosphate (tetrahydrate)	10305-76-7						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Sodium hexametaphosphate	10124-56-8						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Sodium polyphosphate	68915-31-1						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Sodium trimetaphosphate	7785-84-4						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Sodium triphosphate	7758-29-4						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Tetrapotassium phosphate	7320-34-5						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Tetrasodium pyrophosphate	7722-88-5						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Trialuminum sodium tetra decahydrogenooctaoorthophosphate (dihydrate)	15136-87-5						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Tricalcium phosphate	7758-87-4						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Trimagnesium phosphate	7757-87-1						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Tripotassium phosphate	7778-53-2						3.8E+06		3.8E+06	
				4.9E+01	P					1.4E+09			1		~Trisodium phosphate	7601-54-9						3.8E+06		3.8E+06	
				3.0E-04	I	3.0E-04	I	V		1.4E+09			1		Phosphine	7803-51-2						2.3E+01		4.3E+05	2.3E+01
				4.9E+01	P	1.0E-02	I	V		1.4E+09			1		Phosphoric Acid	7664-38-2						3.8E+06		1.4E+07	3.0E+06
				2.0E-05	I			V		1.4E+09	6.9E+03		1		Phosphorus, White	7723-14-0						1.6E+00		1.6E+00	
															Phthalates										
1.4E-02	I	2.4E-06	C	2.0E-02	I					1.4E+09			1	0.1	~Bis(2-ethylhexyl)phthalate	117-81-7	5.0E+03	1.8E+04	1.6E+08	3.9E+03	1.6E+03	6.6E+03		1.3E+03	
1.9E-03	P			2.0E-01	I					1.4E+09			1	0.1	~Butyl Benzyl Phthalate	85-68-7	3.7E+04	1.3E+05		2.9E+04	1.6E+04	6.6E+04		1.3E+04	
				1.0E+00	I					1.4E+09			1	0.1	~Butylphthalyl Butylglycolate	85-70-1					7.8E+04	3.3E+05		6.3E+04	
				1.0E-01	I					1.4E+09			1	0.1	~Dibutyl Phthalate	84-74-2						7.8E+03	3.3E+04	6.3E+03	
				8.0E-01	I					1.4E+09			1	0.1	~Diethyl Phthalate	84-66-2						6.3E+04	2.6E+05	5.1E+04	

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRVT; A = ATSDR; C = Cal EPA; X = APPENDIX PPRVT SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)																										
Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³) ⁻¹	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³)	k _e (y)	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)			
				1.0E-01	I			V		1.4E+09	2.1E+04	1		**Dimethylterephthalate	120-61-6								7.8E+03		7.8E+03	
				1.0E-02	P					1.4E+09		1	0.1	**Octyl Phthalate, di-N-	117-84-0								7.8E+02	3.3E+03	6.3E+02	
				1.0E+00	H					1.4E+09		1	0.1	**Phthalic Acid, P-	100-21-0								7.8E+04	3.3E+05	6.3E+04	
				2.0E+00	I	2.0E-02	C			1.4E+09		1	0.1	**Phthalic Anhydride	85-44-9								1.6E+05	6.6E+05	2.8E+07	1.3E+05
				7.0E-02	I					1.4E+09		1	0.1	Picloram	1918-02-1								5.5E+03	2.3E+04	4.4E+03	
				1.0E-04	X					1.4E+09		1	0.1	Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3								7.8E+00	3.3E+01	6.3E+00	
				9.0E-04	X					1.4E+09		1	0.1	Picric Acid (2,4,6-Trinitrophenol)	88-89-1								7.0E+01	3.0E+02	5.7E+01	
3.0E+01	C	8.6E-03	C	7.0E-06	H					1.4E+09		1	0.1	Pirimiphos, Methyl	29232-93-7	2.3E+00	8.2E+00	4.4E+04	1.8E+00				7.8E+02	3.3E+03	6.3E+02	
				7.0E-06	H					1.4E+09		1	0.1	Polybrominated Biphenyls	59536-65-1								5.5E-01	2.3E+00	4.4E-01	
														Polychlorinated Biphenyls (PCBs)												
7.0E-02	S	2.0E-05	S	7.0E-05	I			V		1.4E+09	7.1E+05	1	0.14	**Aroclor 1016	12674-11-2	9.9E+02	2.5E+03	1.0E+04	6.7E+02				5.5E+00	1.6E+01	4.1E+00	
2.0E+00	S	5.7E-04	S					V		1.4E+09	2.0E+05	1	0.14	**Aroclor 1221	11104-28-2	3.5E+01	8.8E+01	1.0E+02	2.0E+01							
2.0E+00	S	5.7E-04	S					V		1.4E+09	1.1E+05	1	0.14	**Aroclor 1232	11141-16-5	3.5E+01	8.8E+01	5.5E+01	1.7E+01							
2.0E+00	S	5.7E-04	S					V		1.4E+09	5.9E+05	1	0.14	**Aroclor 1242	53469-21-9	3.5E+01	8.8E+01	2.9E+02	2.3E+01							
2.0E+00	S	5.7E-04	S					V		1.4E+09	6.3E+05	1	0.14	**Aroclor 1248	12672-29-6	3.5E+01	8.8E+01	3.1E+02	2.3E+01							
2.0E+00	S	5.7E-04	S	2.0E-05	I			V		1.4E+09	8.4E+05	1	0.14	**Aroclor 1254	11097-69-1	3.5E+01	8.8E+01	4.1E+02	2.4E+01				1.6E+00	4.7E+00	1.2E+00	
2.0E+00	S	5.7E-04	S					V		1.4E+09	1.3E+06	1	0.14	**Aroclor 1260	11096-82-5	3.5E+01	8.8E+01	6.5E+02	2.4E+01							
3.9E+00	E	1.1E-03	E	6.0E-04	X			V		1.4E+09	9.6E+05	1	0.14	**Aroclor 5460	11126-42-4								4.7E+01	1.4E+02	3.5E+01	
				2.3E-05	E	1.3E-03	E	V		1.4E+09	2.4E+06	1	0.14	**Heptachlorobiphenyl, 2,3,3',4,4',5,5'-(PCB 189)	39635-31-9	1.8E+01	4.5E+01	6.0E+02	1.3E+01				1.8E+00	5.5E+00	3.4E+03	1.4E+00
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	1.6E+06	1	0.14	**Hexachlorobiphenyl, 2,3',4,4',5,5'-(PCB 167)	52663-72-6	1.8E+01	4.5E+01	3.9E+02	1.2E+01				1.8E+00	5.5E+00	2.2E+03	1.4E+00
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	1.0E+06	1	0.14	**Hexachlorobiphenyl, 2,3,3',4,4',5'-(PCB 157)	69782-90-7	1.8E+01	4.5E+01	2.6E+02	1.2E+01				1.8E+00	5.5E+00	1.4E+03	1.4E+00
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	1.1E+06	1	0.14	**Hexachlorobiphenyl, 2,3,3',4,4',5'-(PCB 156)	38380-08-4	1.8E+01	4.5E+01	2.7E+02	1.2E+01				1.8E+00	5.5E+00	1.5E+03	1.4E+00
3.9E+03	E	1.1E+00	E	2.3E-08	E	1.3E-06	E	V		1.4E+09	1.6E+06	1	0.14	**Hexachlorobiphenyl, 3,3',4,4',5,5'-(PCB 169)	32774-16-6	1.8E+02	4.5E+02	3.9E+01	1.2E+02				1.8E+03	5.5E+03	2.2E+00	1.4E+03
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	7.3E+05	1	0.14	**Pentachlorobiphenyl, 2',3,4,4',5'-(PCB 123)	65510-44-3	1.8E+01	4.5E+01	1.8E+02	1.2E+01				1.8E+00	5.5E+00	1.0E+03	1.4E+00
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	5.9E+05	1	0.14	**Pentachlorobiphenyl, 2,3',4,4',5'-(PCB 118)	31508-00-6	1.8E+01	4.5E+01	1.5E+02	1.2E+01				1.8E+00	5.5E+00	8.2E+02	1.4E+00
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	6.0E+05	1	0.14	**Pentachlorobiphenyl, 2,3,3',4,4'-(PCB 105)	32598-14-4	1.8E+01	4.5E+01	1.5E+02	1.2E+01				1.8E+00	5.5E+00	8.4E+02	1.4E+00
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	1.1E+06	1	0.14	**Pentachlorobiphenyl, 2,3,4,4',5'-(PCB 114)	74472-37-0	1.8E+01	4.5E+01	2.6E+02	1.2E+01				1.8E+00	5.5E+00	1.5E+03	1.4E+00
1.3E+04	E	3.8E+00	E	7.0E-09	E	4.0E-07	E	V		1.4E+09	7.3E+05	1	0.14	**Pentachlorobiphenyl, 3,3',4,4',5'-(PCB 126)	57465-28-8	5.3E-03	1.4E-02	5.4E-02	3.6E-03				5.5E-04	1.6E-03	3.0E-01	4.1E-04
2.0E+00	I	5.7E-04	I					V		1.4E+09	5.3E+05	1	0.14	**Polychlorinated Biphenyls (high risk)	1336-36-3	3.5E+01	8.8E+01	2.6E+02	2.3E+01							
4.0E-01	I	1.0E-04	I					V		1.4E+09		1	0.14	**Polychlorinated Biphenyls (low risk)	1336-36-3											
7.0E-02	I	2.0E-05	I					V		1.4E+09		1	0.14	**Polychlorinated Biphenyls (lowest risk)	1336-36-3											
1.3E+01	E	3.8E-03	E	7.0E-06	E	4.0E-04	E	V		1.4E+09		1	0.14	**Tetrachlorobiphenyl, 3,3',4,4'-(PCB 77)	32598-13-3	5.3E+00	1.4E+01	1.0E+05	3.8E+00				5.5E-01	1.6E+00	5.7E+05	4.1E-01
3.9E+01	E	1.1E-02	E	2.3E-06	E	1.3E-04	E	V		1.4E+09	5.1E+05	1	0.14	**Tetrachlorobiphenyl, 3,4,4',5'-(PCB 81)	70362-50-4	1.8E+00	4.5E+00	1.3E+01	1.2E+00				1.8E-01	5.5E-01	7.1E+01	1.4E-01
				6.0E-04	I			V		1.4E+09		1	0.1	Polymeric Methylene Diethylene Dithiocyanate (PMDI)	9916-87-9								8.5E+05		8.5E+05	
														Polynuclear Aromatic Hydrocarbons (PAHs)												
				6.0E-02	I			V		1.4E+09	1.4E+05	1	0.13	**Acenaphthene	81-32-9								4.7E+03	1.5E+04	3.6E+03	
				3.0E-01	I			V		1.4E+09	5.2E+05	1	0.13	**Anthracene	120-12-7								2.3E+04	7.6E+04	1.8E+04	
7.3E-01	E	1.1E-04	C					V	M	1.4E+09	4.4E+06	1	0.13	**Benz[a]anthracene	56-55-3	2.1E+01	6.3E+01	4.1E+03	1.6E+01							
1.2E+00	C	1.1E-04	C							1.4E+09		1	0.13	**Benzo[<i>j</i>]fluoranthene	205-82-3	5.8E+01	1.6E+02	3.5E+06	4.2E+01							
7.3E+00	I	1.1E-03	C						M	1.4E+09		1	0.13	**Benzo[<i>a</i>]pyrene	50-32-8	2.1E+00	6.3E+00	1.3E+05	1.6E+00							
7.3E-01	E	1.1E-04	C						M	1.4E+09		1	0.13	**Benzo[<i>b</i>]fluoranthene	205-99-2	2.1E+01	6.3E+01	1.3E+06	1.6E+01							
7.3E-02	E	1.1E-04	C						M	1.4E+09		1	0.13	**Benzo[<i>k</i>]fluoranthene	207-08-9	2.1E+02	6.3E+02	1.3E+06	1.6E+02							
				8.0E-02	I			V		1.4E+09	8.0E+04	1	0.13	**Chloronaphthalene, Beta-	91-58-7								6.3E+03	2.0E+04	4.8E+03	
7.3E-03	E	1.1E-05	C						M	1.4E+09		1	0.13	**Chrysene	218-01-9	2.1E+03	6.3E+03	1.3E+07	1.6E+03							
7.3E+00	E	1.2E-03	C						M	1.4E+09		1	0.13	**Dibenz[<i>a,h</i>]anthracene	53-70-3	2.1E+00	6.3E+00	1.1E+05	1.6E+00							
1.2E+01	C	1.1E-03	C							1.4E+09		1	0.13	**Dibenzo[<i>a,e</i>]pyrene	192-65-4	5.8E+00	1.6E+01	3.5E+05	4.2E+00							
2.5E+02	C	7.1E-02	C						M	1.4E+09		1	0.13	**Dimethylbenz[<i>a</i>]anthracene, 7,12-	57-97-6	6.1E-02	1.8E-01	1.9E+03	4.6E-02							
				4.0E-02	I			V		1.4E+09		1	0.13	**Fluoranthene	206-44-0								3.1E+03	1.0E+04	2.4E+03	
				4.0E-02	I			V		1.4E+09	2.8E+05	1														

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child 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	o m u t a g e n	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)		
				4.0E-03	I				1.4E+09				1	0.1	Prometryn	7287-19-6					3.1E+02	1.3E+03		2.5E+02	
				1.3E-02	I				1.4E+09				1	0.1	Propachlor	1918-16-7					1.0E+03	4.3E+03		8.2E+02	
				5.0E-03	I				1.4E+09				1	0.1	Propanil	709-98-8					3.9E+02	1.6E+03		3.2E+02	
				2.0E-02	I				1.4E+09				1	0.1	Propargite	2312-35-8					1.6E+03	6.6E+03		1.3E+03	
				2.0E-03	I			V	1.1E+05	6.3E+04			1		Propargyl Alcohol	107-19-7					1.6E+02			1.6E+02	
				2.0E-02	I				1.4E+09				1	0.1	Propazine	139-40-2					1.6E+03	6.6E+03		1.3E+03	
				2.0E-02	I				1.4E+09				1	0.1	Propham	122-42-9					1.6E+03	6.6E+03		1.3E+03	
				1.3E-02	I				1.4E+09				1	0.1	Propiconazole	60207-90-1					1.0E+03	4.3E+03		8.2E+02	
				8.0E-03	I	V			3.3E+04	1.4E+09	8.9E+03		1		Propionaldehyde	123-38-6						7.5E+01		7.5E+01	
				1.0E-01	X	1.0E+00	X	V	2.6E+02	1.4E+09	7.0E+03		1		Propyl benzene	103-65-1					7.8E+03		7.3E+03	3.8E+03	
								C	3.5E+02	1.4E+09	7.0E+02		1		Propylene	115-07-1						2.2E+03		2.2E+03	
				2.0E+01	P				1.4E+09				1	0.1	Propylene Glycol	57-55-6					1.6E+06	6.6E+06		1.3E+06	
								A	2.7E-04				1	0.1	Propylene Glycol Dinitrate	6423-43-4							3.9E+05	3.9E+05	
				7.0E-01	H	2.0E+00	I	V	1.1E+05	1.4E+09	7.8E+04		1		Propylene Glycol Monomethyl Ether	107-98-2					5.5E+04	1.6E+05		4.1E+04	
2.4E-01	I	3.7E-06	I			3.0E-02	I	V	7.8E+04	1.4E+09	1.0E+04		1		Propylene Oxide	75-56-9	2.9E+02		7.8E+02	2.1E+02			3.2E+02	3.2E+02	
				7.5E-02	I				1.4E+09				1	0.1	Propylamide	23950-58-5					5.9E+03	2.5E+04		4.7E+03	
				1.0E-03	I			V	5.3E+05	1.4E+09	5.5E+04		1		Pyridine	110-86-1					7.8E+01			7.8E+01	
3.0E+00	I			5.0E-04	I				1.4E+09				1	0.1	Quinalphos	13593-03-8				1.8E+01			3.2E+01		
									1.4E+09				1	0.1	Quinoline	91-22-5	2.3E+01	8.2E+01						3.0E+03	3.0E+03
				9.0E-03	I				1.4E+09				1	0.1	Quizalofop-ethyl	76578-14-8					7.0E+02			5.7E+02	
								A	3.0E-02				1		Refractory Ceramic Fibers	NA							4.3E+07	4.3E+07	
				3.0E-02	I				1.4E+09				1	0.1	Resmethrin	10453-86-8					2.3E+03	9.9E+03		1.9E+03	
				5.0E-02	H			V	1.4E+09	4.7E+05			1		Ronnel	299-84-3					3.9E+03			3.9E+03	
2.2E-01	C	6.3E-05	C						1.4E+09				1	0.1	Rotenone	83-79-4								2.5E+02	
				4.0E-03	I				1.4E+09				1	0.1	Safrole	94-59-7	7.0E+01	2.7E+02	2.2E+06	5.5E+01					3.1E+02
				5.0E-03	I				1.4E+09				1		Selenious Acid	7783-00-8					3.9E+02			3.9E+02	
				5.0E-03	I	2.0E-02	C		1.4E+09				1		Selenium	7782-49-2					3.9E+02		2.8E+07	3.9E+02	
				5.0E-03	C	2.0E-02	C		1.4E+09				1		Selenium Sulfide	7446-34-6					3.9E+02		2.8E+07	3.9E+02	
				9.0E-02	I				1.4E+09				1	0.1	Sethoxydim	74051-80-2					7.0E+03	3.0E+04		5.7E+03	
1.2E-01	H					3.0E-03	C		1.4E+09				1		Silica (crystalline, respirable)	7631-86-9							4.3E+06	4.3E+06	
				5.0E-03	I				1.4E+09		0.04				Silver	7440-22-4					3.9E+02			3.9E+02	
				5.0E-03	I				1.4E+09				1	0.1	Simazine	122-34-9	5.8E+02	2.1E+03		4.5E+02		3.9E+02	1.6E+03		3.2E+02
5.0E-01	C	1.5E-01	C						1.4E+09				1	0.1	Sodium Acetate	61476-59-9					1.0E+03	4.3E+03		8.2E+02	
				4.0E-03	I				1.4E+09				1		Sodium Azide	26628-22-8					3.1E+02			3.1E+02	
				2.0E-02	C	2.0E-04	C	M	1.4E+09				0.025		Sodium Dichromate	19588-01-9	3.1E+01		9.2E+02	3.0E+01		1.6E+03		2.8E+05	1.6E+03
2.7E-01	H			3.0E-02	I				1.4E+09				1	0.1	Sodium Diethyldithiocarbamate	148-18-5	2.6E+02	9.2E+02		2.0E+02				1.8E+07	1.9E+03
				5.0E-02	A	1.3E-02	C		1.4E+09				1		Sodium Fluoride	7781-49-4					3.9E+03			3.9E+03	
				2.0E-05	I				1.4E+09				1	0.1	Sodium Fluoroacetate	62-74-8					1.6E+00	6.6E+00		1.3E+00	
				1.0E-03	H				1.4E+09				1		Sodium Metavanadate	13718-26-8					7.8E+01			7.8E+01	
				8.0E-04	P				1.4E+09				1		Sodium Tungstate	13472-45-2					6.3E+01			6.3E+01	
				8.0E-04	P				1.4E+09				1		Sodium Tungstate Dihydrate	10213-10-2					6.3E+01			6.3E+01	
2.4E-02	H			3.0E-02	I				1.4E+09				1	0.1	Stirofos (Tetrachlorovinphos)	961-11-5	2.9E+03	1.0E+04		2.3E+03				9.9E+03	1.9E+03
5.0E-01	C	1.5E-01	C						1.4E+09				0.025		Strontium Chromate	7789-06-2	3.1E+01		9.2E+02	3.0E+01		1.6E+03		2.8E+05	1.6E+03
				6.0E-01	I				1.4E+09				1		Strontium, Stable	7440-24-6					4.7E+04			4.7E+04	
				3.0E-04	I				1.4E+09				1	0.1	Strychnine	57-24-9					2.3E+01	9.9E+01		1.9E+01	
				2.0E-01	I	1.0E+00	I	V	8.7E+02	1.4E+09	9.4E+03		1		Styrene	100-42-5					1.6E+04		9.7E+03	6.0E+03	
				3.0E-03	P				1.4E+09				1	0.1	Styrene-Acrylonitrile (SAN) Trimer	NA					2.3E+02	9.9E+02		1.9E+02	
				1.0E-03	P	2.0E-03	X		1.4E+09				1	0.1	Sulfolane	126-33-0					7.8E+01	3.3E+02	2.8E+06	6.3E+01	
				8.0E-04	P				1.4E+09				1	0.1	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9					6.3E+01	2.6E+02		5.1E+01	
								C	1.0E-03				1		Sulfur Trioxide	7446-11-9							1.4E+06	1.4E+06	
2.5E-02	I	7.1E-06	I			1.0E-03	C		1.4E+09				1		Sulfuric Acid	7664-93-9							1.4E+06	1.4E+06	
				5.0E-02	H				1.4E+09				1	0.1	Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl ester	140-57-8	2.8E+03	9.9E+03	5.4E+07	2.2E+03				3.9E+03	1.6E+04
				3.0E-02	H				1.4E+09				1	0.1	TCMTB	21564-17-0					2.3E+03	9.9E+03		1.9E+03	
				7.0E-02	I				1.4E+09				1	0.1	Tebuthiuron	34014-18-1					5.5E+03	2.3E+04		4.4E+03	
				2.0E-02	H				1.4E+09				1	0.1	Temephos	3383-96-8					1.6E+03	6.6E+03		1.3E+03	
				1.3E-02	I				1.4E+09				1	0.1	Terbacil	5902-51-2					1.0E+03	4.3E+03		8.2E+02	
				2.5E-05	H			V	3.1E+01	1.4E+09	2.6E+05		1		Terbufos	13071-79-9					2.0E+00			2.0E+00	
				1.0E-03	I				1.4E+09				1	0.1	Terbutryn	886-50-0					7.8E+01	3.3E+02		6.3E+01	
				1.0E-04	I																				

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRVT; A = ATSDR; C = Cal EPA; X = APPENDIX PPRVT SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)														Toxicity and Chemical-specific Information				Contaminant				Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg- day)	k e y	RfC _o (mg/m ³) ⁻¹	k e y	o v e r l o a d	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)					
2.6E-02	I	7.4E-06	I	3.0E-02	I	3.0E-02	I	V		6.8E+02	1.4E+09	5.7E+03	1		Tetrachlorobenzene, 1,2,4,5-	95-94-3					2.3E+01			2.3E+01					
2.0E-01	I	5.8E-05	C	2.0E-02	I		V			1.9E+03	1.4E+09	1.5E+04	1		Tetrachloroethane, 1,1,1,2-	630-20-6	2.7E+03		2.2E+02	2.0E+02	2.3E+03			2.3E+03					
2.1E-03	I	2.6E-07	I	6.0E-03	I	4.0E-02	I	V		1.7E+02	1.4E+09	2.4E+03	1		Tetrachloroethylene	127-18-4	3.3E+04		2.5E+03	2.4E+03	4.7E+02		9.8E+01	8.1E+01					
2.0E+01	H			3.0E-02	I					1.4E+09			0.1	Tetrachlorophenol, 2,3,4,6-	58-90-2					2.3E+03	9.9E+03			1.9E+03					
				5.0E-04	I					1.4E+09			0.1	Tetraethyl Dithiopyrophosphate	3689-24-5						3.9E+01	1.6E+02			3.2E+01				
						8.0E+01	I	V		2.1E+03	1.4E+09	1.2E+03	1		Tetrafluoroethane, 1,1,1,2-	811-97-2							1.0E+05		1.0E+05				
				2.0E-03	P					1.4E+09			0.0007	Tetryl (Trinitrophenylmethylnitramine)	479-45-8						1.6E+02	1.0E+05			1.6E+02				
				2.0E-05	S					1.4E+09			1	Thallic Oxide	1314-32-5						1.6E+00				1.6E+00				
				1.0E-05	X					1.4E+09			1	Thallium (I) Nitrate	10102-45-1						7.8E-01				7.8E-01				
				1.0E-05	X					1.4E+09			1	Thallium (Soluble Salts)	7440-28-0						7.8E-01				7.8E-01				
				1.0E-05	X		V			1.4E+09			1	Thallium Acetate	563-68-8						7.8E-01				7.8E-01				
				2.0E-05	X		V			1.4E+09			1	Thallium Carbonate	6533-73-9						1.6E+00				1.6E+00				
				1.0E-05	X					1.4E+09			1	Thallium Chloride	7791-12-0						7.8E-01				7.8E-01				
				1.0E-05	S					1.4E+09			1	Thallium Selenite	12039-52-0						7.8E-01				7.8E-01				
				2.0E-05	X					1.4E+09			1	Thallium Sulfate	7446-18-6						1.6E+00				1.6E+00				
				1.3E-02	I					1.4E+09			0.1	Thiensusulfuron-methyl	79277-27-3						1.0E+03	4.3E+03			8.2E+02				
				1.0E-02	I					1.4E+09			0.1	Thiobencarb	28249-77-6						7.8E+02	3.3E+03			6.3E+02				
				7.0E-02	X					1.4E+09			0.0075	Thiodiglycol	111-48-8						5.5E+03	3.1E+05			5.4E+03				
				3.0E-04	H					1.4E+09			0.1	Thiofanox	39196-18-4						2.3E+01	9.9E+01			1.9E+01				
				8.0E-02	I					1.4E+09			0.1	Thiophanate, Methyl	23564-05-8						6.3E+03	2.6E+04			5.1E+03				
				5.0E-03	I					1.4E+09			0.1	Thiram	137-26-8						3.9E+02	1.6E+03			3.2E+02				
				6.0E-01	H					1.4E+09			1	Tin	7440-31-5						4.7E+04				4.7E+04				
				1.0E-04	A	V				1.4E+09			1	Titanium Tetrachloride	7550-45-0								1.4E+05			1.4E+05			
				8.0E-02	I	5.0E+00	I	V		8.2E+02	1.4E+09	4.3E+03	1	Toluene	108-88-3						6.3E+03		2.2E+04		4.9E+03				
				1.1E-05	C					1.4E+09		7.6E+05	1	Toluene-2,4-diisocyanate	584-84-9			1.9E+04	1.9E+04				6.4E+00		6.4E+00				
1.8E-01	X			2.0E-04	X					1.4E+09			0.1	Toluene-2,5-diamine	95-70-5	3.9E+02	1.4E+03		3.0E+02		1.6E+01	6.6E+01			1.3E+01				
				1.1E-05	C					1.4E+09		6.3E+05	1	Toluene-2,6-diisocyanate	91-08-7			1.6E+04	1.6E+04				5.3E+00		5.3E+00				
1.6E-02	P	5.1E-05	C			8.0E-06	C	V		1.7E+03	1.4E+09		0.1	Toluidine, o- (Methylaniline, 2-)	95-53-4	4.3E+03	1.5E+04	7.5E+06	3.4E+03										
3.0E-02	P			4.0E-03	X					1.4E+09			0.1	Toluidine, p-	106-49-0	2.3E+03	8.2E+03		1.8E+03		3.1E+02	1.3E+03			2.5E+02				
				3.0E+00	P		V			3.4E-01	1.4E+09	1.1E+03	1	Total Petroleum Hydrocarbons (Aliphatic High)	NA						2.3E+05				2.3E+05				
				6.0E-01	P	V				1.4E+02	1.4E+09	8.3E+02	1	Total Petroleum Hydrocarbons (Aliphatic Low)	NA								5.2E+02		5.2E+02				
				1.0E-02	X	1.0E-01	P	V		6.9E+00	1.4E+09	1.0E+03	1	Total Petroleum Hydrocarbons (Aliphatic Medium)	NA						7.8E+02		1.1E+02		9.6E+01				
				4.0E-02	P					1.4E+09			0.1	Total Petroleum Hydrocarbons (Aromatic High)	NA						3.1E+03	1.3E+04			2.5E+03				
				4.0E-03	P	3.0E-02	P	V		1.8E+03	1.4E+09	3.5E+03	1	Total Petroleum Hydrocarbons (Aromatic Low)	NA						3.1E+02		1.1E+02		8.2E+01				
1.1E+00	I	3.2E-04	I	4.0E-03	P	3.0E-03	P	V		1.4E+09	5.2E+04		1	Total Petroleum Hydrocarbons (Aromatic Medium)	NA						3.1E+02		1.6E+02		1.1E+02				
				7.5E-03	I					1.4E+09			0.1	Toxaphene	8001-35-2	6.3E+01	2.2E+02	1.2E+06	4.9E+01						4.7E+02				
				3.0E-04	A		V			1.4E+09	3.4E+03		1	Tri-n-butyltin	688-73-3						2.3E+01				2.3E+01				
				8.0E+01	X					1.4E+09			0.1	Triacetin	102-76-1						6.3E+06	2.6E+07			5.1E+06				
				3.0E-02	I					1.4E+09			0.1	Triadimefon	43121-43-3						2.3E+03	9.9E+03			1.9E+03				
				1.3E-02	I		V			1.4E+09	3.6E+05		1	Triallate	2303-17-5						1.0E+03				1.0E+03				
				1.0E-02	I					1.4E+09			0.1	Triasulfuron	82097-50-5						7.8E+02	3.3E+03			6.3E+02				
				8.0E-03	I					1.4E+09			0.1	Tribenuron-methyl	101200-48-0						6.3E+02	2.6E+03			5.1E+02				
9.0E-03	P			5.0E-03	I		V			1.4E+09	4.8E+04		1	Tribromobenzene, 1,2,4-	615-54-3						3.9E+02				3.9E+02				
				1.0E-02	P					1.4E+09			0.1	Tributyl Phosphate	126-73-8	7.7E+03	2.7E+04		6.0E+03		7.8E+02	3.3E+03			6.3E+02				
				3.0E-04	P					1.4E+09			0.1	Tributyltin Compounds	NA						2.3E+01	9.9E+01			1.9E+01				
				3.0E-04	I					1.4E+09			0.1	Tributyltin Oxide	56-35-9						2.3E+01	9.9E+01			1.9E+01				
7.0E-02	I			3.0E+01	H	V				9.1E+02	1.4E+09	1.3E+03	1	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1						2.3E+06		4.0E+04		4.0E+04				
				2.0E-02	I					1.4E+09			0.1	Trichloroacetic Acid	76-03-9	9.9E+02	3.5E+03		7.8E+02		1.6E+03	6.6E+03			1.3E+03				
2.9E-02	H									1.4E+09			0.1	Trichloroaniline HCl, 2,4,6-	33663-50-2	2.4E+03	8.5E+03		1.9E+03										
7.0E-03	X			3.0E-05	X					1.4E+09			0.1	Trichloroaniline, 2,4,6-	634-93-5	9.9E+03	3.5E+04		7.8E+03		2.3E+00	9.9E+00			1.9E+00				
				8.0E-04	X		V			1.4E+09	3.2E+04		1	Trichlorobenzene, 1,2,3-	87-61-6						6.3E+01				6.3E+01				
2.9E-02	P			1.0E-02	I	2.0E-03	P	V		4.0E+02	1.4E+09	3.0E+04	1	Trichlorobenzene, 1,2,4-	120-82-1	2.4E+03			2.4E+03		7.8E+02		6.2E+01		5.8E+01				
				2.0E+00	I	5.0E+00	I	V		6.4E+02	1.4E+09	1.7E+03	1	Trichloroethane, 1,1,1-	71-55-6						1.6E+05		8.6E+03		8.1E+03				
5.7E-02	I	1.6E-05	I	4.0E-03	I	2.0E-04	X	V		2.2E+03	1.4E+09	7.2E+03	1	Trichloroethane, 1,1,2-	79-00-5	1.2E+03		1.3E+02	1.1E+02		3.1E+02		1.5E+00		1.5E+00				
4.6E-02	I	4.1E-06	I	5.0E-04	I	2.0E-03	I	V	M	6.9E+02	1.4E+09	2.2E+03	1	Trichloroethylene	79-01-6														

Regional Removal Management Level (RML) Resident Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Child Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg- day)	k e y	RfC _o (mg/m ³)	k e y	o m u t a g e n	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL Child THQ=1 (mg/kg)	Dermal SL Child THQ=1 (mg/kg)	Inhalation SL Child THQ=1 (mg/kg)	Noncarcinogenic SL Child THI=1 (mg/kg)	
				1.0E-01	I				1.4E+09				1	0.1	Trichlorophenol, 2,4,5-	95-95-4					7.8E+03	3.3E+04		6.3E+03
1.1E-02	I	3.1E-06	I	1.0E-03	P				1.4E+09				1	0.1	Trichlorophenol, 2,4,6-	88-06-2	6.3E+03	2.2E+04	1.2E+08	4.9E+03	7.8E+01	3.3E+02		6.3E+01
				1.0E-02	I				1.4E+09				1	0.1	Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5					7.8E+02	3.3E+03		6.3E+02
				8.0E-03	I				1.4E+09				1	0.1	Trichlorophenoxypropionic acid, -2,4,5	93-72-1					6.3E+02	2.6E+03		5.1E+02
3.0E+01	I			5.0E-03	I			V	1.3E+03	1.4E+09	1.5E+04		1		Trichloropropane, 1,1,2-	598-77-6	5.1E-01				3.9E+02			3.9E+02
				4.0E-03	I	3.0E-04	I	V	1.4E+03	1.4E+09	1.6E+04		1		Trichloropropane, 1,2,3-	96-18-4				5.1E-01	3.1E+02		4.9E+00	4.8E+00
				3.0E-03	X	3.0E-04	P	V	3.1E+02	1.4E+09	2.3E+03		1		Trichloropropene, 1,2,3-	96-19-5					2.3E+02		7.3E-01	7.3E-01
				2.0E-02	A				1.4E+09				1	0.1	Tricresyl Phosphate (TCP)	1330-78-5					1.6E+03	6.6E+03		1.3E+03
				3.0E-03	I				1.4E+09				1	0.1	Triidiphane	58138-08-2					2.3E+02	9.9E+02		1.9E+02
				7.0E-03	I	V			2.8E+04	1.4E+09	1.6E+04		1		Triethylamine	121-44-8							1.2E+02	1.2E+02
				2.0E+00	P				1.4E+09				1	0.1	Triethylene Glycol	112-27-6					1.6E+05	6.6E+05		1.3E+05
7.7E-03	I			2.0E+01	P	V			4.8E+03	1.4E+09	7.1E+02		1		Trifluoroethane, 1,1,1-	420-46-2						1.5E+04	1.5E+04	
				7.5E-03	I	V			1.4E+09	5.1E+05			1		Trifluralin	1582-09-8	9.0E+03			9.0E+03	5.9E+02			5.9E+02
2.0E-02	P			1.0E-02	P				1.4E+09				1	0.1	Trimethyl Phosphate	512-56-1	3.5E+03	1.2E+04		2.7E+03	7.8E+02	3.3E+03		6.3E+02
				5.0E-03	P	V			2.9E+02	1.4E+09	9.4E+03		1		Trimethylbenzene, 1,2,3-	526-73-8							4.9E+01	4.9E+01
				7.0E-03	P	V			2.2E+02	1.4E+09	7.9E+03		1		Trimethylbenzene, 1,2,4-	95-63-6							5.8E+01	5.8E+01
				1.0E-02	X			V	1.8E+02	1.4E+09	6.6E+03		1		Trimethylbenzene, 1,3,5-	108-67-8					7.8E+02			7.8E+02
				1.0E-02	X			V	3.0E+01	1.4E+09	1.0E+03		1		Trimethylpentene, 2,4,4-	25167-70-8					7.8E+02			7.8E+02
				3.0E-02	I				1.4E+09				1	0.019	Trinitrobenzene, 1,3,5-	99-35-4					2.3E+03	5.2E+04		2.2E+03
3.0E-02	I			5.0E-04	I				1.4E+09				1	0.032	Trinitrotoluene, 2,4,6-	118-96-7	2.3E+03	2.6E+04		2.1E+03	3.9E+01	5.2E+02		3.6E+01
				2.0E-02	P				1.4E+09				1	0.1	Triphenylphosphine Oxide	791-28-6					1.6E+03	6.6E+03		1.3E+03
				2.0E-02	A				1.4E+09				1	0.1	Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8					1.6E+03	6.6E+03		1.3E+03
				1.0E-02	X				1.4E+09				1	0.1	Tris(1-chloro-2-propyl)phosphate	13674-84-5					7.8E+02	3.3E+03		6.3E+02
2.3E+00	C	6.6E-04	C					V	4.7E+02	1.4E+09	9.0E+05		1		Tris(2,3-dibromopropyl)phosphate	126-72-7	3.0E+01		3.8E+02	2.8E+01	3.0E+01			3.0E+01
2.0E-02	P			7.0E-03	P				1.4E+09				1	0.1	Tris(2-chloroethyl)phosphate	115-96-8	3.5E+03	1.2E+04		2.7E+03	5.5E+02	2.3E+03		4.4E+02
3.2E-03	P			1.0E-01	P				1.4E+09				1	0.1	Tris(2-ethylhexyl)phosphate	78-42-2	2.2E+04	7.7E+04		1.7E+04	7.8E+03	3.3E+04		6.3E+03
				8.0E-04	P				1.4E+09				1		Tungsten	7440-33-7					6.3E+01			6.3E+01
				3.0E-03	I	4.0E-05	A		1.4E+09				1		Uranium (Soluble Salts)	NA					2.3E+02		5.7E+04	2.3E+02
1.0E+00	C	2.9E-04	C					M	1.4E+09				1	0.1	Urethane	51-79-6	1.5E+01	6.0E+01	4.8E+05	1.2E+01	7.0E+02		9.9E+03	6.6E+02
		8.3E-03	P	9.0E-03	I	7.0E-06	P		1.4E+09				0.026		Vanadium Pentoxide	1314-62-1			4.6E+04	4.6E+04	3.9E+02		1.4E+05	3.9E+02
				5.0E-03	S	1.0E-04	A		1.4E+09				0.026		Vanadium and Compounds	7440-62-2					7.8E+01			7.8E+01
				1.0E-03	I			V	1.4E+09	1.2E+05			1		Vernolate	1929-77-7					2.0E+03			1.6E+03
				2.5E-02	I				1.4E+09				1	0.1	Vindozolin	50471-44-8					7.8E+04	8.2E+03		9.1E+02
				1.0E+00	H	2.0E-01	I	V	2.8E+03	1.4E+09	4.4E+03		1		Vinyl Acetate	108-05-4					7.8E+04			9.1E+02
7.2E-01	I	4.4E-06	I	3.2E-05	H	3.0E-03	I	V	2.5E+03	1.4E+09	1.4E+03		1		Vinyl Bromide	593-60-2			1.2E+01	1.2E+01	2.3E+02		1.0E+02	4.3E+00
				3.0E-04	I	1.0E-01	I	V	3.9E+03	1.4E+09	9.6E+02		1		Vinyl Chloride	75-01-4	9.4E+00		1.6E+01	5.9E+00	2.3E+02			7.0E+01
				3.0E-04	I				1.4E+09				1	0.1	Warfarin	81-81-2					2.3E+01	9.9E+01		1.9E+01
				2.0E-01	S	1.0E-01	S	V	3.9E+02	1.4E+09	5.6E+03		1		Xylene, p-	106-42-3					1.6E+04		5.8E+02	5.6E+02
				2.0E-01	S	1.0E-01	S	V	3.9E+02	1.4E+09	5.5E+03		1		Xylene, m-	108-38-3					1.6E+04		5.7E+02	5.5E+02
				2.0E-01	S	1.0E-01	S	V	4.3E+02	1.4E+09	6.5E+03		1		Xylene, o-	95-47-6					1.6E+04		6.7E+02	6.5E+02
				2.0E-01	I	1.0E-01	I	V	2.6E+02	1.4E+09	5.7E+03		1		Xylenes	1330-20-7					1.6E+04		6.0E+02	5.8E+02
				3.0E-04	I				1.4E+09				1		Zinc Phosphide	1314-84-7					2.3E+01			2.3E+01
				3.0E-01	I				1.4E+09				1		Zinc and Compounds	7440-66-6					2.3E+04			2.3E+04
				5.0E-02	I				1.4E+09				1	0.1	Zineb	12122-67-7					3.9E+03	1.6E+04		3.2E+03
				8.0E-05	X				1.4E+09				1		Zirconium	7440-67-7					6.3E+00			6.3E+00

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1					
SFO (mg/kg-day) ¹	ke y	IUR (ug/m ³) ¹	ke y	RfD _o (mg/kg-day)	ke y	RfC _o (mg/m ³)	ke y	o I	muta- gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
2.3E+02	I	6.7E-02	I	3.0E-03	I				M	1.34	1	1	Yes	Benztidine	92-87-5	1.1E-02	5.0E-01		1.1E-02	6.0E+01	3.0E+03		5.9E+01	
				4.0E+00	I					1.87	1	1	Yes	Benzoic Acid	65-85-0				1.1E-02	8.0E+04	1.2E+06		7.5E+04	
1.3E+01	I			1.0E-01	P				V	3.9	1	1	Yes	Benzotrchloride	98-07-7	6.0E-01	6.0E-01		3.0E-01	2.0E+03	8.9E+04		2.0E+03	
1.7E-01	I	4.9E-05	C	2.0E-03	P	1.0E-03	P	V		2.3	1	1	Yes	Benzyl Alcohol	100-51-6				3.0E-01	4.0E+01	3.2E+02	2.1E+00	2.0E+00	
		2.4E-03	I	2.0E-03	I	2.0E-05	I				0.007	1	Yes	Benzyl Chloride	100-44-7	4.6E+01	3.4E+02	1.1E+01	8.9E+00	4.0E+01	3.2E+02	2.1E+00	2.0E+00	
				9.0E-03	P					4.48	1	0.9	Yes	Beryllium and compounds	7440-41-7				3.9E+02	4.0E+01	6.4E+01		2.5E+01	4.0E+00
				1.5E-02	I					6	1	0	Yes	Bifenox	42576-02-3				3.9E+02	1.8E+02	2.3E+02		1.0E+02	
														Biphenthrin	82657-04-3				3.0E+02	3.0E+02		3.0E+02		
8.0E-03	I			5.0E-01	I	4.0E-04	X	V		4.01	1	1	Yes	Biphenyl, 1,1'-	92-52-4	9.7E+02	6.5E+02		3.9E+02	1.0E+04	7.3E+03	8.3E-01	8.3E-01	
				4.0E-02	I			V		2.48	1	1	Yes	Bis(2-chloro-1-methylethyl) ether	108-60-1				3.9E+02	8.0E+02	6.5E+03		7.1E+02	
				3.0E-03	P					1.3	1	1	Yes	Bis(2-chloroethoxy)methane	111-91-1				3.9E+02	6.0E+01	3.0E+03		5.9E+01	
1.1E+00	I	3.3E-04	I						V	1.29	1	1	Yes	Bis(2-chloroethyl)ether	111-44-4	7.1E+00	2.7E+02	1.7E+00	1.4E+00					
2.2E+02	I	6.2E-02	I						V	0.57	1	1	Yes	Bis(chloromethyl)ether	542-88-1	3.5E-02	3.4E+00	9.1E-03	7.2E-03					
				5.0E-02	I					3.32	1	1	Yes	Bisphenol A	80-05-7				7.2E-03	1.0E+03	3.2E+03		7.7E+02	
				2.0E-01	I	2.0E-02	H				1	1	Yes	Boron And Borates Only	7440-42-8				7.2E-03	4.0E+03	9.1E+05		4.0E+03	
				2.0E+00	P	2.0E-02	P	V		1.16	1	1	Yes	Boron Trichloride	10294-34-5				7.2E-03	4.0E+04	9.1E+06	4.2E+01	4.2E+01	
				4.0E-02	C	1.3E-02	C	V		0.22	1	1	Yes	Boron Trifluoride	7637-07-2				7.2E-03	8.0E+02	1.8E+05	2.7E+01	2.6E+01	
7.0E-01	I			4.0E-03	I						1	1	Yes	Bromate	15541-45-4	1.1E+01	2.1E+03		1.1E+01	8.0E+01	1.8E+04		8.0E+01	1.0E+01
2.0E+00	X	6.0E-04	X						V	1.92	1	1	Yes	Bromo-2-chloroethane, 1-	107-04-0	3.9E+00	5.7E+01	9.4E-01	7.4E-01					
				8.0E-03	I	6.0E-02	I	V		2.99	1	1	Yes	Bromobenzene	108-86-1				7.4E-01	1.6E+02	5.4E+02	1.3E+02	6.2E+01	
				4.0E-02	X	V				1.41	1	1	Yes	Bromochloromethane	74-97-5				7.4E-01			8.3E+01	8.3E+01	
6.2E-02	I	3.7E-05	C	2.0E-02	I				V	2	1	1	Yes	Bromodichloromethane	75-27-4	1.3E+02	1.9E+03	1.5E+01	1.3E+01	4.0E+02	6.5E+03	6.2E+03	3.8E+02	8.0E+01(F)
7.9E-03	I	1.1E-06	I	2.0E-02	I				V	2.4	1	1	Yes	Bromoform	75-25-2	9.9E+02	1.4E+04	5.1E+02	3.3E+02	4.0E+02	6.2E+03		3.8E+02	8.0E+01(F)
				1.4E-03	I	5.0E-03	I	V		1.19	1	1	Yes	Bromomethane	74-83-9				3.3E+02	2.8E+01	1.0E+03	1.0E+01	7.5E+00	
				5.0E-03	H				V	5.21	1	0.8	Yes	Bromophos	2104-96-3				3.3E+02	1.0E+02	5.5E+01		3.5E+01	
				2.0E-02	I				V	2.8	1	0.9	Yes	Bromoxynil	1689-84-5				3.3E+02	4.0E+02	1.8E+03		3.3E+02	
				2.0E-02	I				V	5.4	1	0.8	Yes	Bromoxynil Octanoate	1689-99-2				3.3E+02	4.0E+02	2.1E+02		1.4E+02	
3.4E+00	C	3.0E-05	I			2.0E-03	I	V		1.99	1	1	Yes	Butadiene, 1,3-	106-99-0	2.3E+00	1.6E+01	1.9E+01	1.8E+00			4.2E+00	4.2E+00	
				1.0E-01	I				V	0.88	1	1	Yes	Butanol, N	71-36-3				1.8E+00	2.0E+03	1.0E+05		2.0E+03	
				2.0E+00	P	3.0E+01	P	V		0.61	1	1	Yes	Butyl alcohol, sec	78-92-2				1.8E+00	4.0E+04	3.0E+06	6.3E+04	2.4E+04	
				5.0E-02	I				V	4.15	1	1	Yes	Butylate	7008-41-5				1.8E+00	1.0E+03	8.5E+02		4.6E+02	
2.0E-04	C	5.7E-08	C						V	3.5	1	0.8	Yes	Butylated hydroxyanisole	25013-16-5	3.9E+04	2.5E+04		1.5E+04					
3.6E-03	P			3.0E-01	P					5.1	1	1	Yes	Butylated hydroxytoluene	123-37-0	2.2E+03	4.0E+02		3.4E+02	6.0E+03	1.2E+03		1.0E+03	
				5.0E-02	P				V	4.38	1	1	No	Butylbenzene, n-	104-51-8				3.4E+02	1.0E+03			1.0E+03	
				1.0E-01	X				V	4.57	1	1	No	Butylbenzene, sec-	135-98-8				3.4E+02	2.0E+03			2.0E+03	
				1.0E-01	X				V	4.11	1	1	Yes	Butylbenzene, tert-	98-06-6				3.4E+02	2.0E+03	1.1E+03		6.9E+02	
				2.0E-02	A					0.36	1	1	Yes	Cacodylic Acid	75-60-5				3.4E+02	4.0E+02	6.7E+04		4.0E+02	
		1.8E-03	I	1.0E-03	I	1.0E-05	A				0.025	1		Cadmium (Diet)	7440-43-9				3.4E+02	4.0E+02			4.0E+02	
		1.8E-03	I	5.0E-04	I	1.0E-05	A				0.05	1	Yes	Cadmium (Water)	7440-43-9				3.4E+02	1.0E+01	1.1E+02		9.2E+00	5.0E+00
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C	M			0.025	1	Yes	Calcium Chromate	13765-19-0	5.0E+00	2.3E+01		4.1E+00	4.0E+02	2.3E+03		3.4E+02	
				5.0E-01	I	2.2E-03	C			-0.19	1	1	Yes	Caprolactam	105-60-2				4.1E+00	1.0E+04	9.0E+05		9.9E+03	
1.5E-01	C	4.3E-05	C	2.0E-03	I					3.8	1	0.9	Yes	Captafol	2425-06-1	5.2E+01	1.8E+02		4.0E+01	4.0E+01	1.5E+02		3.2E+01	
2.3E-03	C	6.6E-07	C	1.3E-01	I					2.8	1	1	Yes	Captan	133-06-2	3.4E+03	3.6E+04		3.1E+03	2.6E+03	3.0E+04		2.4E+03	
				1.0E-01	I					2.36	1	1	Yes	Carbaryl	63-25-2				3.1E+03	2.0E+03	3.4E+04		1.8E+03	
				5.0E-03	I					2.32	1	1	Yes	Carbofuran	1563-66-2				3.1E+03	1.0E+02	1.4E+03		9.4E+01	4.0E+01
				1.0E-01	I	7.0E-01	I	V		1.94	1	1	Yes	Carbon Disulfide	75-15-0				3.1E+03	2.0E+03	2.0E+04	1.5E+03	8.1E+02	
7.0E-02	I	6.0E-06	I	4.0E-03	I	1.0E-01	I	V		2.83	1	1	Yes	Carbon Tetrachloride	56-23-5	1.1E+02	4.3E+02	9.4E+01	4.6E+01	8.0E+01	3.4E+02	2.1E+02	4.9E+01	5.0E+00
				1.0E-01	P	V				-1.33	1	1	Yes	Carbonyl Sulfide	463-58-1				4.6E+01			2.1E+02	2.1E+02	
				1.0E-02	I					5.57	1	0.8	Yes	Carbosulfan	55285-14-8				4.6E+01	2.0E+02	6.9E+01		5.1E+01	
				1.0E-01	I					2.14	1	1	Yes	Carboxin	5234-68-4				4.6E+01	2.0E+03	4.1E+04		1.9E+03	
				9.0E-04	I						1	1	Yes	Ceric oxide	1306-38-3				4.6E+01					
				1.0E-01	I				V	0.99	1	1	Yes	Chloral Hydrate	302-17-0				4.6E+01	2.0E+03	1.5E+05		2.0E+03	
				1.5E-02	I					1.9	1	1	Yes	Chloramben	133-90-4				4.6E+01	3.0E+02	7.4E+03		2.9E+02	
4.0E-01	H									2.22	1	1	Yes	Chloranil	118-75-2	1.9E+01	3.5E+02		1.8E+01					
3.5E-01	I	1.0E-04	I	5.0E-04	I	7.0E-04	I	V		6.16	1	0.7	Yes	Chlordane	12789-03-6									

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information												Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	ke y	IUR (ug/m ³) ⁻¹	ke y	RfD _o (mg/kg-day)	ke y	RfC _o (mg/m ³)	ke y	o I	muta- gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
1.0E+01	I	4.6E-03	C	3.0E-04	I					5.41	1	0.8	Yes	Chlordecone (Kepone)	143-50-0	7.8E-01	6.5E-01		3.5E-01	6.0E+00	5.4E+00		2.9E+00	
		7.0E-04	A							3.81	1	0.9	Yes	Chlorfenvinphos	470-90-6					1.4E+01	5.6E+01		1.1E+01	
		2.0E-02	I							2.5	1	1	Yes	Chlorimuron, Ethyl- Chlorine	90982-32-4 7782-50-5					4.0E+02	1.5E+04		3.9E+02	
		1.0E-01	I	1.5E-04	A	V				0.85	1	1	Yes	Chlorine Dioxide	10049-04-4					2.0E+03	4.6E+05	3.0E-01	3.0E-01	
		3.0E-02	I	2.0E-04	I	V					1	1	Yes	Chlorite (Sodium Salt)	7758-19-2					6.0E+02	1.4E+05	4.2E-01	4.2E-01	
		3.0E-02	I								1	1	Yes	Chloro-1,1-difluoroethane, 1-	75-68-3					6.0E+02	1.4E+05		6.0E+02	1.0E+03
		5.0E+01	I	V						2.05	1	1	Yes	Chloro-1,3-butadiene, 2-	126-99-8			1.9E+00	1.9E+00	4.0E+02	1.8E+03	4.2E+01	3.7E+01	
4.6E-01	H	3.0E-04	I	2.0E-02	H	2.0E-02	I	V		2.53	1	1	Yes	Chloro-2-methylaniline HCl, 4-	3165-93-3	1.7E+01	5.1E+04		1.7E+01	6.0E+01	5.6E+02		5.4E+01	
1.0E-01	P	7.7E-05	C	3.0E-03	X					2.27	1	1	Yes	Chloro-2-methylaniline, 4-	95-69-2	7.8E+01	6.6E+02		7.0E+01					
2.7E-01	X								V	0.09	1	1	Yes	Chloroacetaldehyde, 2-	107-20-0	2.9E+01	4.6E+03		2.9E+01					6.0E+01
				3.0E-05	I					0.22	1	1	Yes	Chloroacetic Acid	79-11-8									
										1.93	1	1	Yes	Chloroacetophenone, 2-	532-27-4									
2.0E-01	P			4.0E-03	I					1.83	1	1	Yes	Chloroaniline, p-	106-47-8	3.9E+01	5.9E+02		3.7E+01	8.0E+01	1.3E+03		7.6E+01	
				2.0E-02	I	5.0E-02	P	V		2.84	1	1	Yes	Chlorobenzene	108-90-7					4.0E+02	1.3E+03	1.0E+02	7.8E+01	1.0E+02
1.1E-01	C	3.1E-05	C	2.0E-02	I					4.74	1	0.8	Yes	Chlorobenzilate	510-15-6	7.1E+01	5.6E+01		3.1E+01	4.0E+02	3.5E+02		1.9E+02	
				3.0E-02	X					2.65	1	1	Yes	Chlorobenzoic Acid, p-	74-11-3					6.0E+02	3.4E+03		5.1E+02	
				3.0E-03	P	3.0E-01	P	V		3.6	1	1	Yes	Chlorobenzotrifluoride, 4-	98-56-6					6.0E+01	9.3E+01	6.3E+02	3.5E+01	
				4.0E-02	P					2.64	1	1	Yes	Chlorobutane, 1-	109-69-3					8.0E+02	3.1E+03		6.4E+02	
				5.0E+01	I	V				1.08	1	1	Yes	Chlorodifluoromethane	75-45-6							1.0E+05	1.0E+05	
				2.0E-02	P					0.03	1	1	Yes	Chloroethanol, 2-	107-07-3					4.0E+02	7.7E+04		4.0E+02	
3.1E-02	C	2.3E-05	I	1.0E-02	I	9.8E-02	A	V		1.97	1	1	Yes	Chloroform	67-66-3	2.5E+02	2.9E+03	2.4E+01	2.2E+01	2.0E+02	2.5E+03	2.0E+02	9.7E+01	8.0E+01(F)
				9.0E-02	I	V				0.91	1	1	Yes	Chloromethane	74-87-3							1.9E+02	1.9E+02	
2.4E+00	C	6.9E-04	C						V	0.32	1	1	Yes	Chloromethyl Methyl Ether	107-30-2	3.2E+00	3.7E+02	8.1E-01	6.5E-01	6.0E+01	6.4E+02		5.5E+01	
3.0E-01	P			3.0E-03	P	1.0E-05	X			2.24	1	1	Yes	Chloronitrobenzene, o-	88-73-3	2.6E+01	2.6E+02		2.4E+01					
6.0E-02	P			7.0E-04	P	2.0E-03	P			2.39	1	1	Yes	Chloronitrobenzene, p-	100-00-5	1.3E+02	1.0E+03		1.2E+02	1.4E+01	1.2E+02		1.3E+01	
				5.0E-03	I				V	2.15	1	1	Yes	Chlorophenol, 2-	95-57-8					1.0E+02	1.0E+03		9.1E+01	
				4.0E-04	C	V				2.09	1	1	Yes	Chloropicrin	76-05-2							8.3E-01	8.3E-01	
3.1E-03	C	8.9E-07	C	1.5E-02	I					3.05	1	0.9	Yes	Chlorothalcnil	1897-45-6	2.5E+03	1.6E+04		2.2E+03	3.0E+02	2.1E+03		2.6E+02	
				2.0E-02	I				V	3.42	1	1	Yes	Chlorotoluene, o-	95-49-8					4.0E+02	5.8E+02		2.4E+02	
				2.0E-02	X				V	3.33	1	1	Yes	Chlorotoluene, p-	106-43-4					4.0E+02	6.6E+02		2.5E+02	
2.4E+02	C	6.9E-02	C							-1.02	1	1	Yes	Chlorozotocin	54749-90-5	3.2E-02	1.0E+02		3.2E-02					
				2.0E-01	I					3.51	1	0.9	Yes	Chlorpropham	101-21-3					4.0E+03	9.8E+03		2.8E+03	
				1.0E-03	A					4.96	1	0.8	Yes	Chlorpyrifos	2921-88-2					2.0E+01	1.5E+01		8.4E+00	
				1.0E-02	H					4.31	1	0.9	Yes	Chlorpyrifos Methyl	5598-13-0					2.0E+02	2.9E+02		1.2E+02	
				5.0E-02	I					2	1	1	Yes	Chlorsulfuron	64902-72-3					1.0E+03	5.7E+04		9.9E+02	
				1.0E-02	I					4.28	1	0.9	Yes	Chlorthal-dimethyl	1861-32-1					2.0E+02	3.3E+02		1.2E+02	
				8.0E-04	H					5.8	1	0.8	Yes	Chlorthiophos	60238-56-4					1.6E+01	3.4E+00		2.8E+00	
				1.5E+00	I					0.013	1	1	Yes	Chromium(III), Insoluble Salts	16065-83-1					3.0E+04	8.9E+04		2.2E+04	
5.0E-01	J	8.4E-02	S	3.0E-03	I	1.0E-04	I	M		0.025	1	1	Yes	Chromium(VI)	18540-29-9	5.0E+00	1.2E+01		3.5E+00	6.0E+01	1.7E+02		4.4E+01	
				1.3E-02	I					3.1	1	0.9	Yes	Chromium, Total	7440-47-3					2.6E+02	2.1E+03		2.3E+02	1.0E+02
				9.0E-03	P	3.0E-04	P	6.0E-06	P		1	1	Yes	Clofentezine	74115-24-5					6.0E+00	3.4E+03		6.0E+00	
				6.2E-04	I				V	1	0			Cobalt	7440-48-4									
				4.0E-02	H				V	1	1	1	Yes	Coke Oven Emissions	8007-45-2					8.0E+02	1.8E+05		8.0E+02	1.3E+03
				5.0E-02	I	6.0E-01	C			1.96	1	1	Yes	Copper	7440-50-8					1.0E+03	1.2E+04		9.3E+02	
				5.0E-02	I	6.0E-01	C			1.95	1	1	Yes	Cresol, o-	95-48-7					1.0E+03	1.2E+04		9.3E+02	
				1.0E-01	A	6.0E-01	C			1.94	1	1	Yes	Cresol, p-	106-44-5					2.0E+03	2.5E+04		1.9E+03	
				1.0E-01	A					3.1	1	1	Yes	Cresol, p-chloro-m-	59-50-7					2.0E+03	5.2E+03		1.4E+03	
				1.0E-01	A	6.0E-01	C			1.95	1	0.9	Yes	Cresols	1319-77-3					2.0E+03	6.7E+03		1.5E+03	
1.9E+00	H			1.0E-03	P				V	0.6	1	1	Yes	Crotonaldehyde, trans-	123-73-9	4.1E+00	2.7E+02		4.0E+00	2.0E+01	1.5E+03		2.0E+01	
				1.0E-01	I	4.0E-01	I	V		3.66	1	1	Yes	Cumene	98-82-8					2.0E+03	1.9E+03	8.3E+02	4.5E+02	
2.2E-01	C	6.3E-05	C							-1.73	1	1	Yes	Cupferron	135-20-6	3.5E+01	1.3E+06		3.5E+01					
8.4E-01	H			2.0E-03	H					2.22	1	1	Yes	Cyanazine	21725-46-2	9.3E+00	1.6E+02		8.8E+00	4.0E+01	7.6E+02		3.8E+01	

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)													Toxicity and Chemical-specific Information				Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1			
SFO (mg/kg-day) ¹	k _e IUR (ug/m ³ -day) ¹	k _e RfD _o (mg/kg-day)	k _e RfC _o (mg/m ³ -day)	k _e v _o muta- gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)							
		5.0E-02	I	2.0E-01	I	V	2.13	1	1	Yes	Dichloroethylene, 1,1-								75-35-4	1.0E+03	8.5E+03	4.2E+02	2.8E+02	7.0E+00		
		2.0E-03	I		V		1.86	1	1	Yes	Dichloroethylene, 1,2-cis-				4.0E+01	3.6E+02			156-59-2	4.0E+01	3.6E+02			7.0E+01		
		2.0E-02	I		V		2.09	1	1	Yes	Dichloroethylene, 1,2-trans-				4.0E+02	3.6E+03			156-60-5	4.0E+02	3.6E+03			1.0E+02		
		3.0E-03	I				3.06	1	1	Yes	Dichlorophenol, 2,4-				6.0E+01	1.9E+02			120-83-2	6.0E+01	1.9E+02			4.6E+01		
		1.0E-02	I				2.81	1	1	Yes	Dichlorophenoxy Acetic Acid, 2,4-				2.0E+02	1.4E+03			94-75-7	2.0E+02	1.4E+03			1.7E+02		
		8.0E-03	I				3.53	1	0.9	Yes	Dichlorophenoxybutyric Acid, 4-(2,4-				1.6E+02	4.8E+02			94-82-6	1.6E+02	4.8E+02			1.2E+02		
3.6E-02	C	1.0E-05	C	9.0E-02	A	4.0E-03	I	V	1.98	1	1	Yes	Dichloropropane, 1,2-	78-87-5	2.2E+02	2.4E+03	5.6E+01	4.4E+01		1.8E+03	2.2E+04	8.3E+00	8.3E+00	5.0E+00		
		2.0E-02	P		V		2	1	1	Yes	Dichloropropane, 1,3-				4.0E+02	4.6E+03			142-28-9	4.0E+02	4.6E+03			3.7E+02		
		3.0E-03	I				0.78	1	1	Yes	Dichloropropanol, 2,3-				6.0E+01	5.0E+03			616-23-9	6.0E+01	5.0E+03			5.9E+01		
1.0E-01	I	4.0E-06	I	3.0E-02	I	2.0E-02	I	V	2.04	1	1	Yes	Dichloropropene, 1,3-	542-75-6	7.8E+01	7.8E+02	1.4E+02	4.7E+01		6.0E+02	6.6E+03	4.2E+01	3.9E+01			
2.9E-01	I	8.3E-05	C	5.0E-04	I	5.0E-04	I		1.43	1	1	Yes	Dichlorvos	62-73-7	2.7E+01	1.4E+03		2.6E+01		1.0E+01	5.6E+02			9.9E+00		
		1.0E-04	I				0	1	1	Yes	Dicrotophos				2.0E+00	1.1E+03			141-66-2	2.0E+00	1.1E+03			2.0E+00		
		8.0E-02	P	3.0E-04	X	V	3.16	1	1	Yes	Dicyclopentadiene				1.6E+03	3.5E+03	6.3E-01		77-73-6	1.6E+03	3.5E+03	6.3E-01		6.3E-01		
1.6E+01	I	4.6E-03	I	5.0E-05	I		5.4	1	0.8	Yes	Dieldrin		4.9E-01	2.7E-01					60-57-1	1.0E+00	6.1E-01			3.8E-01		
		3.0E-04	C		5.0E-03	I			1	0									NA							
		2.0E-03	P	2.0E-04	P		-1.43	1	1	Yes	Diethanolamine				4.0E+01	8.4E+04			111-42-2	4.0E+01	8.4E+04			4.0E+01		
		3.0E-02	P	1.0E-04	P		0.56	1	1	Yes	Diethylene Glycol Monobutyl Ether				6.0E+02	8.7E+04			112-34-5	6.0E+02	8.7E+04			6.0E+02		
		6.0E-02	P	3.0E-04	P		-0.54	1	1	Yes	Diethylene Glycol Monoethyl Ether				1.2E+03	7.8E+05			111-90-0	1.2E+03	7.8E+05			1.2E+03		
		1.0E-03	P		V		0.05	1	1	Yes	Diethylformamide				2.0E+01	4.3E+03			617-84-5	2.0E+01	4.3E+03			2.0E+01		
3.5E+02	C	1.0E-01	C				5.07	1	0.9	Yes	Diethylstilbestrol	2.2E-02	6.6E-03		5.1E-03				56-53-1	2.2E-02	6.6E-03			5.1E-03		
		8.0E-02	I				0.65	1	1	Yes	Difenzoquat				1.6E+03	7.3E+05			43222-48-6	1.6E+03	7.3E+05			1.6E+03		
		2.0E-02	I				3.88	1	0.9	Yes	Diflubenzuron	4.0E+02			4.0E+02	1.0E+03			35367-38-5	4.0E+02	1.0E+03			2.9E+02		
4.4E-02	C	1.3E-05	C	4.0E+01	I	V	0.75	1	1	Yes	Diffuoroethane, 1,1-	1.8E+02	2.3E+02	4.3E+01	3.0E+01				75-37-6			8.3E+04		8.3E+04		
					V		3.58	1	1	Yes	Dihydrosafrole								94-58-6							
				7.0E-01	P	V	1.52	1	1	Yes	Diisopropyl Ether								108-20-3			1.5E+03		1.5E+03		
		8.0E-02	I		V		1.03	1	1	Yes	Diisopropyl Methylphosphonate				1.6E+03	1.3E+05			1445-75-6	1.6E+03	1.3E+05			1.6E+03		
		2.0E-02	I				-0.17	1	1	Yes	Dimethipin				4.0E+02	2.4E+05			55299-64-7	4.0E+02	2.4E+05			4.0E+02		
		2.0E-04	I				0.78	1	1	Yes	Dimethoate				4.0E+00	6.4E+02			60-51-5	4.0E+00	6.4E+02			4.0E+00		
1.6E+00	P						1.81	1	1	Yes	Dimethoxybenzidine, 3,3'	4.9E+00	1.6E+02		4.7E+00				119-50-4	4.9E+00	1.6E+02			4.7E+00		
1.7E-03	P		6.0E-02	P			-0.61	1	1	Yes	Dimethyl methylphosphonate	4.6E+03	2.8E+06		4.6E+03				756-79-6	4.6E+03	2.8E+06			4.6E+03		
4.6E+00	C	1.3E-03	C				4.58	1	1	Yes	Dimethylamino azobenzene [p-]	1.7E+00	7.2E-01		5.0E-01				60-11-7	1.7E+00	7.2E-01			5.0E-01		
5.8E-01	H						2.17	1	1	Yes	Dimethylaniline HCl, 2,4-	1.3E+01	5.2E+04		1.3E+01				21435-96-4	1.3E+01	5.2E+04			1.3E+01		
2.0E-01	P		2.0E-03	X			1.68	1	1	Yes	Dimethylaniline, 2,4-	3.9E+01	7.1E+02		3.7E+01				95-68-1	3.9E+01	7.1E+02			3.7E+01		
			2.0E-03	I		V	2.31	1	1	Yes	Dimethylaniline, N,N-				4.0E+01	8.0E+02			121-69-7	4.0E+01	8.0E+02			3.8E+01		
1.1E+01	P						2.34	1	1	Yes	Dimethylbenzidine, 3,3'	7.1E-01	8.5E+00		6.5E-01				119-93-7	7.1E-01	8.5E+00			6.5E-01		
		1.0E-01	P	3.0E-02	I	V	-1.01	1	1	Yes	Dimethylformamide				2.0E+03	1.8E+06	6.3E+01		68-12-2	2.0E+03	1.8E+06	6.3E+01		6.1E+01		
		1.0E-04	X	2.0E-06	X	V	-1.19	1	1	Yes	Dimethylhydrazine, 1,1-	2.0E+00	3.5E+03	4.2E-03	2.0E+00	3.5E+03	4.2E-03		57-14-7	2.0E+00	3.5E+03	4.2E-03		4.2E-03		
5.5E+02	C	1.6E-01	C		V		-0.54	1	1	Yes	Dimethylhydrazine, 1,2-	1.4E-02	5.0E+00	3.5E-03	2.8E-03				540-73-8	1.4E-02	5.0E+00	3.5E-03	2.8E-03			
		2.0E-02	I				2.3	1	1	Yes	Dimethylphenol, 2,4-				4.0E+02	3.1E+03			105-67-9	4.0E+02	3.1E+03			3.6E+02		
		6.0E-04	I				2.36	1	1	Yes	Dimethylphenol, 2,6-				1.2E+01	8.5E+01			576-26-1	1.2E+01	8.5E+01			1.1E+01		
4.5E-02	C	1.3E-05	C	1.0E-03	I		2.23	1	1	Yes	Dimethylphenol, 3,4-	1.7E+02	6.5E+02	4.3E+01	3.3E+01				95-65-8	1.7E+02	6.5E+02	4.3E+01	3.3E+01			
		8.0E-05	X		V		2.58	1	1	Yes	Dimethylvinylchloride				2.0E+01	1.7E+02			513-37-1	2.0E+01	1.7E+02			1.8E+01		
							2.13	1	1	Yes	Dinitro-o-cresol, 4,6-				1.6E+00	2.6E+01			534-52-1	1.6E+00	2.6E+01			1.5E+00		
		2.0E-03	I				4.12	1	0.9	Yes	Dinitro-o-cyclohexyl Phenol, 4,6-				4.0E+01	5.3E+01			131-89-5	4.0E+01	5.3E+01			2.3E+01		
		1.0E-04	P				1.69	1	1	Yes	Dinitrobenzene, 1,2-				2.0E+00	5.3E+01			528-29-0	2.0E+00	5.3E+01			1.9E+00		
		1.0E-04	I				1.49	1	1	Yes	Dinitrobenzene, 1,3-				2.0E+00	7.3E+01			99-65-0	2.0E+00	7.3E+01			2.0E+00		
		1.0E-04	P				1.46	1	1	Yes	Dinitrobenzene, 1,4-				2.0E+00	7.6E+01			100-25-4	2.0E+00	7.6E+01			2.0E+00		
6.8E-01	I			2.0E-03	I		1.67	1	1	Yes	Dinitrophenol, 2,4-				4.0E+01	1.2E+03			51-28-5	4.0E+01	1.2E+03			3.9E+01		
							2.18	1	1	Yes	Dinitrotoluene Mixture, 2,4/2,6-	1.1E+01	1.5E+02		1.1E+01				NA	1.1E+01	1.5E+02			1.1E+01		
3.1E-01	C	8.9E-05	C	2.0E-03	I		1.98	1	1	Yes	Dinitrotoluene, 2,4-	2.5E+01	4.3E+02		2.4E+01				121-14-2	2.5E+01	4.3E+02			3.8E+01		
1.5E+00	P		3.0E-04	X			2.1	1	1	Yes	Dinitrotoluene, 2,6-	5.2E+00	7.4E+01		4.9E+00				606-20-2	5.2E+00	7.4E+01			5.7E+00		
			2.0E-03	S			1.84	1	1	Yes	Dinitrotoluene, 2-Amino-4,6-				4.0E+01	1.0E+03			35572-78-2	4.0E+01	1.0E+03			3.9E+01		
		2.0E-03	S				1.84	1	1	Yes	Dinitrotoluene, 4-Amino-2,6-				4.0E+01	1.0E+03			19406-51-0	4.0E+01	1.0E+03			3.9E+01		
4.5E-0																										

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1						
SFO (mg/kg-day) ¹	ky	IUR (ug/m ³) ¹	ky	RfD _o (mg/kg-day)	ky	RfC _o (mg/m ³)	ky	vo	muta-	LOGP	GIABS	FA	In	EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
															Dinoseb	88-85-7					2.0E+01	5.4E+01		1.5E+01	7.0E+00
1.0E-01	I	5.0E-06	I	3.0E-02	I	3.0E-02	I	V		-0.27	1	1	Yes		Dioxane, 1,4-	123-91-1	7.8E+01	2.3E+04	1.1E+02	4.6E+01	6.0E+02	1.9E+05	6.3E+01	5.7E+01	
6.2E+03	I	1.3E+00	I							8.21	1	0	No		Dioxins										
															~Hexachlorodibenzo-p-dioxin, Mixture	NA	1.3E-03			1.3E-03					
1.3E+05	C	3.8E+01	C	7.0E-10	I	4.0E-08	C	V		6.8	1	0.5	No		~TCDD, 2,3,7,8-	1746-01-6	6.0E-05		1.5E-05	1.2E-05	1.4E-05		8.3E-05	1.2E-05	3.0E-05
										2.17	1	1	Yes		Diphenamid	957-51-7					6.0E+02	4.2E+03		5.3E+02	
										2.4	1	1	Yes		Diphenyl Sulfone	127-63-9					1.6E+01	2.0E+02		1.5E+01	
8.0E-01	I	2.2E-04	I	2.5E-02	I					3.5	1	1	Yes		Diphenylamine	122-39-4					5.0E+02	8.4E+02		3.1E+02	
										2.94	1	1	Yes		Diphenylhydrazine, 1,2-	122-66-7	9.7E+00	3.9E+01		7.8E+00					
										-4.6	1	1	No		Diquat	85-00-7					4.4E+01			4.4E+01	2.0E+01
7.1E+00	C	1.4E-01	C							4.9	1	1	No		Direct Black 38	1937-37-7	1.1E+00			1.1E+00					
7.4E+00	C	1.4E-01	C							2.6	1	1	No		Direct Blue 6	2602-46-2	1.1E+00			1.1E+00					
6.7E+00	C	1.4E-01	C							-6.53	1	1	No		Direct Brown 95	16071-86-6	1.2E+00			1.2E+00					
										4.0E-05	I			Yes	Disulfoton	298-04-4					8.0E-01	1.3E+00		5.0E-01	
										1.0E-02	I		V	Yes	Dithiane, 1,4-	505-29-3					2.0E+02	1.6E+04		2.0E+02	
										2.0E-03	I			Yes	Diuron	330-54-1					4.0E+01	3.6E+02		3.6E+01	
										4.0E-03	I			Yes	Dodine	2439-10-3					8.0E+01	1.1E+04		8.0E+01	
										2.5E-02	I		V	Yes	EPTC	759-94-4					5.0E+02	1.5E+03		3.8E+02	
										6.0E-03	I		V	Yes	Endosulfan	115-29-7					1.2E+02	6.3E+02		1.0E+02	
										2.0E-02	I			Yes	Endothall	145-73-3					4.0E+02	8.5E+03		3.8E+02	1.0E+02
										3.0E-04	I			Yes	Endrin	72-20-8					6.0E+00	3.7E+00		2.3E+00	2.0E+00
9.9E-03	I	1.2E-06	I	6.0E-03	P	1.0E-03	I	V		0.45	1	1	Yes		Epichlorohydrin	106-89-8	7.9E+02	7.9E+04	4.7E+02	2.9E+02	1.2E+02	1.3E+04	2.1E+00	2.0E+00	
										2.0E-02	I	V		Yes	Epoxybutane, 1,2-	106-88-7							4.2E+01	4.2E+01	
										4.0E-02	P			Yes	Ethanol, 2-(2-methoxyethoxy)-	111-77-3					8.0E+02	3.9E+05		8.0E+02	
										5.0E-03	I			Yes	Ethephon	16672-87-0					1.0E+02	4.2E+04		1.0E+02	
										5.0E-04	I			Yes	Ethion	563-12-2					1.0E+01	7.7E+00		4.3E+00	
										1.0E-01	P	6.0E-02	P	V	Yes	Ethoxyethanol Acetate, 2-	111-15-9				2.0E+03	2.3E+05	1.3E+02	1.2E+02	
										9.0E-02	P	2.0E-01	I	V	Yes	Ethoxyethanol, 2-	110-30-5				1.8E+03	6.3E+05	4.2E+02	3.4E+02	
										9.0E-01	I	7.0E-02	P	V	Yes	Ethyl Acetate	141-78-6				1.8E+04	1.2E+06	1.5E+02	1.4E+02	
										5.0E-03	P	8.0E-03	P	V	Yes	Ethyl Acrylate	140-38-5				1.0E+02	3.0E+03	1.7E+01	1.4E+01	
										1.0E+01	I	V		Yes	Ethyl Chloride (Chloroethane)	75-00-3							2.1E+04	2.1E+04	
										2.0E-01	I		V	Yes	Ethyl Ether	60-29-7					4.0E+03	2.0E+05		3.9E+03	
										3.0E-01	P	V		Yes	Ethyl Methacrylate	97-63-2							6.3E+02	6.3E+02	
										1.0E-05	I			Yes	Ethyl-p-nitrophenyl Phosphonate	2104-64-5					2.0E-01	1.6E-01		8.9E-02	
1.1E-02	C	2.5E-06	C	1.0E-01	I	1.0E+00	I	V		3.15	1	1	Yes		Ethylbenzene	100-41-4	7.1E+02	1.2E+03	2.2E+02	1.5E+02	2.0E+03	3.8E+03	2.1E+03	8.1E+02	7.0E+02
										7.0E-02	P			Yes	Ethylene Cyanohydrin	109-78-4				1.4E+03	1.1E+06		1.4E+03		
										9.0E-02	P		V	No	Ethylene Diamine	107-15-3					1.8E+03			1.8E+03	
										2.0E+00	I	4.0E-01	C	Yes	Ethylene Glycol	107-21-1					4.0E+04	5.7E+07		4.0E+04	
										1.0E-01	I	1.6E+00	I	Yes	Ethylene Glycol Monobutyl Ether	111-76-2					2.0E+03	1.4E+05		2.0E+03	
3.1E-01	C	8.8E-05	C							3.0E-02	C	V		Yes	Ethylene Oxide	75-21-8	2.5E+01	5.4E+03	6.4E+00	5.1E+00			6.3E+01	6.3E+01	
4.5E-02	C	1.3E-05	C	8.0E-05	I					-0.66	1	1	Yes		Ethylene Thiourea	96-45-7	1.7E+02	1.0E+05		1.7E+02	1.6E+00	1.0E+03		1.6E+00	
6.5E+01	C	1.9E-02	C							-0.28	1	1	Yes		Ethyleneimine	151-56-4	1.2E-01	2.5E+01	3.0E-02	2.4E-02					
										3.0E+00	I			Yes	Ethylphthalyl Ethyl Glycolate	84-72-0					6.0E+04	1.5E+06		5.8E+04	
										2.5E-04	I			Yes	Fenamiphos	22224-92-6					5.0E+00	3.4E+01		4.4E+00	
										2.5E-02	I			Yes	Fenpropathrin	39515-41-8					5.0E+02	7.3E+01		6.4E+01	
										2.5E-02	I			No	Fenvalerate	51630-58-1					5.0E+02			5.0E+02	
										1.3E-02	I			Yes	Fluometuron	2164-17-2					2.6E+02	3.4E+03		2.4E+02	
										4.0E-02	C	1.3E-02	C	Yes	Fluoride	16984-48-8					8.0E+02	1.8E+05		8.0E+02	
										6.0E-02	I	1.3E-02	C	Yes	Fluorine (Soluble Fluoride)	7782-41-4					1.2E+03	2.7E+05		1.2E+03	4.0E+03
										8.0E-02	I			Yes	Fluridone	59756-60-4					1.6E+03	1.4E+04		1.4E+03	
										2.0E-02	I			Yes	Flurprimidol	56425-91-3					4.0E+02	2.4E+03		3.4E+02	
										7.0E-04	I			Yes	Flusilazole	85509-19-9					1.4E+01	5.0E+01		1.1E+01	
										6.0E-02	I			Yes	Flutolanil	66332-96-5					1.2E+03	4.5E+03		9.5E+02	
										1.0E-02	I			No	Fluvalinate	69409-94-5					2.0E+02			2.0E+02	

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)													Toxicity and Chemical-specific Information				Contaminant				Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k e y	IUR (ug/m ³) ⁻¹	k e y	RfD _o (mg/kg-day)	k e y	RfC _o (mg/m ³)	k e y	v o l u t i l e	muta- gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL Child THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)				
3.5E-03	I			1.0E-01	I					2.85	1	1	Yes	Folpet	133-07-3	2.2E+03	2.1E+04		2.0E+03	2.0E+03	2.1E+04		1.8E+03					
1.9E-01	I			2.0E-03	I					2.9	1	1	Yes	Fomesafen	72178-02-0	4.1E+01	9.1E+02		3.9E+01									
		1.3E-05	I	2.0E-01	I	9.8E-03	A	V		3.94	1	0.9	Yes	Fonofos	944-22-9					4.0E+01	6.3E+01	2.0E+01	2.4E+01					
				9.0E-01	P	3.0E-04	X	V		0.35	1	1	Yes	Formaldehyde	50-00-0			4.3E+01	4.3E+01				2.0E+01					
				3.0E+00	I					-0.54	1	1	Yes	Formic Acid	64-18-6					1.8E+04	6.4E+06	6.3E-01	6.3E-01					
										-2.4	1	1	No	Fosetyl-AL	39148-24-8					6.0E+04			6.0E+04					
				1.0E-03	X			V		4.12	1	1	Yes	Furans														
				1.0E-03	I			V		1.34	1	1	Yes	~Dibenzofuran	132-64-9					2.0E+01	1.3E+01		7.9E+00					
				9.0E-01	I	2.0E+00	I	V		0.46	1	1	Yes	~Furan	110-00-9					2.0E+01	4.8E+02		1.9E+01					
														~Tetrahydrofuran	109-99-9					1.8E+04	1.7E+06	4.2E+03	3.4E+03					
3.8E+00	H			3.0E-03	I	5.0E-02	H	V		-0.04	1	1	Yes	Furazolidone	67-45-8	2.1E+00	1.0E+03		2.0E+00									
1.5E+00	C	4.3E-04	C							0.41	1	1	Yes	Furfural	98-01-1					6.0E+01	7.1E+03	1.0E+02	3.8E+01					
										1.8	1	1	Yes	Furium	531-82-8	5.2E+00	1.9E+02		5.1E+00									
3.0E-02	I	8.6E-06	C							4.38	1	0.9	Yes	Furmecyclo	60568-05-0	2.6E+02	2.0E+02		1.1E+02									
				4.0E-04	I					-4.81	1	1	No	Glufosinate, Ammonium	77182-82-2								8.0E+00					
						8.0E-05	C			-0.33	1	1	Yes	Glutaraldehyde	111-30-8								8.0E+00					
				4.0E-04	I	1.0E-03	H	V		-0.12	1	1	Yes	Glycidyl	765-34-4					8.0E+00	1.8E+03	2.1E+00	1.7E+00					
				1.0E-01	I					-3.4	1	1	No	Glyphosate	1071-83-6					2.0E+03			2.0E+03	7.0E+02				
				1.0E-02	X			V		-1.63	1	1	Yes	Guanidine	113-00-8					2.0E+02	4.2E+05		2.0E+02					
				2.0E-02	P					-3.56	1	1	No	Guanidine Chloride	50-01-1					4.0E+02			4.0E+02					
				5.0E-05	I					4.07	1	0.9	Yes	Haloxypol, Methyl	69806-40-2					1.0E+00	3.1E+00		7.6E-01					
4.5E+00	I	1.3E-03	I	5.0E-04	I			V		6.1	1	0.8	Yes	Heptachlor	76-44-8	1.7E+00	2.3E-01	4.3E-01	1.4E-01	1.0E+01	1.5E+00		1.3E+00	4.0E-01				
9.1E+00	I	2.6E-03	I	1.3E-05	I			V		4.98	1	0.8	Yes	Heptachlor Epoxide	1024-57-3	8.6E-01	7.1E-01	2.2E-01	1.4E-01	2.6E-01	2.4E-01		1.2E-01	2.0E-01				
				2.0E-03	I			V		6.07	1	0.7	No	Hexabromobenzene	87-82-1					4.0E+01			4.0E+01					
				2.0E-04	I					1	0	No	Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BDF-153)	68631-49-2					4.0E+00			4.0E+00						
1.6E+00	I	4.6E-04	I	8.0E-04	I			V		5.73	1	0.9	No	Hexachlorobenzene	118-74-1	4.9E+00		1.2E+00	9.8E-01	1.6E+01			1.6E+01	1.0E+00				
7.8E-02	I	2.2E-05	I	1.0E-03	P			V		4.78	1	0.9	Yes	Hexachlorobutadiene	87-68-3	1.0E+02	4.4E+01	2.6E+01	1.4E+01	2.0E+01	9.5E+00		6.5E+00					
6.3E+00	I	1.8E-03	I	8.0E-03	A					3.8	1	0.9	Yes	Hexachlorocyclohexane, Alpha-	319-34-6	1.2E+00	1.8E+00		7.2E-01	1.6E+02	2.5E+02		9.7E+01					
1.8E+00	I	5.3E-04	I							3.78	1	0.9	Yes	Hexachlorocyclohexane, Beta-	319-35-7	4.3E+00	6.1E+00		2.5E+00									
1.1E+00	C	3.1E-04	C	3.0E-04	I					3.72	1	0.9	Yes	Hexachlorocyclohexane, Gamma (Lindane)	58-89-9	7.1E+00	1.0E+01		4.2E+00	6.0E+00	9.3E+00		3.6E+00	2.0E-01				
1.8E+00	I	5.1E-04	I							4.14	1	0.9	Yes	Hexachlorocyclohexane, Technical	608-73-1	4.3E+00	6.1E+00		2.5E+00									
				6.0E-03	I	2.0E-04	I	V		5.04	1	0.9	Yes	Hexachlorocyclopentadiene	77-47-4					1.2E+02	4.2E+01	4.2E-01	4.1E-01	5.0E+01				
4.0E-02	I	1.1E-05	C	7.0E-04	I	3.0E-02	I	V		4.14	1	1	Yes	Hexachloroethane	67-72-1	1.9E+02	1.7E+02	5.1E+01	3.3E+01	1.4E+01	1.4E+01	6.3E+01	6.2E+00					
				3.0E-04	I					7.54	1	0	No	Hexachlorophene	70-30-4					6.0E+00			6.0E+00					
1.1E-01	I			3.0E-03	I					0.87	1	1	Yes	Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	7.1E+01	8.6E+03		7.0E+01	6.0E+01	8.0E+03		6.0E+01					
						1.0E-05	I	V		3.2	1	1	Yes	Hexamethylene Diisocyanate, 1,6-	822-06-0							2.1E-02	2.1E-02					
				4.0E-04	P					0.28	1	1	Yes	Hexamethylphosphoramide	680-31-9					8.0E+00	2.0E+03		8.0E+00					
				7.0E-01	I	V				3.9	1	1	Yes	Hexane, N-	110-54-3								1.5E+03					
				2.0E+00	P					0.08	1	1	Yes	Hexanedioic Acid	124-04-9					4.0E+04	1.1E+07		4.0E+04					
				5.0E-03	I	3.0E-02	I	V		1.38	1	1	Yes	Hexanone, 2-	591-78-6					1.0E+02	2.8E+03	6.3E+01	3.8E+01					
				3.3E-02	I					1.85	1	1	Yes	Hexazinone	51235-04-2					6.6E+02	2.4E+04		6.4E+02					
				2.5E-02	I					5.57	1	0.8	Yes	Hexythiazox	78587-05-0					5.0E+02	1.4E+02		1.1E+02					
				3.0E-04	I					2.31	1	1	Yes	Hydramethylnon	67485-29-4					6.0E+00	5.1E+02		5.9E+00					
3.0E+00	I	4.9E-03	I			3.0E-05	P	V		-2.07	1	1	Yes	Hydrazine	302-01-2	2.6E+00	1.1E+04	1.1E-01	1.1E-01				6.3E-02	6.3E-02				
3.0E+00	I	4.9E-03	I								1	1	Yes	Hydrazine Sulfate	10034-93-2	2.6E+00	4.9E+02		2.6E+00									
						2.0E-02	I	V			1	1	Yes	Hydrogen Chloride	7647-01-0							4.2E+01	4.2E+01					
				4.0E-02	C	1.4E-02	C	V		0.23	1	1	Yes	Hydrogen Fluoride	7664-39-3					8.0E+02	1.8E+05	2.9E+01	2.8E+01					
						2.0E-03	I	V		0.23	1	1	Yes	Hydrogen Sulfide	7783-06-4							4.2E+00	4.2E+00					
6.0E-02	P			4.0E-02	P					0.59	1	1	Yes	Hydroquinone	123-31-9	1.3E+02	1.2E+04		1.3E+02	8.0E+02	7.9E+04		7.9E+02					
				1.3E-02	I					3.82	1	0.9	Yes	Imazalil	35554-44-0					2.6E+02	6.8E+02		1.9E+02					
				2.5E-01	I					1.86	1	1	Yes	Imazaquin	81335-37-7					5.0E+03	2.6E+05		4.9E+03					
				2.5E-01	I					1.49	1	1	Yes	Imazethapyr	81335-77-5					5.0E+03	7.2E+04		4.7E+03					
				1.0E-02	A					2.49	1	1	Yes	Iodine	7553-56-2					2.0E+02	4.6E+04		2.0E+02					
				4.0E-02	I					3	1	0.9	Yes	Iprodione	36734-19-7					8.0E+02	9.1E+03		7.4E+02					

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1						
SFO (mg/kg-day) ¹	ky	IUR (ug/m ³) ¹	ky	RfD _o (mg/kg-day)	ky	RfC _o (mg/m ³)	ky	vo	muta- gen	LOGP	GIABS	FA	In	EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL THQ=1 (ug/L)	Dermal SL THQ=1 (ug/L)	Inhalation SL THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)
				7.0E-01	P						1	1	Yes	Iron	7439-89-6					1.4E+04	3.2E+06			1.4E+04	
9.5E-04	I			3.0E-01	I				V	0.76	1	1	Yes	Isobutyl Alcohol	78-83-1					6.0E+03	3.6E+05			5.9E+03	
				2.0E-01	I	2.0E+00	C			1.7	1	1	Yes	Isophorone	78-59-1	8.2E+03	1.6E+05		7.8E+03	4.0E+03	8.6E+04			3.8E+03	
				1.5E-02	I				V	5.8	1	0.8	Yes	Isopropalin	33820-53-0					3.0E+02	4.6E+01			4.0E+01	
				2.0E+00	P	2.0E-01	P	V		0.05	1	1	Yes	Isopropanol	67-63-0					4.0E+04	6.5E+06	4.2E+02		4.1E+02	
				1.0E-01	I					0.27	1	1	Yes	Isopropyl Methyl Phosphonic Acid	1832-54-8					2.0E+03	3.9E+05			2.0E+03	
				5.0E-02	I					3.94	1	0.9	Yes	Isoxaben	82558-50-7					1.0E+03	2.7E+03			7.3E+02	
						3.0E-01	A	V		8	1	0	No	JP-7	NA								6.3E+02	6.3E+02	
				2.0E-03	I					4.81	1	0.9	Yes	Lactofen	77501-63-4					4.0E+01	6.7E+01			2.5E+01	
														Lead Compounds											
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C		M		0.025	1	Yes	~Lead Chromate	7758-97-6	5.0E+00	2.3E+01		4.1E+00	4.0E+02	2.3E+03			3.4E+02	
8.5E-03	C	1.2E-05	C								1	0.8	Yes	~Lead Phosphate	7446-27-7	9.2E+02	1.7E+05		9.1E+02						
8.5E-03	C	1.2E-05	C							-0.08	1	1	Yes	~Lead acetate	301-04-2	9.2E+02	9.1E+05		9.2E+02						
8.5E-03	C	1.2E-05	C								1	1	Yes	~Lead and Compounds	7439-92-1									1.5E+01	1.5E+01
				1.0E-07	I				V	-4	1	1	No	~Lead subacetate	1335-32-6	9.2E+02			9.2E+02	2.0E-03	3.8E-03			1.3E-03	
										4.15	1	0.9	Yes	~Tetraethyl Lead	78-00-2										
5.0E-06	P								V	2.56	1	1	Yes	Lewisite	541-25-3					1.0E-01	9.1E-01			9.0E-02	
2.0E-03	I									3.2	1	0.9	Yes	Linuron	330-55-2					4.0E+01	2.0E+02			3.3E+01	
2.0E-03	P										1	1	Yes	Lithium	7439-93-2					4.0E+01	9.1E+03			4.0E+01	
5.0E-04	I									3.25	1	1	Yes	MCPA	94-74-6					1.0E+01	3.0E+01			7.5E+00	
1.0E-02	I									2.79	1	0.9	Yes	MCPB	94-81-5					2.0E+02	5.5E+02			1.5E+02	
1.0E-03	I									3.13	1	1	Yes	MCPD	93-65-2					2.0E+01	7.1E+01			1.6E+01	
2.0E-02	I									2.36	1	1	Yes	Malathion	121-75-5					4.0E+02	1.1E+04			3.9E+02	
1.0E-01	I	7.0E-04	C							1.62	1	1	Yes	Maleic Anhydride	108-31-6					2.0E+03	3.8E+04			1.9E+03	
5.0E-01	I									-0.84	1	1	Yes	Maleic Hydrazide	123-33-1					1.0E+04	8.9E+06			1.0E+04	
1.0E-04	P									-0.6	1	1	Yes	Malononitrile	109-77-3					2.0E+00	9.2E+02			2.0E+00	
3.0E-02	H									1.33	1	0.9	Yes	Mancozeb	8018-01-7					6.0E+02	4.9E+03			5.4E+02	
5.0E-03	I									0.62	1	1	Yes	Maneb	12427-38-2					1.0E+02	3.6E+03			9.8E+01	
1.4E-01	I	5.0E-05	I								1	1	Yes	Manganese (Diet)	7439-96-5										
2.4E-02	S	5.0E-05	I							0.04	1	1	Yes	Manganese (Non-diet)	7439-96-5					4.8E+02	4.4E+03			4.3E+02	
9.0E-05	H									1.04	1	1	Yes	Mepfosolan	950-10-7					1.8E+00	2.5E+02			1.8E+00	
3.0E-02	I									-2.82	1	1	No	Mepiquat Chloride	24307-26-4					6.0E+02				6.0E+02	
														Mercury Compounds											
3.0E-04	I	3.0E-04	S							-0.22	0.07	1	Yes	~Mercuric Chloride (and other Mercury salts)	7487-94-7					6.0E+00	9.6E+01			5.7E+00	2.0E+00
				3.0E-04	I	V				0.62	1	1	Yes	~Mercury (elemental)	7439-97-6								6.3E-01	6.3E-01	2.0E+00
1.0E-04	I										1	1	Yes	~Methyl Mercury	22967-92-6					2.0E+00	4.6E+02			2.0E+00	
8.0E-05	I									0.71	1	1	Yes	~Phenylmercuric Acetate	62-38-4					1.6E+00	5.7E+02			1.6E+00	
3.0E-05	I								V	7.67	1	0.3	No	Merphos	150-50-5					6.0E-01				6.0E-01	
3.0E-05	I									5.7	1	0.9	Yes	Merphos Oxide	78-48-8					6.0E-01	9.9E-02			8.5E-02	
6.0E-02	I									1.65	1	1	Yes	Metalaxyl	57837-19-1					1.2E+03	6.4E+04			1.2E+03	
1.0E-04	I	3.0E-02	P	V						0.68	1	1	Yes	Methacrylonitrile	126-98-7					2.0E+00	1.3E+02	6.3E+01		1.9E+00	
5.0E-05	I									-0.8	1	1	Yes	Methamidophos	10265-92-6					1.0E+00	1.0E+03			1.0E+00	
2.0E+00	I	2.0E+01	I	V						-0.77	1	1	Yes	Methanol	67-56-1					4.0E+04	1.8E+07	4.2E+04		2.0E+04	
1.0E-03	I									2.2	1	1	Yes	Methodathion	950-37-8					2.0E+01	5.8E+02			1.9E+01	
2.5E-02	I									0.6	1	1	Yes	Methomyl	16752-77-5					5.0E+02	6.8E+04			5.0E+02	
4.9E-02	C	1.4E-05	C							1.47	1	1	Yes	Methoxy-5-nitroaniline, 2-	99-59-2	1.6E+02	5.4E+03		1.5E+02						
5.0E-03	I									5.08	1	0.8	Yes	Methoxychlor	72-43-5					1.0E+02	5.9E+01			3.7E+01	4.0E+01
8.0E-03	P	1.0E-03	P	V						0.1	1	1	Yes	Methoxyethanol Acetate, 2-	110-49-6					1.6E+02	3.5E+04	2.1E+00		2.1E+00	
5.0E-03	P	2.0E-02	I	V						-0.77	1	1	Yes	Methoxyethanol, 2-	109-86-4					1.0E+02	6.3E+04	4.2E+01		2.9E+01	
1.0E+00	X									0.18	1	1	Yes	Methyl Acetate	79-20-9					2.0E+04	2.9E+06			2.0E+04	
				2.0E-02	P	V				0.8	1	1	Yes	Methyl Acrylate	96-33-3								4.2E+01	4.2E+01	
6.0E-01	I	5.0E+00	I	V						0.29	1	1	Yes	Methyl Ethyl Ketone (2-Butanone)	78-93-3					1.2E+04	1.5E+06	1.0E+04		5.6E+03	
1.0E-03	X	1.0E-03	P	2.0E-05	X	V				-1.05	1	1	Yes	Methyl Hydrazine	60-34-4			5.6E-01	5.6E-01	2.0E+01	1.5E+04	4.2E-02		4.2E-02	
				3.0E+00	I	V				1.31	1	1	Yes	Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1								6.3E+03	6.3E+03	

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

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SFO (mg/kg-day) ¹	k _e y	IUR (ug/m ³) ¹	k _e y	RfD _o (mg/kg-day)	k _e y	RfC _o (mg/m ³)	k _e y	v	muta- gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL THQ=1 (ug/L)	Dermal SL Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)					
						1.0E-03	C	V		0.79	1	1	Yes	Methyl Isocyanate	624-83-9							2.1E+00	2.1E+00						
		1.4E+00				7.0E-01	I	V		1.38	1	1	Yes	Methyl Methacrylate	80-62-6					2.8E+04	7.7E+05	1.5E+03	1.4E+03						
		2.5E-04								2.86	1	1	Yes	Methyl Parathion	298-00-0					5.0E+00	4.1E+01		4.5E+00						
		6.0E-02	X							-0.7	1	1	Yes	Methyl Phosphonic Acid	993-13-5					1.2E+03	1.2E+06		1.2E+03						
		6.0E-03	H	4.0E-02	H	V				3.44	1	0.8	Yes	Methyl Styrene (Mixed Isomers)	25013-15-4					1.2E+02	4.3E+01	8.3E+01	2.3E+01						
9.9E-02	C	2.8E-05	C							-0.66	1	1	Yes	Methyl methanesulfonate	66-27-3	7.9E+01	4.8E+04		7.9E+01										
1.8E-03	C	2.6E-07	C			3.0E+00	I	V		0.94	1	1	Yes	Methyl tert-Butyl Ether (MTBE)	1634-04-4	4.3E+03	2.0E+05	2.2E+03	1.4E+03			6.3E+03	6.3E+03						
		3.0E-04	X							-2.06	1	1	Yes	Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2					6.0E+00	5.9E+04		6.0E+00						
9.0E-03	P			2.0E-02	X					1.87	1	1	Yes	Methyl-5-Nitroaniline, 2-	99-55-8	8.7E+02	1.4E+04		8.2E+02	4.0E+02	7.3E+03		3.8E+02						
8.3E+00	C	2.4E-03	C							-0.92	1	1	Yes	Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	9.4E-01	1.1E+03		9.4E-01										
1.3E-01	C	3.7E-05	C							1.62	1	1	Yes	Methylaniline Hydrochloride, 2-	636-21-5	6.0E+01	3.9E+05		6.0E+01										
		1.0E-02	A							-1.18	1	1	Yes	Methylarsonic acid	124-58-3					2.0E+02	3.6E+05		2.0E+02						
		2.0E-04	X							1	0	No		Methylbenzene,1,4-diamine monohydrochloride, 2-	74612-12-7					4.0E+00		4.0E+00							
1.0E-01	X			3.0E-04	X						1	0	No	Methylbenzene-1,4-diamine sulfate, 2-	615-50-9	7.8E+01			7.8E+01	6.0E+00		6.0E+00							
2.2E+01	C	6.3E-03	C						M	6.42	1	0.8	No	Methylcholanthrene, 3-	56-49-5	1.1E-01			1.1E-01										
2.0E-03	I	1.0E-08	I	6.0E-03	I	6.0E-01	I	V	M	1.25	1	1	Yes	Methylene Chloride	75-09-2	1.3E+03	3.5E+04	2.0E+04	1.1E+03	1.2E+02	3.7E+03	1.3E+03	1.1E+02	5.0E+00					
1.0E-01	P	4.3E-04	C	2.0E-03	P				M	3.91	1	0.9	Yes	Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	2.5E+01	4.3E+01		1.6E+01	4.0E+01	7.5E+01		2.6E+01						
4.6E-02	I	1.3E-05	C							4.37	1	1	Yes	Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	1.7E+02	6.7E+01		4.8E+01										
1.6E+00	C	4.6E-04	C			2.0E-02	C			1.59	1	1	Yes	Methylenbisbenzenamine, 4,4'-	101-77-9	4.9E+00	1.7E+02		4.7E+00										
				6.0E-04	I					5.22	1	0.9	Yes	Methylenediphenyl Diisocyanate	101-68-8					1.4E+03	1.7E+03		7.8E+02						
		7.0E-02	H					V		3.48	1	1	Yes	Methylstyrene, Alpha-	98-83-9					3.0E+03	2.6E+04		2.7E+03						
		1.5E-01	I							3.13	1	1	Yes	Metolachlor	51218-45-2														
		2.5E-02	I							1.7	1	1	Yes	Metribuzin	21087-64-9					5.0E+02	1.8E+04		4.9E+02						
		2.5E-01	I							2.2	1	1	Yes	Metsulfuron-methyl	74223-64-6					5.0E+03	2.4E+05		4.9E+03						
		3.0E+00	P					V		6.1	1	1	No	Mineral oils	8012-95-1					6.0E+04		6.0E+04							
1.8E+01	C	5.1E-03	C	2.0E-04	I			V		6.89	1	0.5	No	Mirex	2385-85-5	4.3E-01		1.1E-01	8.8E-02	4.0E+00		4.0E+00							
		2.0E-03	I							3.21	1	1	Yes	Molinate	7212-67-1					4.0E+01	1.2E+02		3.0E+01						
		5.0E-03	I							1	1	1	Yes	Molybdenum	7439-98-7					1.0E+02	2.3E+04		1.0E+02						
		1.0E-01	I							1	1	1	Yes	Monochloramine	10599-90-3					2.0E+03	4.6E+05		2.0E+03	4.0E+03					
		2.0E-03	P							1.66	1	1	Yes	Monomethylaniline	100-61-8					4.0E+01	7.5E+02		3.8E+01						
		2.5E-02	I							2.94	1	1	Yes	Myclobutanil	88671-89-0					5.0E+02	4.7E+03		4.5E+02						
		3.0E-04	X							4.04	1	0.9	Yes	N,N'-Diphenyl-1,4-benzenediamine	74-31-7					6.0E+00	8.9E+00		3.6E+00						
		2.0E-03	I					V		1.38	1	1	Yes	Naled	300-76-5					4.0E+01	6.8E+03		4.0E+01						
		3.0E-02	X	1.0E-01	P	V				1	0	No		Naphtha, High Flash Aromatic (HFAN)	64742-95-6					6.0E+02		2.1E+02	1.5E+02						
1.8E+00	C	0.0E+00	C							2.28	1	1	Yes	Naphthylamine, 2-	91-59-8	4.3E+00	3.6E+01		3.9E+00	2.0E+03	9.0E+03		1.6E+03						
		1.0E-01	I							3.36	1	0.9	Yes	Napropamide	15299-99-7					2.2E+02	6.8E+05		2.2E+02						
		2.6E-04	C	1.1E-02	C	1.4E-05	C			-1.38	1	1	Yes	Nickel Acetate	373-02-4					2.2E+02	1.4E+06		2.2E+02						
		2.6E-04	C	1.1E-02	C	1.4E-05	C			-2.12	1	1	Yes	Nickel Carbonate	3333-67-3					2.2E+02		2.9E-02	2.9E-02						
		2.6E-04	C	1.1E-02	C	1.4E-05	C	V			1	0	Yes	Nickel Carbonyl	13463-39-3			2.2E+00	2.2E+00	2.2E+02		2.0E+03	2.0E+02						
		2.6E-04	C	1.1E-02	C	1.4E-05	C			0.04	1	1	Yes	Nickel Hydroxide	12054-48-7					2.2E+02	2.0E+03		2.0E+02						
		2.6E-04	C	1.1E-02	C	2.0E-05	C			0.04	1	1	Yes	Nickel Oxide	1313-99-1					2.2E+02	2.0E+03		2.0E+02						
		2.4E-04	I	1.1E-02	C	1.4E-05	C			0.04	0	Yes		Nickel Refinery Dust	NA				2.2E+02	1.0E+04		2.2E+02							
		2.6E-04	C	2.0E-02	I	9.0E-05	A			0.04	1	1	Yes	Nickel Soluble Salts	7440-02-0					4.0E+02	1.8E+04		3.9E+02						
1.7E+00	C	4.8E-04	I	1.1E-02	C	1.4E-05	C			0.04	1	1	Yes	Nickel Sub sulfide	12035-72-2	4.6E+00	1.7E+02		4.5E+00	2.2E+02	1.0E+04		2.2E+02						
		2.6E-04	C	1.1E-02	C	1.4E-05	C			1	0	Yes		Nickelocene	1271-28-9					2.2E+02			2.2E+02						
		1.6E+00	I							1	1	1	Yes	Nitrate	14797-55-8					3.2E+04	7.3E+06		3.2E+04	1.0E+04					
		1.0E-01	I							1	0	Yes		Nitrate + Nitrite (as N)	NA					2.0E+03	4.6E+05		2.0E+03	1.0E+04					
		1.0E-02	X	5.0E-05	X					1.85	1	1	Yes	Nitrite	14797-65-0					2.0E+02	3.4E+03		1.9E+02	1.0E+03					
2.0E-02	P			4.0E-03	P	6.0E-03	P			1.39	1	1	Yes	Nitroaniline, 2-	88-74-4	3.9E+02	1.2E+04		3.8E+02	8.0E+01	2.8E+03		7.8E+01						
		4.0E-05	I	2.0E-03	I	9.0E-03	I	V		1.85	1	1	Yes	Nitrobenzene	98-95-3			1.4E+01	1.4E+01	4.0E+01	6.2E+02	1.9E+01	1.3E+01						
		3.0E+03	P							-4.56	1	1	No	Nitrocellulose	9004-70-0				6.0E+07			6.0E+07							
		7.0E-02	H							-0.47	1	1	Yes	Nitrofurantoin	67-20-9					1.4E+03	1.6E+06		1.4E+03						
1.3E+00	C	3.7E-04	C							0.23	1	1	Yes	Nitrofurazone	59-87-0	6.0E+00	1.7E+03		6.0E+00										

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1					
SFO (mg/kg-day) ¹	ke (ug/m ³) ¹	IUR (ug/m ³) ¹	ke (mg/kg-day)	RfD _o (mg/kg-day)	ke (mg/m ³)	ke (mg/m ³)	ke (mg/m ³)	ke (mg/m ³)	ke (mg/m ³)	ke (mg/m ³)	ke (mg/m ³)	ke (mg/m ³)	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL THQ=1 (ug/L)	Dermal SL THQ=1 (ug/L)	Inhalation SL THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)	
1.7E-02	P		1.0E-04	P									Nitroglycerin	55-63-0	4.6E+02	1.8E+04		4.5E+02	2.0E+00	8.7E+01		2.0E+00		
			1.0E-01	I									Nitroguanidine	556-88-7					2.0E+03	1.8E+06		2.0E+03		
		8.8E-06	P		5.0E-03	P	V						Nitromethane	75-52-5			6.4E+01	6.4E+01			1.0E+01	1.0E+01		
		2.7E-03	H		2.0E-02	I	V						Nitropropane, 2-	79-46-9			2.1E-01	2.1E-01			4.2E+01	4.2E+01		
2.7E+01	C	7.7E-03	C					M	0.23	1	1	Yes	Nitroso-N-ethylurea, N-	759-73-9	9.3E-02	1.5E+01		9.2E-02						
1.2E+02	C	3.4E-02	C					M	-0.03	1	1	Yes	Nitroso-N-methylurea, N-	684-93-5	2.1E-02	4.6E+00		2.1E-02						
5.4E+00	I	1.6E-03	I					V	2.63	1	1	Yes	Nitroso-di-N-butylamine, N-	924-16-3	1.4E+00	7.9E+00	3.5E-01	2.7E-01						
7.0E+00	I	2.0E-03	C						1.36	1	1	Yes	Nitroso-di-N-propylamine, N-	621-64-7	1.1E+00	3.5E+01		1.1E+00						
2.8E+00	I	8.0E-04	C						-1.28	1	1	Yes	Nitrosodiethanolamine, N-	1116-54-7	2.8E+00	8.1E+03		2.8E+00						
1.5E+02	I	4.3E-02	I					M	0.48	1	1	Yes	Nitrosodiethylamine, N-	55-18-5	1.7E-02	1.7E+00		1.7E-02						
5.1E+01	I	1.4E-02	I	8.0E-06	P	4.0E-05	X	V	M	-0.57	1	1	Yes	Nitrosodimethylamine, N-	62-75-9	4.9E-02	2.0E+01	1.4E-02	1.1E-02	1.6E-01	7.4E+01	8.3E-02	5.5E-02	
4.9E-03	I	2.6E-06	C						3.13	1	1	Yes	Nitrosodiphenylamine, N-	86-30-6	1.6E+03	5.2E+03		1.2E+03						
2.2E+01	I	6.3E-03	C					V	0.04	1	1	Yes	Nitrosomethylethylamine, N-	10595-95-6	3.5E-01	6.4E+01	8.9E-02	7.1E-02						
6.7E+00	C	1.9E-03	C						-0.44	1	1	Yes	Nitrosomorpholine [N-]	59-89-2	1.2E+00	5.3E+02		1.2E+00						
9.4E+00	C	2.7E-03	C						0.36	1	1	Yes	Nitrosopiperidine [N-]	100-75-4	8.3E-01	1.1E+02		8.2E-01						
2.1E+00	I	6.1E-04	I						-0.19	1	1	Yes	Nitrosopyrrolidine, N-	930-55-2	3.7E+00	1.0E+03		3.7E+00						
2.2E-01	P		1.0E-04	X					2.45	1	1	Yes	Nitrotoluene, m-	99-08-1					2.0E+00	1.4E+01		1.7E+00		
1.6E-02	P		9.0E-04	P				V	2.3	1	1	Yes	Nitrotoluene, o-	88-72-2	3.5E+01	2.8E+02		3.1E+01	1.8E+01	1.5E+02		1.6E+01		
			4.0E-03	P					2.37	1	1	Yes	Nitrotoluene, p-	99-99-0	4.9E+02	3.4E+03		4.3E+02	8.0E+01	6.2E+02		7.1E+01		
			3.0E-04	X	2.0E-02	P	V		5.65	1	1	No	Nonane, n-	111-84-2					6.0E+00		4.2E+01	5.3E+00		
			4.0E-02	I					2.3	1	1	Yes	Norflurazon	27314-13-2					8.0E+02	2.0E+04		7.7E+02		
			3.0E-03	I					8.71	1	0.3	No	Octabromodiphenyl Ether	32536-52-0					6.0E+01			6.0E+01		
			5.0E-02	I					0.16	1	1	Yes	Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0				1.0E+03	6.3E+05		1.0E+03			
			2.0E-03	H					-1.01	1	1	Yes	Octamethylpyrophosphoramide	152-16-9				4.0E+01	1.4E+05		4.0E+01			
			5.0E-02	I					3.73	1	0.9	Yes	Oryzalin	19044-88-3				1.0E+03	4.1E+03		8.1E+02			
			5.0E-03	I					4.8	1	0.8	Yes	Oxadiazon	19666-30-9				1.0E+02	9.0E+01		4.7E+01			
			2.5E-02	I					-0.47	1	1	Yes	Oxamyl	23135-22-0				5.0E+02	5.1E+05		5.0E+02	2.0E+02		
			3.0E-03	I					4.73	1	0.8	Yes	Oxyfluorfen	42874-03-3				6.0E+01	6.7E+01		3.2E+01			
			1.3E-02	I					3.2	1	0.9	Yes	Paclotrazol	76738-62-0				2.6E+02	1.7E+03		2.3E+02			
			4.5E-03	I					-4.5	1	1	No	Paraquat Dichloride	1910-42-5				9.0E+01			9.0E+01			
			6.0E-03	H					3.83	1	0.9	Yes	Parathion	56-98-2				1.2E+02	3.0E+02		8.6E+01			
			5.0E-02	H				V	3.83	1	1	Yes	Pebulate	1114-71-2				1.0E+03	1.3E+03		5.6E+02			
			4.0E-02	I					5.2	1	0.9	Yes	Pendimethalin	40487-42-1				8.0E+02	2.4E+02		1.8E+02			
			2.0E-03	I				V	6.84	1	0.6	No	Pentabromodiphenyl Ether	32534-81-9				4.0E+01			4.0E+01			
			1.0E-04	I					7.66	1	0.6	No	Pentabromodiphenyl ether, 2,2',4,4',5'- (BDE-99)	60348-60-9				2.0E+00			2.0E+00			
			8.0E-04	I				V	5.17	1	0.9	Yes	Pentachlorobenzene	608-93-5				1.6E+01	3.9E+00		3.2E+00			
9.0E-02	P							V	3.22	1	1	Yes	Pentachloroethane	76-01-7	8.7E+01	2.5E+02		6.5E+01						
2.6E-01	H		3.0E-03	I				V	4.64	1	0.9	Yes	Pentachloronitrobenzene	82-68-8	3.0E+01	2.0E+01		1.2E+01	6.0E+01	4.4E+01		2.6E+01		
4.0E-01	I	5.1E-06	C		5.0E-03	I			5.12	1	0.9	Yes	Pentachlorophenol	87-86-5	1.9E+01	5.2E+00		4.1E+00	1.0E+02	2.9E+01		2.3E+01	1.0E+00	
4.0E-03	X		2.0E-03	P					2.38	1	1	Yes	Pentaerythritol tetranitrate (PETN)	78-11-5	1.9E+03	4.3E+04		1.9E+03	4.0E+01	9.6E+02		3.9E+01		
			1.0E+00	P	V				3.39	1	1	Yes	Pentane, n-	109-66-0							2.1E+03	2.1E+03		
			7.0E-04	I						1	1	Yes	Perchlorates						1.4E+01	3.2E+03		1.4E+01		
			7.0E-04	I						1	1	Yes	~Ammonium Perchlorate	7790-98-9					1.4E+01	3.2E+03		1.4E+01		
			7.0E-04	I						1	1	Yes	~Lithium Perchlorate	7791-03-9					1.4E+01	3.2E+03		1.4E+01		
			7.0E-04	I						1	1	Yes	~Perchlorate and Perchlorate Salts	14797-73-0					1.4E+01	3.2E+03		1.4E+01	1.5E+01(F)	
			7.0E-04	I						1	1	Yes	~Potassium Perchlorate	7778-74-7					1.4E+01	1.6E+03		1.4E+01		
			7.0E-04	I						1	1	Yes	~Sodium Perchlorate	7601-89-0					1.4E+01	3.2E+03		1.4E+01		
			2.0E-02	P				V	2.41	1	1	Yes	Perfluorobutane Sulfonate	375-73-5				4.0E+02	8.3E+03		3.8E+02			
			5.0E-02	I					6.5	1	0.6	No	Permethrin	52645-53-1				1.0E+03			1.0E+03			
2.2E-03	C	6.3E-07	C						1.58	1	1	Yes	Phenacetin	62-44-2	3.5E+03	1.1E+05		3.4E+03						
			2.5E-01	I					3.59	1	0.9	Yes	Phenmedipham	13684-63-4				5.0E+03	1.9E+04		4.0E+03			
			3.0E-01	I	2.0E-01	C			1.46	1	1	Yes	Phenol	108-95-2				6.0E+03	1.4E+05		5.8E+03			
			4.0E-03	I					1.52	1	1	Yes	Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1				8.0E+01	3.6E+03		7.8E+01			
			5.0E-04	X					4.15	1	1	Yes	Phenothiazine	92-84-2				1.0E+01	7.6E+00		4.3E+00			

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1			
SFO (mg/kg-day) ¹	k _e IUR (ug/m ³) ¹	k _e RfD _o (mg/kg-day)	k _e RfC _o (mg/m ³)	k _e v _o muta- gen	LOGP	GIABS	FA	In	EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL THQ=1 (ug/L)	Dermal SL THQ=1 (ug/L)	Inhalation SL THQ=1 (ug/L)	Noncarcinogenic SL THI=1 (ug/L)	MCL (ug/L)		
4.7E-02	H	6.0E-03	I		-0.33	1	1	Yes		Phenylenediamine, m-	108-45-2					1.2E+02	4.8E+04		1.2E+02			
										Phenylenediamine, o-	95-54-5	1.7E+02	2.9E+04		1.6E+02							
1.9E-03	H	1.9E-01	H		-0.3	1	1	Yes		Phenylenediamine, p-	106-50-3					3.8E+03	1.4E+06		3.8E+03			
										Phenylphenol, 2-	90-43-7	4.0E+03	1.2E+04		3.0E+03							
		2.0E-04	H		3.56	1	0.9	Yes		Phorate	298-02-2					4.0E+00	1.2E+01		3.0E+00			
										Phosgene	75-44-5											
		2.0E-02	I	3.0E-04	I	V				Phosmet	732-11-6					4.0E+02	5.3E+03		3.7E+02			
		4.9E+01	P							Phosphates, Inorganic												
		4.9E+01	P							~Aluminum metaphosphate	13776-88-0					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Ammonium polyphosphate	68333-79-9					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Calcium pyrophosphate	7790-76-3					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Diammonium phosphate	7783-28-0					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Dicalcium phosphate	7757-93-9					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Dimagnesium phosphate	7782-75-4					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Dipotassium phosphate	7758-11-4					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Disodium phosphate	7558-79-4					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Monoaluminum phosphate	13530-50-2					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Monoammonium phosphate	7722-76-1					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Monocalcium phosphate	7758-23-8					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Monomagnesium phosphate	7757-86-0					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Monopotassium phosphate	7778-77-0					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Monosodium phosphate	7558-80-7					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Polyphosphoric acid	8017-16-1					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Potassium tripolyphosphate	13845-36-8					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Sodium acid pyrophosphate	7758-16-9					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Sodium aluminum phosphate (acidic)	7785-88-8					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Sodium aluminum phosphate (anhydrous)	10279-59-1					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Sodium aluminum phosphate (tetrahydrate)	10305-76-7					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Sodium hexametaphosphate	10124-56-8					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Sodium polyphosphate	68915-31-1					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Sodium trimetaphosphate	7785-84-4					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Sodium tripolyphosphate	7758-29-4					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Tetrapotassium phosphate	7320-34-5					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Tetrasodium pyrophosphate	7722-88-5					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Trialuminum sodium tetra decahydrogenoctaorthophosphate (dihydrate)	15136-87-5					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Tricalcium phosphate	7758-87-4					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Trimagnesium phosphate	7757-87-1					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Tripotassium phosphate	7778-53-2					9.7E+05	2.2E+08		9.7E+05			
		4.9E+01	P							~Trisodium phosphate	7601-54-9					9.7E+05	2.2E+08		9.7E+05			
		3.0E-04	I	3.0E-04	I	V				Phosphine	7803-51-2					6.0E+00	1.4E+03	6.3E-01	5.7E-01			
		4.9E+01	P	1.0E-02	I					Phosphoric Acid	7664-38-2					9.7E+05	2.2E+08		9.7E+05			
		2.0E-05	I		V					Phosphorus, White	7723-14-0					4.0E-01	9.1E+01		4.0E-01			
										Phthalates												
1.4E-02	I	2.4E-06	C	2.0E-02	I					~Bis(2-ethylhexyl)phthalate	117-81-7	5.6E+02			5.6E+02	4.0E+02			4.0E+02	6.0E+00		
1.9E-03	P	2.0E-01	I							~Butyl Benzyl Phthalate	85-68-7	4.1E+03	2.7E+03		1.6E+03	4.0E+03	2.9E+03		1.7E+03			
		1.0E+00	I							~Butylphthalyl Butylglycolate	85-70-1					2.0E+04	4.1E+04		1.3E+04			
		1.0E-01	I							~Dibutyl Phthalate	84-74-2					2.0E+03	1.6E+03		9.0E+02			
		8.0E-01	I							~Diethyl Phthalate	84-66-2					1.6E+04	2.0E+05		1.5E+04			
		1.0E-01	I	V						~Dimethylterephthalate	120-61-6					2.0E+03	2.7E+04		1.9E+03			
		1.0E-02	P							~Octyl Phthalate, di-N-	117-84-0					2.0E+02			2.0E+02			
		1.0E+00	H							~Phthalic Acid, P-	100-21-0					2.0E+04	3.3E+05		1.9E+04			
		2.0E+00	I	2.0E-02	C					~Phthalic Anhydride	85-44-9					4.0E+04	1.1E+06		3.9E+04			
		7.0E-02	I							Picloram	1918-02-1					1.4E+03	4.3E+04		1.4E+03	5.0E+02		
		1.0E-04	X							Picramic Acid (2-Amino-4,6-dinitrophenol)	96-91-3					2.0E+00	2.1E+02		2.0E+00			

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)														Toxicity and Chemical-specific Information				Contaminant				Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1			
SFO (mg/kg-day) ¹	k e y	IUR (ug/m ³) ¹	k e y	RfD _o (mg/kg-day)	k e y	RfC _o (mg/m ³)	k e y	muta- gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL THQ=1 (ug/L)	Dermal SL THQ=1 (ug/L)	Inhalation SL THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)						
				9.0E-04	X				1.44	1	1	Yes	Picric Acid (2,4,6-Trinitrophenol)	88-89-1					1.8E+01	1.2E+03		1.8E+01							
3.0E+01	C	8.6E-03	C	1.0E-02	I	7.0E-06	H		4.2	1	0.9	Yes	Primingophos, Methyl	29232-93-7					2.0E+02	3.1E+02		1.2E+02							
				7.0E-06	H					1	0	No	Polybrominated Biphenyls	59536-65-1	2.6E-01			2.6E-01	1.4E-01			1.4E-01							
													Polychlorinated Biphenyls (PCBs)																
7.0E-02	S	2.0E-05	S	7.0E-05	I		V		5.69	1	0	No	~Aroclor 1016	12674-11-2	1.1E+02		2.8E+01	2.2E+01	1.4E+00			1.4E+00							
2.0E+00	S	5.7E-04	S				V		4.65	1	1	Yes	~Aroclor 1221	11104-28-2	3.9E+00	1.2E+00	9.8E-01	4.7E-01											
2.0E+00	S	5.7E-04	S				V		4.4	1	1	Yes	~Aroclor 1232	11141-16-5	3.9E+00	1.2E+00	9.8E-01	4.7E-01											
2.0E+00	S	5.7E-04	S				V		6.34	1	0.7	No	~Aroclor 1242	53469-21-9	3.9E+00		9.8E-01	7.8E-01											
2.0E+00	S	5.7E-04	S				V		6.2	1	0	No	~Aroclor 1248	12672-29-6	3.9E+00		9.8E-01	7.8E-01											
2.0E+00	S	5.7E-04	S	2.0E-05	I		V		6.5	1	0.5	No	~Aroclor 1254	11097-69-1	3.9E+00		9.8E-01	7.8E-01	4.0E-01			4.0E-01							
2.0E+00	S	5.7E-04	S				V		7.55	1	0	No	~Aroclor 1260	11096-82-5	3.9E+00		9.8E-01	7.8E-01											
3.9E+00	E	1.1E-03	E	6.0E-04	X		V		6.34	1	0.7	No	~Aroclor 5460	11126-42-4					1.2E+01			1.2E+01							
				2.3E-05	E	1.3E-03	E	V	8.27	1	0	No	~Heptachlorobiphenyl, 2,3,3',4,4',5,5' (PCB 189)	39635-31-9	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01							
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.5	1	0	No	~Hexachlorobiphenyl, 2,3',4,4',5,5' (PCB 167)	52663-72-6	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01							
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.6	1	0	No	~Hexachlorobiphenyl, 2,3,3',4,4',5' (PCB 157)	69782-90-7	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01							
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.6	1	0	No	~Hexachlorobiphenyl, 2,3,3',4,4',5' (PCB 156)	38380-08-4	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01							
3.9E+03	E	1.1E+00	E	2.3E-08	E	1.3E-06	E	V	7.41	1	0.1	No	~Hexachlorobiphenyl, 3,3',4,4',5,5' (PCB 169)	32774-16-6	2.0E-03		4.9E-04	4.0E-04	4.7E-04		2.8E-03	4.0E-04							
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	6.98	1	0.4	No	~Pentachlorobiphenyl, 2',3,4,4',5' (PCB 123)	65510-44-3	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01							
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	7.12	1	0.3	No	~Pentachlorobiphenyl, 2,3',4,4',5' (PCB 118)	31508-00-6	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01							
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	6.79	1	0.5	No	~Pentachlorobiphenyl, 2,3,3',4,4' (PCB 105)	32598-14-4	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01							
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V	6.98	1	0.4	No	~Pentachlorobiphenyl, 2,3,4,4',5' (PCB 114)	74472-37-0	2.0E+00		4.9E-01	4.0E-01	4.7E-01		2.8E+00	4.0E-01							
1.3E+04	E	3.8E+00	E	7.0E-09	E	4.0E-07	E	V	6.98	1	0.4	No	~Pentachlorobiphenyl, 3,3',4,4',5' (PCB 126)	57465-28-8	6.0E-04		1.5E-04	1.2E-04	1.4E-04		8.3E-04	1.2E-04							
2.0E+00	I	5.7E-04	I				V		7.1	1	0.7	No	~Polychlorinated Biphenyls (high risk)	1336-36-3															
4.0E-01	I	1.0E-04	I				V		7.1	1	0.7	No	~Polychlorinated Biphenyls (low risk)	1336-36-3	1.9E+01		5.6E+00	4.4E+00					5.0E-01						
7.0E-02	I	2.0E-05	I				V		7.1	1	0.7	No	~Polychlorinated Biphenyls (lowest risk)	1336-36-3															
1.3E+01	E	3.8E-03	E	7.0E-06	E	4.0E-04	E	V	6.63	1	0.6	No	~Tetrachlorobiphenyl, 3,3',4,4' (PCB 77)	32598-13-3	6.0E-01			6.0E-01	1.4E-01			1.4E-01							
3.9E+01	E	1.1E-02	E	2.3E-06	E	1.3E-04	E	V	6.34	1	0.7	No	~Tetrachlorobiphenyl, 3,4,4',5' (PCB 81)	70263-50-4	2.0E-01		4.9E-02	4.0E-02	4.7E-02		2.8E-01	4.0E-02							
				6.0E-04	I				10.46	1	0	No	Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-81-9															
				6.0E-02	I		V		3.92	1	1	Yes	Polynuclear Aromatic Hydrocarbons (PAHs)																
				3.0E-01	I		V		4.45	1	1	Yes	~Acenaphthene	83-32-9					1.2E+03	9.6E+02		5.3E+02							
7.3E-01	E	1.1E-04	C				V	M	5.76	1	1	No	~Anthracene	120-12-7					6.0E+03	2.5E+03		1.8E+03							
1.2E+00	C	1.1E-04	C				V	M	6.11	1	0.9	No	~Benzo[a]anthracene	56-55-3	3.4E+00		1.8E+00	1.2E+00											
7.3E+00	I	1.1E-03	C				V	M	6.13	1	1	No	~Benzo[j]fluoranthene	205-82-3	6.5E+00			6.5E+00					2.0E-01						
7.3E-01	E	1.1E-04	C				V	M	5.78	1	1	No	~Benzo[a]pyrene	50-32-8	3.4E-01			3.4E-01											
7.3E-02	E	1.1E-04	C				V	M	6.11	1	0.9	No	~Benzo[b]fluoranthene	205-99-2	3.4E+00			3.4E+00											
				8.0E-02	I		V		3.9	1	1	Yes	~Benzo[k]fluoranthene	207-08-9	3.4E+01			3.4E+01											
							V		5.81	1	1	No	~Chloronaphthalene, Beta-	91-58-7					1.6E+03	1.4E+03		7.5E+02							
7.3E-03	E	1.1E-05	C				V	M	5.81	1	1	No	~Chrysene	218-01-9	3.4E+02			3.4E+02											
7.3E+00	E	1.2E-03	C				V	M	6.75	1	0.6	No	~Dibenzo[a,h]anthracene	53-70-3	3.4E-01			3.4E-01											
1.2E+01	C	1.1E-03	C				V	M	7.71	1	0.3	No	~Dibenzo[a,e]pyrene	192-65-4	6.5E-01			6.5E-01											
2.5E+02	C	7.1E-02	C				V	M	5.8	1	0.9	No	~Dimethylbenz(a)anthracene, 7,12-	57-97-6	1.0E-02			1.0E-02											
				4.0E-02	I		V		5.16	1	1	No	~Fluoranthene	206-44-0					8.0E+02			8.0E+02							
				4.0E-02	I		V		4.18	1	1	Yes	~Fluorene	86-73-7					8.0E+02	4.6E+02		2.9E+02							
7.3E-01	E	1.1E-04	C				V	M	6.7	1	0.6	No	~Indeno[1,2,3-cd]pyrene	193-39-5	3.4E+00			3.4E+00											
2.9E-02	P			7.0E-02	A		V		3.87	1	1	Yes	~Methylnaphthalene, 1-	90-12-0	2.7E+02	2.0E+02		1.1E+02	1.4E+03	1.1E+03		6.2E+02							
				4.0E-03	I		V		3.86	1	1	Yes	~Methylnaphthalene, 2-	91-57-6					8.0E+01	6.5E+01		3.6E+01							
1.2E+00	C	3.4E-05	C	2.0E-02	I	3.0E-03	I	V	3.3	1	1	Yes	~Naphthalene	91-20-3			1.7E+01	1.7E+01	4.0E+02	7.0E+02	6.3E+00	6.1E+00							
				3.0E-02	I		V		4.75	1	0.9	Yes	~Nitropyrene, 4-	57835-92-4	6.5E+00	2.7E+00		1.9E+00											
							V		4.88	1	1	Yes	~Pyrene	129-00-0					6.0E+02	1.5E+02		1.2E+02							
1.5E-01	I			2.0E-02	P				-0.3297	1	1	Yes	Potassium Perfluorobutane Sulfonate	29420-49-3					4.0E+02	2.8E+05		4.0E+02							
				9.0E-03	I		V		4.1	1	0.9	Yes	Prochloraz	67747-09-5	5.2E+01	1.4E+02		3.8E+01	1.8E+02	5.1E+02		1.3E+02							
				6.0E-03	H		V		5.58	1	0.8	Yes	Profluralin	26399-36-0					1.2E+02	3.3E+01		2.6E+01							
				1.5E-02	I		V		2.99	1	1	Yes	Prometon	1610-18-0					3.0E+02	1.6E+03		2.5E+02							
				4.0E-03	I		V		3.51	1	0.9	Yes	Prometryn	7287-19-6					8.0E+01	2.3E+02		6.0E+01							

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1					
SFO (mg/kg-day) ¹	k e y	IUR (ug/m ³) ¹	k e y	RfD _o (mg/kg-day)	k e y	RfC _o (mg/m ³)	k e y	v o l a t i l e	m u t a g e n	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL THQ=1 (ug/L)	Dermal SL THQ=1 (ug/L)	Inhalation SL THQ=1 (ug/L)	Noncarcinogenic SL THI=1 (ug/L)	MCL (ug/L)
		1.3E-02	I							2.18	1	1	Yes	Propachlor	1918-16-7					2.6E+02	4.3E+03		2.5E+02	
		5.0E-03	I							3.07	1	1	Yes	Propanil	709-98-8					1.0E+02	4.4E+02		8.2E+01	
		2.0E-02	I							5	1	0.8	Yes	Propargite	2312-35-8					4.0E+02	2.7E+02		1.6E+02	
		2.0E-03	I					V		-0.38	1	1	Yes	Propargyl Alcohol	107-19-7					4.0E+01	1.2E+04		4.0E+01	
		2.0E-02	I							2.93	1	1	Yes	Propazine	139-40-2					4.0E+02	2.4E+03		3.4E+02	
		2.0E-02	I							2.6	1	1	Yes	Propham	122-42-9					4.0E+02	2.8E+03		3.5E+02	
		1.3E-02	I							3.72	1	0.9	Yes	Propiconazole	60207-90-1					2.6E+02	1.1E+03		2.1E+02	
						8.0E-03	I	V		0.59	1	1	Yes	Propionaldehyde	123-38-6							1.7E+01	1.7E+01	
		1.0E-01	X			1.0E+00	X	V		3.69	1	1	Yes	Propyl benzene	103-65-1					2.0E+03	1.8E+03		6.6E+02	
						3.0E+00	C	V		1.77	1	1	Yes	Propylene	115-07-1							6.3E+03	6.3E+03	
		2.0E+01	P							-0.92	1	1	Yes	Propylene Glycol	57-55-6					4.0E+05	3.2E+08		4.0E+05	
						2.7E-04	A			1.83	1	1	Yes	Propylene Glycol Dinitrate	6423-43-4									
		7.0E-01	H			2.0E+00	I	V		-0.49	1	1	Yes	Propylene Glycol Monomethyl Ether	107-98-2					1.4E+04	3.9E+06	4.2E+03	3.2E+03	
2.4E-01	I	3.7E-06	I			3.0E-02	I	V		0.03	1	1	Yes	Propylene Oxide	75-56-9	3.2E+01	4.7E+03	1.5E+02	2.7E+01			6.3E+01	6.3E+01	
		7.5E-02	I							3.43	1	0.9	Yes	Propylamide	23950-58-5					1.5E+03	5.5E+03		1.2E+03	
		1.0E-03	I					V		0.65	1	1	Yes	Pyridine	110-86-1					2.0E+01	1.5E+03		2.0E+01	
		5.0E-04	I							4.44	1	0.9	Yes	Quinalphos	13593-03-8					1.0E+01	1.0E+01		5.1E+00	
3.0E+00	I									2.03	1	1	Yes	Quinoline	91-22-5	2.6E+00	2.9E+01		2.4E+00					
		9.0E-03	I							4.28	1	0.9	Yes	Quizalofop-ethyl	76578-14-8					1.8E+02	3.8E+02		1.2E+02	
						3.0E-02	A				1	0	Yes	Refractory Ceramic Fibers	NA									
		3.0E-02	I							6.14	1	0.7	Yes	Resmethrin	10453-86-8					6.0E+02	7.6E+01		6.7E+01	
		5.0E-02	H					V		4.88	1	0.8	Yes	Ronnel	299-84-3					1.0E+03	6.8E+02		4.1E+02	
2.2E-01	C	6.3E-05	C							4.1	1	0.9	Yes	Rotenone	83-79-4					8.0E+01	2.6E+02		6.1E+01	
		5.0E-03	I							1	1	1	Yes	Safrole	94-59-7	1.1E+01	6.0E+01		9.6E+00					
										1	1	1	Yes	Selenious Acid	7783-00-8					1.0E+02	2.3E+04		1.0E+02	
		5.0E-03	I			2.0E-02	C			1	1	1	Yes	Selenium	7782-49-2					1.0E+02	2.3E+04		1.0E+02	5.0E+01
		5.0E-03	C			2.0E-02	C			1	1	1	Yes	Selenium Sulfide	7446-34-6					1.0E+02	2.3E+04		1.0E+02	
		9.0E-02	I							4.38	1	0.9	Yes	Sethoxydim	74051-80-2					1.8E+03	2.4E+03		1.0E+03	
						3.0E-03	C			1	1	1	Yes	Silica (crystalline, respirable)	7631-86-9									
1.2E-01	H					5.0E-03	I			0.04	1	1	Yes	Silver	7440-22-4					1.0E+02	1.5E+03		9.4E+01	
		5.0E-03	I							2.18	1	1	Yes	Simazine	127-34-9	6.5E+01	9.3E+02		6.1E+01	1.0E+02	1.6E+03		9.4E+01	4.0E+00
		1.3E-02	I							0.37	1	1	Yes	Sodium Acifluorfen	62476-59-9					2.6E+02	2.1E+05		2.6E+02	
		4.0E-03	I							1	1	1	Yes	Sodium Azide	26628-22-8					8.0E+01	1.8E+04		8.0E+01	
5.0E-01	C	1.5E-01	C			2.0E-02	C	2.0E-04	C	M	0.025	1	Yes	Sodium Dichromate	10588-01-9	5.0E+00	2.3E+01		4.1E+00	4.0E+02	2.3E+03		3.4E+02	
2.7E-01	H					3.0E-02	I			-1.43	1	1	Yes	Sodium Diethyldithiocarbamate	148-18-5	2.9E+01	8.5E+04		2.9E+01	6.0E+02	1.9E+06		6.0E+02	
		5.0E-02	A			1.3E-02	C			1	1	1	Yes	Sodium Fluoride	7681-49-4					1.0E+03	2.3E+05		1.0E+03	
		2.0E-05	I							-3.78	1	1	No	Sodium Fluoroacetate	62-74-8					4.0E-01			4.0E-01	
		1.0E-03	H							1	1	1	Yes	Sodium Metavanadate	13718-26-8					2.0E+01	4.6E+03		2.0E+01	
		8.0E-04	P							1	1	1	Yes	Sodium Tungstate	13472-45-2					1.6E+01	3.6E+03		1.6E+01	
		8.0E-04	P							1	1	1	Yes	Sodium Tungstate Dihydrate	10213-10-2					1.6E+01	3.6E+03		1.6E+01	
2.4E-02	H					3.0E-02	I			3.53	1	0.9	Yes	Stirofos (Tetrachlorovinphos)	961-11-5	3.2E+02	1.9E+03		2.8E+02	6.0E+02	3.8E+03		5.2E+02	
5.0E-01	C	1.5E-01	C			2.0E-02	C	2.0E-04	C	M	0.025	1	Yes	Strontium Chromate	7789-06-2	5.0E+00	2.3E+01		4.1E+00	4.0E+02	2.3E+03		3.4E+02	
						6.0E-01	I			1	1	1	Yes	Strontium, Stable	7440-24-6					1.2E+04	2.7E+06		1.2E+04	
		3.0E-04	I							1.93	1	1	Yes	Strychnine	57-24-9					6.0E+00	3.2E+02		5.9E+00	
		2.0E-01	I			1.0E+00	I	V		2.95	1	1	Yes	Styrene	100-42-5					4.0E+03	1.0E+04	2.1E+03	1.2E+03	1.0E+02
		3.0E-03	P							3.1	1	1	Yes	Styrene-Acrylonitrile (SAN) Trimer	NA				6.0E+01	2.4E+02		4.8E+01		
		1.0E-03	P			2.0E-03	X			-0.77	1	1	Yes	Sulfolane	126-33-0					2.0E+01	1.7E+04		2.0E+01	
		8.0E-04	P							3.9	1	0.9	Yes	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9					1.6E+01	3.5E+01		1.1E+01	
						1.0E-03	C	V		1	1	1	Yes	Sulfur Trioxide	7446-11-9							2.1E+00	2.1E+00	
						1.0E-03	C			1	1	1	Yes	Sulfuric Acid	7664-93-9									
2.5E-02	I	7.1E-06	I			5.0E-02	H			4.82	1	0.8	Yes	Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl ester	140-57-8	3.1E+02	2.3E+02		1.3E+02	1.0E+03	8.2E+02		4.5E+02	
						3.0E-02	H			3.3	1	0.9	Yes	TCMTB	21564-17-0					6.0E+02	2.4E+03		4.8E+02	
		7.0E-02	I							1.79	1	1	Yes	Tebuthiuron	34014-18-1					1.4E+03	4.7E+04		1.4E+03	
		2.0E-02	H							5.96	1	0.7	No	Temephos	3383-96-8					4.0E+02			4.0E+02	

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)													Toxicity and Chemical-specific Information				Contaminant				Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1			
SFO (mg/kg-day) ¹	k e y	IUR (ug/m ³) ¹	k e y	RfD _o (mg/kg-day)	k e y	RfC _o (mg/m ³)	k e y	v o l u t i l e	m u t a g e n	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL THQ=1 (ug/L)	Dermal SL THQ=1 (ug/L)	Inhalation SL THQ=1 (ug/L)	Noncarcinogenic SL THI=1 (ug/L)	MCL (ug/L)				
				1.3E-02	I					1.89	1	1	Yes	Terbacil	5902-51-2					2.6E+02	7.0E+03		2.5E+02					
				2.5E-05	H					4.48	1	0.9	Yes	Terbufos	13071-79-9					5.0E-01	4.5E-01		2.4E-01					
				1.0E-03	I					3.74	1	0.9	Yes	Terbutryn	886-50-0					2.0E+01	4.1E+01		1.3E+01					
				1.0E-04	I					6.77	1	0.6	No	Tetrabromodiphenyl ether, 2,2',4,4'-(BDE-47)	5436-43-1					2.0E+00			2.0E+00					
				3.0E-04	I					4.64	1	1	Yes	Tetrachlorobenzene, 1,2,4,5-	95-94-3					6.0E+00	2.4E+00		1.7E+00					
2.6E-02	I	7.4E-06	I	3.0E-02	I					2.93	1	1	Yes	Tetrachloroethane, 1,1,1,2-	630-20-6	3.0E+02	1.1E+03	7.6E+01	5.7E+01	6.0E+02	2.4E+03		4.8E+02					
2.0E-01	I	5.8E-05	C	2.0E-02	I					2.39	1	1	Yes	Tetrachloroethane, 1,1,2,2-	79-34-5	3.9E+01	3.3E+02	9.7E+00	7.6E+00	4.0E+02	3.6E+03		3.6E+02					
2.1E-03	I	2.6E-07	I	6.0E-03	I	4.0E-02	I	V		3.4	1	1	Yes	Tetrachloroethylene	127-18-4	3.7E+03	6.5E+03	2.2E+03	1.1E+03	1.2E+02	2.3E+02	8.3E+01	4.1E+01	5.0E+00				
2.0E+01	H			3.0E-02	I					4.45	1	0.9	Yes	Tetrachlorophenol, 2,3,4,6-	58-90-2					6.0E+02	3.9E+02		2.4E+02					
				5.0E-04	I					3.99	1	0.9	Yes	Tetraethyl Dithiopyrophosphate	3689-24-5					1.0E+01	2.4E+01		7.1E+00					
				8.0E+01	I	V				1.68	1	1	Yes	Tetrafluoroethane, 1,1,1,2-	811-97-2							1.7E+05	1.7E+05					
				2.0E-03	P					1.64	1	1	Yes	Tetryl (Trinitrophenylmethylnitramine)	479-45-8					4.0E+01	2.5E+03		3.9E+01					
				2.0E-05	S						1	0.9	Yes	Thallic Oxide	1314-32-5					4.0E-01	9.1E+01		4.0E-01					
				1.0E-05	X						1	1	Yes	Thallium (I) Nitrate	10102-45-1					2.0E-01	4.6E+01		2.0E-01					
				1.0E-05	X						1	1	Yes	Thallium (Soluble Salts)	7440-28-0					2.0E-01	4.6E+01		2.0E-01	2.0E+00				
				1.0E-05	X		V			-0.17	1	1	Yes	Thallium Acetate	563-68-8					2.0E-01	1.7E+02		2.0E-01					
				2.0E-05	X		V			-0.86	1	1	Yes	Thallium Carbonate	6533-73-9					4.0E-01	3.7E+03		4.0E-01					
				1.0E-05	X						1	1	Yes	Thallium Chloride	7791-12-0					2.0E-01	4.6E+01		2.0E-01					
				1.0E-05	S						1	1	Yes	Thallium Selenite	12039-52-0					2.0E-01	4.6E+01		2.0E-01					
				2.0E-05	X						1	0.9	Yes	Thallium Sulfate	7446-18-6					4.0E-01	9.1E+01		4.0E-01					
				1.3E-02	I					1.56	1	1	Yes	Thifensulfuron-methyl	79277-27-3					2.6E+02	3.5E+04		2.6E+02					
				1.0E-02	I					3.4	1	0.9	Yes	Thiobencarb	28249-77-6					2.0E+02	7.7E+02		1.6E+02					
				7.0E-02	X					-0.63	1	1	Yes	Thiodiglycol	111-48-8					1.4E+03	9.7E+05		1.4E+03					
				3.0E-04	H					2.16	1	1	Yes	Thiofanox	39196-18-4					6.0E+00	4.4E+01		5.3E+00					
				8.0E-02	I					1.4	1	1	Yes	Thiophanate, Methyl	23564-05-8					1.6E+03	2.1E+05		1.6E+03					
				5.0E-03	I					1.73	1	1	Yes	Thiram	13726-8					1.0E+02	4.0E+03		9.8E+01					
				6.0E-01	H						1	1	Yes	Tin	7440-31-5					1.2E+04	2.7E+06		1.2E+04					
				1.0E-04	A	V					1	1	Yes	Titanium Tetrachloride	7550-45-0								2.1E-01	2.1E-01				
				8.0E-02	I	5.0E+00	I	V		2.73	1	1	Yes	Toluene	10888-3					1.6E+03	5.3E+03	1.0E+04	1.1E+03	1.0E+03				
				1.1E-05	C	8.0E-06	C	V		3.74	1	1	Yes	Toluene-2,4-diisocyanate	58484-9			5.1E+01	5.1E+01			1.7E-02	1.7E-02					
1.8E-01	X			2.0E-04	X					0.16	1	1	Yes	Toluene-2,5-diamine	95-70-5	4.3E+01	8.2E+03		4.3E+01	4.0E+00	8.3E+02		4.0E+00					
				1.1E-05	C	8.0E-06	C	V		3.74	1	1	Yes	Toluene-2,6-diisocyanate	91-08-7			5.1E+01	5.1E+01			1.7E-02	1.7E-02					
1.6E-02	P	5.1E-05	C							1.32	1	1	Yes	Toluidine, o- (Methylaniline, 2-)	95-53-4	4.9E+02	1.4E+04		4.7E+02									
3.0E-02	P			4.0E-03	X					1.39	1	1	Yes	Toluidine, p-	106-49-0	2.6E+02	6.8E+03		2.5E+02	8.0E+01	2.3E+03		7.7E+01					
				3.0E+00	P		V			6.1	1	1	No	Total Petroleum Hydrocarbons (Aliphatic High)	NA					6.0E+04			6.0E+04					
				6.0E-01	P	V				3.9	1	1	Yes	Total Petroleum Hydrocarbons (Aliphatic Low)	NA							1.3E+03	1.3E+03					
				1.0E-02	X	1.0E-01	P	V		5.65	1	1	No	Total Petroleum Hydrocarbons (Aliphatic Medium)	NA					2.0E+02		2.1E+02	1.0E+02					
				4.0E-02	P					5.16	1	1	No	Total Petroleum Hydrocarbons (Aromatic High)	NA					8.0E+02			8.0E+02					
				4.0E-03	P	3.0E-02	P	V		2.13	1	1	Yes	Total Petroleum Hydrocarbons (Aromatic Low)	NA					8.0E+01	6.1E+02	6.3E+01	3.3E+01					
				4.0E-03	P	3.0E-03	P	V		3.58	1	1	Yes	Total Petroleum Hydrocarbons (Aromatic Medium)	NA					8.0E+01	9.0E+01	6.3E+00	5.5E+00					
1.1E+00	I	3.2E-04	I							5.9	1	0.8	No	Toxaphene	8001-35-2	7.1E+00			7.1E+00					3.0E+00				
				7.5E-03	I					7.56	1	0.5	No	Tralometrin	66841-25-6								1.5E+02					
				3.0E-04	A		V			4.1	1	0.9	Yes	Tri-n-butyltin	688-73-3					6.0E+00	9.9E+00		3.7E+00					
				8.0E+01	X					0.25	1	1	Yes	Triacetin	102-76-1					1.6E+06	5.3E+08		1.6E+06					
				3.0E-02	I					2.77	1	1	Yes	Triadimefon	43121-43-3					6.0E+02	6.9E+03		5.5E+02					
				1.3E-02	I		V			4.6	1	0.9	Yes	Triallate	2303-17-5					2.6E+02	2.2E+02		1.2E+02					
				1.0E-02	I					1.1	1	1	Yes	Triasulfuron	82097-50-5					2.0E+02	6.0E+04		2.0E+02					
				8.0E-03	I					0.78	1	1	Yes	Tribenuron-methyl	101200-48-0					1.6E+02	5.0E+03		1.6E+02					
				5.0E-03	I		V			4.66	1	0.9	Yes	Tribromobenzene, 1,2,4-	615-54-3					1.0E+02	8.1E+01		4.5E+01					
9.0E-03	P			1.0E-02	P					4	1	0.9	Yes	Tributyl Phosphate	126-73-8	8.7E+02	1.3E+03		5.2E+02	2.0E+02	3.3E+02		1.2E+02					
				3.0E-04	P						1	0	No	Tributyltin Compounds	NA					6.0E+00			6.0E+00					
				3.0E-04	I					4.05	1	1	Yes	Tributyltin Oxide	56-35-9					6.0E+00	9.5E+01		5.7E+00					
				3.0E+01	I	3.0E+01	H	V		3.16	1	1	Yes	Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1					6.0E+05	1.9E+06	6.3E+04	5.5E+04					

Regional Removal Management Level (RML) Resident Tapwater Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)													Toxicity and Chemical-specific Information				Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer CHILD Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	ke y	IUR (ug/m ³) ⁻¹	ke y	RfD _o (mg/kg-day)	ke y	RfC _o (mg/m ³)	ke y	o I	muta- gen	LOGP	GIABS	FA	In EPD?	Analyte	CAS No.	Ingestion SL TR=1E-04 (ug/L)	Dermal SL TR=1E-04 (ug/L)	Inhalation SL TR=1E-04 (ug/L)	Carcinogenic SL TR=1E-04 (ug/L)	Ingestion SL THQ=1 (ug/L)	Dermal Child THQ=1 (ug/L)	Inhalation SL Child THQ=1 (ug/L)	Noncarcinogenic SL Child THI=1 (ug/L)	MCL (ug/L)		
7.0E-02	I			2.0E-02	I					1.33	1	1	Yes	Trichloroacetic Acid	76-03-9	1.1E+02	4.6E+03		1.1E+02	4.0E+02	1.8E+04		3.9E+02	6.0E+01		
2.9E-02	H									-0.67	1	1	Yes	Trichloroaniline HCl, 2,4,6-	33663-50-2	2.7E+02	3.7E+05		2.7E+02							
7.0E-03	X			3.0E-05	X					3.52	1	1	Yes	Trichloroaniline, 2,4,6-	634-93-5	1.1E+03	2.0E+03		7.1E+02	6.0E-01	1.2E+00		4.0E-01			
				8.0E-04	X			V		4.05	1	1	Yes	Trichlorobenzene, 1,2,3-	87-61-6					1.6E+01	1.3E+01		7.0E+00			
2.9E-02	P			1.0E-02	I	2.0E-03	P	V		4.02	1	1	Yes	Trichlorobenzene, 1,2,4-	120-82-1	2.7E+02	2.0E+02		1.2E+02	2.0E+02	1.6E+02	4.2E+00	4.0E+00	7.0E+01		
				2.0E+00	I	5.0E+00	I	V		2.49	1	1	Yes	Trichloroethane, 1,1,1-	71-55-6					4.0E+04	2.5E+05	1.0E+04	8.0E+03	2.0E+02		
5.7E-02	I	1.6E-05	I	4.0E-03	I	2.0E-04	X	V		1.89	1	1	Yes	Trichloroethane, 1,1,2-	79-00-5	1.4E+02	2.0E+03	3.5E+01	2.8E+01	8.0E+01	1.3E+03	4.2E-01	4.1E-01	5.0E+02		
4.6E-02	I	4.1E-06	I	5.0E-04	I	2.0E-03	I	V	M	2.42	1	1	Yes	Trichloroethylene	79-01-6	1.2E+02	7.4E+02	9.6E+01	4.9E+01	1.0E+01	6.9E+01	4.2E+00	2.8E+00	5.0E+00		
				3.0E-01	I			V		2.53	1	1	Yes	Trichlorofluoromethane	75-69-4					6.0E+03	3.6E+04		5.2E+03			
				1.0E-01	I					3.72	1	1	Yes	Trichlorophenol, 2,4,5-	95-95-4					2.0E+03	2.9E+03		1.2E+03			
1.1E-02	I	3.1E-06	I	1.0E-03	P					3.69	1	1	Yes	Trichlorophenol, 2,4,6-	88-06-2	7.1E+02	9.8E+02		4.1E+02	2.0E+01	3.0E+01		1.2E+01			
				1.0E-02	I					3.31	1	0.9	Yes	Trichlorophenoxyacetic Acid, 2,4,5-	93-76-5					2.0E+02	8.7E+02		1.6E+02			
				8.0E-03	I					3.8	1	0.9	Yes	Trichlorophenoxypropionic acid, -2,4,5	93-72-1					1.6E+02	3.6E+02		1.1E+02	5.0E+01		
				5.0E-03	I			V		2.43	1	1	Yes	Trichloropropane, 1,1,2-	598-77-6					1.0E+02	7.5E+02		8.8E+01			
3.0E+01	I			4.0E-03	I	3.0E-04	I	V	M	2.27	1	1	Yes	Trichloropropane, 1,2,3-	96-18-4	8.4E-02	7.3E-01		7.5E-02	8.0E+01	7.7E+02	6.3E-01	6.2E-01			
				3.0E-03	X	3.0E-04	P	V		2.78	1	1	Yes	Trichloropropene, 1,2,3-	96-19-5					6.0E+01	2.6E+02	6.3E-01	6.2E-01			
				2.0E-02	A					5.11	1	0.8	Yes	Tricresyl Phosphate (TCP)	1330-78-5					4.0E+02	2.6E+02		1.6E+02			
				3.0E-03	I					5.18	1	0.8	Yes	Tridiphane	58138-08-2					6.0E+01	2.6E+01		1.8E+01			
				7.0E-03	I	V				1.45	1	1	Yes	Triethylamine	121-44-8							1.5E+01		1.5E+01		
				2.0E+00	P					-1.75	1	1	Yes	Triethylene Glycol	112-27-6					4.0E+04	1.8E+08		4.0E+04			
				2.0E+01	P	V				1.74	1	1	Yes	Trifluoroethane, 1,1,1-	420-46-2							4.2E+04		4.2E+04		
7.7E-03	I			7.5E-03	I			V		5.34	1	0.8	Yes	Trifluralin	1582-09-8	1.0E+03	3.4E+02		2.6E+02	1.5E+02	5.5E+01		4.0E+01			
2.0E-02	P			1.0E-02	P					-0.65	1	1	Yes	Trimethyl Phosphate	512-56-1	3.9E+02	2.8E+05		3.9E+02	2.0E+02	1.6E+05		2.0E+02			
				5.0E-03	P	V				3.66	1	1	Yes	Trimethylbenzene, 1,2,3-	526-73-8							1.0E+01	1.0E+01			
				7.0E-03	P	V				3.63	1	1	Yes	Trimethylbenzene, 1,2,4-	95-63-6							1.5E+01		1.5E+01		
				1.0E-02	X			V		3.42	1	1	Yes	Trimethylbenzene, 1,3,5-	108-67-8					2.0E+02	2.8E+02		1.2E+02			
				1.0E-02	X			V		4.08	1	1	Yes	Trimethylpentene, 2,4,4-	25167-70-8					2.0E+02	9.6E+01		6.5E+01			
				3.0E-02	I					1.18	1	1	Yes	Trinitrobenzene, 1,3,5-	99-35-4					6.0E+02	4.7E+04		5.9E+02			
3.0E-02	I			5.0E-04	I					1.6	1	1	Yes	Trinitrotoluene, 2,4,6-	118-96-7	2.6E+02	1.1E+04		2.5E+02	1.0E+01	4.5E+02		9.8E+00			
				2.0E-02	P					2.83	1	1	Yes	Triphenylphosphine Oxide	791-28-6					4.0E+02	3.8E+03		3.6E+02			
				2.0E-02	A					3.65	1	0.9	Yes	Tris(1,3-Dichloro-2-propyl)phosphate	13674-87-8					4.0E+02	3.2E+03		3.6E+02			
2.3E+00	C	6.6E-04	C	1.0E-02	X					2.59	1	1	Yes	Tris(1-chloro-2-propyl)phosphate	13674-84-5					2.0E+02	3.8E+03		1.9E+02			
2.0E-02	P			7.0E-03	P					1.44	1	1	Yes	Tris(2,3-dibromopropyl)phosphate	12672-7	3.4E+00		8.5E-01	6.8E-01					1.4E+02		
				1.0E-01	P					9.49	1	0	No	Tris(2-chloroethyl)phosphate	11596-8	3.9E+02	3.0E+04		3.8E+02	1.4E+02	1.2E+04		1.4E+02			
3.2E-03	P			1.0E-01	P						1	1	Yes	Tris(2-ethylhexyl)phosphate	78-42-2	2.4E+03		2.4E+03		2.0E+03			2.0E+03			
				8.0E-04	P						1	1	Yes	Tungsten	7440-33-7					1.6E+01	3.6E+03		1.6E+01			
				3.0E-03	I	4.0E-05	A				1	1	Yes	Uranium (Soluble Salts)	NA					6.0E+01	1.4E+04		6.0E+01	3.0E+01		
1.0E+00	C	2.9E-04	C	8.3E-03	P				M	-0.15	1	1	Yes	Urethane	51-79-6	2.5E+00	6.1E+02		2.5E+00							
				9.0E-03	I	7.0E-06	P			0.026	1	1	Yes	Vanadium Pentoxide	1314-62-1					1.8E+02	1.1E+03		1.5E+02			
				5.0E-03	S	1.0E-04	A			0.026	1	1	Yes	Vanadium and Compounds	7440-62-2					1.0E+02	6.0E+02		8.6E+01			
				1.0E-03	I			V		3.84	1	1	Yes	Vernolate	1929-77-7					2.0E+01	2.5E+01		1.1E+01			
				2.5E-02	I					3.1	1	0.9	Yes	Vinclozolin	50471-44-8					5.0E+02	3.7E+03		4.4E+02			
				1.0E+00	H	2.0E-01	I	V		0.73	1	1	Yes	Vinyl Acetate	108-05-4					2.0E+04	1.4E+06	4.2E+02	4.1E+02			
7.2E-01	I	4.4E-06	I	3.0E-03	I	1.0E-01	I	V	M	1.57	1	1	Yes	Vinyl Bromide	593-60-2			1.8E+01	1.8E+01	6.0E+01	8.9E+02	2.1E+02	6.3E+00	2.0E+00		
				3.0E-04	I					2.7	1	1	Yes	Vinyl Chloride	75-01-4	2.1E+00	2.8E+01	3.4E+01	1.9E+00	6.0E+01	8.4E+01		4.4E+01			
				2.0E-01	S	1.0E-01	S	V		3.15	1	1	Yes	Warfarin	81-81-2					6.0E+00			5.6E+00			
				2.0E-01	S	1.0E-01	S	V		3.2	1	1	Yes	Xylene, p-	106-42-3					4.0E+03	7.6E+03	2.1E+02	1.9E+02			
				2.0E-01	S	1.0E-01	S	V		3.2	1	1	Yes	Xylene, m-	108-38-3					4.0E+03	7.1E+03	2.1E+02	1.9E+02			
				2.0E-01	S	1.0E-01	S	V		3.12	1	1	Yes	Xylene, o-	95-47-6					4.0E+03	8.0E+03	2.1E+02	1.9E+02			
				2.0E-01	I	1.0E-01	I	V		3.16	1	1	Yes	Xylenes	1330-20-7					4.0E+03	7.5E+03	2.1E+02	1.9E+02	1.0E+04		
				3.0E-04	I						1	1	Yes	Zinc Phosphide	1314-84-7					6.0E+00	2.3E+03		6.0E+00			
				3.0E-01	I						1	1	Yes	Zinc and Compounds	7440-66-6					6.0E+03	2.3E+06		6.0E+03			
				5.0E-02	I					1.3	1	1	Yes	Zineb	12122-67-7					1.0E+03	9.7E+04		9.9E+02			
				8.0E-05	X						1	1	Yes	Zirconium	7440-67-7					1.6E+00	3.6E+02		1.6E+00			

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)																													
Toxicity and Chemical-specific Information															Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1								
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³) ⁻¹	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³)	k _e (y)	o ₁	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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)					
8.7E-03	I	2.2E-06	I	4.0E-03	I	9.0E-03	I	V		1.1E+05	1.4E+09	8.7E+03	1	0.1	Acephate	30560-19-1	3.8E+04	8.9E+04	4.9E+03	2.6E+04	4.7E+03	1.1E+04	3.4E+02	3.3E+03					
				2.0E-02	I						1.4E+09		1	0.1	Acetaldehyde	75-07-0								3.4E+02					
													1		Acetochlor	34256-82-1					2.3E+04	5.5E+04		1.6E+04					
				9.0E-01	I	3.1E+01	A	V		1.1E+05	1.4E+09	1.4E+04	1		Acetone	67-64-1					1.1E+06		1.8E+06	6.7E+05					
											1.4E+09		1	0.1	Acetone Cyanohydrin	75-86-5							1.2E+07						
											1.3E+05	1.4E+09	1.3E+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		2.5E+03	1.4E+09	6.0E+04	1		Acetophenone	98-86-2					1.2E+05			1.2E+05					
											1.4E+09		1	0.1	Acetylaminofluorene, 2-Acrolein	53-96-3 107-02-8	8.6E+01	2.0E+02	1.3E+06	6.0E+01	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			1.4E+09		1	0.1	Acrylamide	79-06-1	6.5E+02	1.5E+03	1.7E+07	4.6E+02	2.3E+03	5.5E+03	3.6E+07	1.6E+03					
				5.0E-01	I	1.0E-03	I	V		1.1E+05	1.4E+09	9.5E+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		1.1E+04	1.4E+09	7.7E+03	1		Acrylonitrile	107-13-1	6.1E+02		1.4E+02	1.1E+02	4.7E+04		6.7E+01	6.7E+01					
											1.4E+09		1	0.1	Adiponitrile	111-69-3							3.6E+07						
5.6E-02	C			1.0E-02	I						1.4E+09		1	0.1	Alachlor	15972-60-8	5.8E+03	1.4E+04		4.1E+03	1.2E+04	2.8E+04		8.2E+03					
				1.0E-03	I						1.4E+09		1	0.1	Aldicarb	116-06-3					1.2E+03	2.8E+03		8.2E+02					
				1.0E-03	I						1.4E+09		1	0.1	Aldicarb Sulfone	1646-88-4					1.2E+03	2.8E+03		8.2E+02					
1.7E+01	I	4.9E-03	I	3.0E-05	I			V			1.4E+09	1.7E+06	1		Aldicarb sulfoxide	1646-87-3													
											1.4E+09		1		Aldrin	309-00-2	1.9E+01		4.3E+02	1.8E+01	3.5E+01			3.5E+01					
2.1E-02	C	6.0E-06	C	5.0E-03	I	1.0E-04	X	V		1.1E+05	1.4E+09	3.4E+04	1		Allyl Alcohol	107-18-6					5.8E+03		1.5E+01	1.5E+01					
											1.4E+09	1.6E+03	1		Allyl Chloride	107-05-1	1.6E+04		3.2E+02	3.2E+02			6.9E+00	6.9E+00					
				1.0E+00	P	5.0E-03	P				1.4E+09		1		Aluminum	7429-90-5					1.2E+06		3.0E+07	1.1E+06					
				4.0E-04	I						1.4E+09		1		Aluminum Phosphide	20859-73-8					4.7E+02			4.7E+02					
2.1E+01	C	6.0E-03	C	9.0E-03	I						1.4E+09		1	0.1	Ametryn	834-12-8					1.1E+04	2.5E+04		7.4E+03					
											1.4E+09		1	0.1	Aminobiphenyl, 4-	92-67-1	1.6E+01	3.7E+01	2.8E+05	1.1E+01									
				8.0E-02	P						1.4E+09		1	0.1	Aminophenol, m-	591-27-5					9.3E+04	2.2E+05		6.6E+04					
				2.0E-02	P						1.4E+09		1	0.1	Aminophenol, p-	123-30-8					2.3E+04	5.5E+04		1.6E+04					
				2.5E-03	I						1.4E+09		1	0.1	Amitraz	33089-61-1					2.9E+03	6.9E+03		2.1E+03					
						1.0E-01	I	V					1		Ammonia	7664-41-7								2.3E+05					
				2.0E-01	I						1.4E+09		1		Ammonium Sulfamate	7773-06-0									2.3E+05				
						3.0E-03	X	V		1.4E+04	1.4E+09	2.6E+04	1		Amyl Alcohol, tert-	75-85-4							3.4E+02	3.4E+02					
5.7E-03	I	1.6E-06	C	7.0E-03	P	1.0E-03	I				1.4E+09		1	0.1	Aniline	62-53-3	5.7E+04	1.4E+05	1.0E+09	4.0E+04	8.2E+03	1.9E+04	6.0E+06	5.7E+03					
4.0E-02	P			2.0E-03	X						1.4E+09		1	0.1	Anthraquinone, 9,10-	84-65-1	8.2E+03	1.9E+04		5.7E+03	2.3E+03	5.5E+03		1.6E+03					
				4.0E-04	I						1.4E+09	0.15			Antimony (metalloid)	7440-36-0					4.7E+02			4.7E+02					
				5.0E-04	H						1.4E+09	0.15			Antimony Pentoxide	1314-60-9					5.8E+02			5.8E+02					
				4.0E-04	H						1.4E+09	0.15			Antimony Tetroxide	1332-81-6					4.7E+02			4.7E+02					
						2.0E-04	I				1.4E+09	0.15			Antimony Trioxide	1309-64-4							1.2E+06	1.2E+06					
1.5E+00	I	4.3E-03	I	3.0E-04	I	1.5E-05	C				1.4E+09		1	0.03	Arsenic, Inorganic	7440-38-2	3.6E+02	1.7E+03	3.9E+05	3.0E+02	5.8E+02	2.8E+03	8.9E+04	4.8E+02					
				3.5E-06	C	5.0E-05	I				1.4E+09		1		Arsine	7784-42-1					4.1E+00		3.0E+05	4.1E+00					
				5.0E-02	I						1.4E+09		1	0.1	Asulam	3337-71-1					5.8E+04	1.4E+05		4.1E+04					
2.3E-01	C	3.5E-02	I	3.5E-02	I						1.4E+09		1	0.1	Atrazine	1912-24-9	1.4E+03	3.4E+03		1.0E+03	4.1E+04	9.7E+04			2.9E+04				
8.8E-01	C	2.5E-04	C	4.0E-04	I						1.4E+09		1	0.1	Auramine	492-80-8	3.7E+02	8.8E+02	6.7E+06	2.6E+02									
											1.4E+09		1	0.1	Avermectin B1	65195-55-3					4.7E+02	1.1E+03		3.3E+02					
1.1E-01	I	3.1E-05	I	3.0E-03	A	1.0E-02	A				1.4E+09		1	0.1	Azinphos-methyl	86-50-0					3.5E+03	8.3E+03	6.0E+07	2.5E+03					
								V			1.4E+09	5.2E+05	1		Azobenzene	103-33-3	3.0E+03		2.1E+04	2.6E+03									
				1.0E+00	P	7.0E-06	P				1.4E+09		1	0.1	Azodicarbonamide	123-77-3					1.2E+06	2.8E+06	4.2E+04	4.0E+04					
5.0E-01	C	1.5E-01	C	2.0E-01	I	5.0E-04	H				1.4E+09	0.07			Barium	7440-39-3					2.3E+05		3.0E+06	2.2E+05					
				2.0E-02	C	2.0E-04	C	M			1.4E+09	0.025			Barium Chromate	10294-40-3	6.5E+02		1.1E+04	6.2E+02	2.3E+04		1.2E+06	2.3E+04					
				3.0E-01	I		V				1.4E+09	3.1E+05	1		Benfluralin	1861-40-1					3.5E+05			3.5E+05					
				5.0E-02	I						1.4E+09		1	0.1	Benomyl	17804-35-2					5.8E+04	1.4E+05		4.1E+04					
				2.0E-01	I						1.4E+09		1	0.1	Bensulfuron-methyl	83055-99-6					2.3E+05	5.5E+05		1.6E+05					
				3.0E-02	I						1.4E+09		1	0.1	Bentazon	25057-89-0					3.5E+04	8.3E+04		2.5E+04					
4.0E-03	P			1.0E-01	I			V		1.2E+03	1.4E+09	2.3E+04	1		Benzaldehyde	100-52-7	8.2E+04			8.2E+04	1.2E+05			1.2E+05					
5.5E-02	I	7.8E-06	I	4.0E-03	I	3.0E-02	I	V		1.8E+03	1.4E+09	3.5E+03	1		Benzene	71-43-2	5.9E+03		5.6E+02	5.1E+02	4.7E+03		4.6E+02	4.2E+02					
1.0E-01	X			3.0E-04	X						1.4E+09		1	0.1	Benzenediamine-2-methyl sulfate, 1,4-	6369-59-1	3.3E+03	7.7E+03		2.3E+03	3.5E+02	8.3E+02		2.5E+02					
2.3E+02	I	6.7E-02	I	1.0E-03	P			V		1.3E+03	1.4E+09	1.9E+04	1		Benzenethiol	108-98-5					1.2E+03			1.2E+03					
				3.0E-03	I			M			1.4E+09		1	0.1	Benzidine	92-87-5	1.4E+00	3.4E+00	2.5E+04	1.0E+00	3.5E+03	8.3E+03		2.5E+03					
				4.0E+00	I						1.4E+09		1																

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)																													
Toxicity and Chemical-specific Information															Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1								
SFO (mg/kg-day) ⁻¹	k _e IUR (ug/m ³) ⁻¹ y	k _e RfD _o (mg/kg-day)	k _e RfC _i (mg/m ³) y	k _e 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-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	Ingestion SL THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THH=1 (mg/kg)										
1.7E-01	I 4.9E-05	C 2.0E-03	P 1.0E-03	P V	1.5E+03	1.4E+09	2.6E+04	1		Benzyl Chloride	100-44-7	1.9E+03		6.4E+02	4.8E+02	2.3E+03		1.1E+02	1.1E+02										
	2.4E-03	I 2.0E-03	I 2.0E-05	I		1.4E+09			0.007	Beryllium and compounds	7440-41-7			6.9E+05	6.9E+05	2.3E+03		1.2E+05	2.3E+03										
		9.0E-03	P			1.4E+09			1	Bifenox	42576-02-3					1.1E+04	2.5E+04		7.4E+03										
		1.5E-02	I			1.4E+09			1	Biphenthrin	82657-04-3					1.8E+04	4.1E+04		1.2E+04										
8.0E-03	I	5.0E-01	I 4.0E-04	X V		1.4E+09	1.1E+05	1		Biphenyl, 1,1'-	92-52-4	4.1E+04			4.1E+04	5.8E+05		2.0E+02	2.0E+02										
		4.0E-02	I	V		1.0E+03	1.4E+09	3.5E+04	1	Bis(2-chloro-1-methylethyl) ether	108-60-1					4.7E+04			4.7E+04										
		3.0E-03	P			1.4E+09			1	Bis(2-chloroethoxy)methane	111-91-1					3.5E+03	8.3E+03		2.5E+03										
1.1E+00	I 3.3E-04	I		V		5.1E+03	1.4E+09	4.3E+04	1	Bis(2-chloroethyl)ether	111-44-4	3.0E+02		1.6E+02	1.0E+02														
2.2E+02	I 6.2E-02	I		V		4.2E+03	1.4E+09	1.9E+03	1	Bis(chloromethyl)ether	542-88-1	1.5E+00		3.7E-02	3.6E-02														
		5.0E-02	I			1.4E+09			1	Bisphenol A	80-05-7					5.8E+04	1.4E+05		4.1E+04										
		2.0E-01	I 2.0E-02	H		1.4E+09			1	Boron And Borates Only	7440-42-8					2.3E+05		1.2E+08	2.3E+05										
		2.0E+00	P 2.0E-02	P V		1.4E+09			1	Boron Trichloride	10294-34-5					2.3E+06		1.2E+08	2.3E+06										
		4.0E-02	C 1.3E-02	C V		1.4E+09			1	Boron Trifluoride	7637-07-2					4.7E+04		7.7E+07	4.7E+04										
7.0E-01	I	4.0E-03	I			1.4E+09			1	Bromate	15541-45-4	4.7E+02			4.7E+02	4.7E+03			4.7E+03										
2.0E+00	X 6.0E-04	X		V		2.4E+03	1.4E+09	5.9E+03	1	Bromo-2-chloroethane, 1-	107-04-0	1.6E+02		1.2E+01	1.1E+01														
		8.0E-03	I 6.0E-02	I V		6.8E+02	1.4E+09	8.4E+03	1	Bromobenzene	108-86-1					9.3E+03		2.2E+03	1.8E+03										
			4.0E-02	X V		4.0E+03	1.4E+09	3.6E+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		9.3E+02	1.4E+09	4.0E+03	1	Bromodichloromethane	75-27-4	5.3E+03		1.3E+02	1.3E+02	2.3E+04			2.3E+04										
7.9E-03	I 1.1E-06	I 2.0E-02	I	V		9.2E+02	1.4E+09	9.7E+03	1	Bromoform	75-25-2	4.1E+04		1.1E+04	8.6E+03	2.3E+04			2.3E+04										
		1.4E-03	I 5.0E-03	I V		3.6E+03	1.4E+09	1.4E+03	1	Bromomethane	74-83-9					1.6E+03		3.1E+01	3.0E+01										
		5.0E-03	H	V		1.4E+09	1.2E+05		1	Bromophos	2104-96-3					5.8E+03			5.8E+03										
		2.0E-02	I			1.4E+09			1	Bromoxynil	1689-84-5					2.3E+04	5.5E+04		1.6E+04										
		2.0E-02	I	V		1.4E+09	4.7E+05		1	Bromoxynil Octanoate	1689-99-2					2.3E+04			2.3E+04										
3.4E+00	C 3.0E-05	I	2.0E-03	I V		6.7E+02	1.4E+09	8.7E+02	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		7.6E+03	1.4E+09	3.0E+04	1	Butanol, N-	71-36-3					1.2E+05			1.2E+05										
		2.0E+00	P 3.0E+01	P V		2.1E+04	1.4E+09	2.9E+04	1	Butyl alcohol, sec-	78-92-2					2.3E+06		3.8E+06	1.5E+06										
		5.0E-02	I	V		1.4E+09	8.6E+04		1	Butylate	2008-41-5					5.8E+04			5.8E+04										
2.0E-04	C 5.7E-08	C		V		1.4E+09			1	Butylated hydroxyanisole	25013-16-5	1.6E+06	3.9E+06	2.9E+10	1.1E+06														
3.6E-03	P	3.0E-01	P			1.4E+09			1	Butylated hydroxytoluene	128-37-0	9.1E+04	2.1E+05		6.4E+04	3.5E+05	8.3E+05		2.5E+05										
		5.0E-02	P	V		1.1E+02	1.4E+09	8.1E+03	1	Butylbenzene, n-	104-51-8					5.8E+04			5.8E+04										
		1.0E-01	X	V		1.5E+02	1.4E+09	7.4E+03	1	Butylbenzene, sec-	135-98-8					1.2E+05			1.2E+05										
		1.0E-01	X	V		1.8E+02	1.4E+09	7.4E+03	1	Butylbenzene, tert	98-06-6					1.2E+05			1.2E+05										
		2.0E-02	A	V		1.4E+09			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		1.4E+09			0.025	Cadmium (Diet)	7440-39-9			9.3E+05	9.3E+05	1.2E+03	6.9E+03	6.0E+04	9.8E+02										
	1.8E-03	I 5.0E-04	I 1.0E-05	A		1.4E+09			0.05	Cadmium (Water)	7440-39-9					1.2E+03	6.9E+03	6.0E+04	9.8E+02										
5.0E-01	C 1.5E-01	C 2.0E-02	C 2.0E-04	C M		1.4E+09			0.025	Calcium Chromate	13765-19-0	6.5E+02		1.1E+04	6.2E+02	2.3E+04		1.2E+06	2.3E+04										
		5.0E-01	I 2.2E-03	C		1.4E+09			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			1.4E+09			1	Captafol	2425-06-1	2.2E+03	5.2E+03	3.9E+07	1.5E+03	2.3E+03	5.5E+03		1.6E+03										
2.3E-03	C 6.6E-07	C 1.3E-01	I			1.4E+09			1	Captan	133-06-2	1.4E+05	3.4E+05	2.5E+09	1.0E+05	1.5E+05	3.6E+05		1.1E+05										
		1.0E-01	I			1.4E+09			1	Carbaryl	63-25-2					1.2E+05	2.8E+05		8.2E+04										
		5.0E-03	I			1.4E+09			1	Carbofuran	1563-66-2					5.8E+03	1.4E+04		4.1E+03										
		1.0E-01	I 7.0E-01	I V		7.4E+02	1.4E+09	1.2E+03	1	Carbon Disulfide	75-15-0					1.2E+05		3.6E+03	3.5E+03										
7.0E-02	I 6.0E-06	I 4.0E-03	I 1.0E-01	I V		4.6E+02	1.4E+09	1.5E+03	1	Carbon Tetrachloride	56-23-5	4.7E+03		3.1E+02	2.9E+02	4.7E+03		6.5E+02	5.7E+02										
			1.0E-01	P V		5.9E+03	1.4E+09	6.5E+02	1	Carbonyl Sulfide	463-58-1							2.8E+02	2.8E+02										
		1.0E-02	I			1.4E+09			1	Carbosulfan	55285-14-8					1.2E+04	2.8E+04		8.2E+03										
		1.0E-01	I			1.4E+09			1	Carboxin	5234-68-4					1.2E+05	2.8E+05		8.2E+04										
			9.0E-04	I		1.4E+09			1	Ceric oxide	1306-38-3							5.4E+06	5.4E+06										
		1.0E-01	I	V		1.4E+09	1.5E+05		1	Chloral Hydrate	302-17-0					1.2E+05			1.2E+05										
		1.5E-02	I			1.4E+09			1	Chloramben	133-90-4					1.8E+04	4.1E+04		1.2E+04										
4.0E-01	H					1.4E+09			1	Chloranil	118-75-2	8.1E+02	1.9E+03		5.7E+02														
3.5E-01	I 1.0E-04	I 5.0E-04	I 7.0E-04	I V		1.4E+09	1.5E+06		1	Chlordane	12789-03-6	9.3E+02	5.5E+03	1.9E+04	7.7E+02	5.8E+02	3.4E+03	4.7E+03	4.5E+02										
1.0E+01	I 4.6E-03	C 3.0E-04	I			1.4E+09			1	Chlordecone (Kepone)	143-50-0	3.3E+01	7.7E+01	3.6E+05	2.3E+01	3.5E+02	8.3E+02		2.5E+02										
		7.0E-04	A			1.4E+09			1	Chlorfenvinphos	470-90-6					8.2E+02	1.9E+03		5.7E+02										
		2.0E-02	I			1.4E+09			1	Chlorimuron, Ethyl-	90982-32-4					2.3E+04	5.5E+04		1.6E+04										
		1.0E-01	I 1.5E-04	A V		2.8E+03	1.4E+09	1.2E+03	1	Chlorine	7782-50-5					1.2E+05		7.8E-01	7.8E-01										
		3.0E-02	I 2.0E-04	I V		1.4E+09			1	Chlorine Dioxide	10049-04-4					3.5E+04		1.2E+06	3.4E+04										
		3.0E-02	I			1.4E+09			1	Chlorite (Sodium Salt)	7758-19-2					3.5E+04			3.5E+04										
			5.0E+01	I V		1.2E+03	1.4E+09	1.0E+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		7.9E+02	1.4E+09	1.1E+03	1	Chloro-1,3-butadiene, 2-	126-99-8			4.4E+00	4.4E+00	2.3E+04		9.4E+01	9.4E+01										

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)																								
Toxicity and Chemical-specific Information															Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k e	IUR (ug/m ³) ⁻¹	k e	Rfd _o (mg/kg- day)	k e	RfC _i (mg/m ³) ⁻¹	k e	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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)	
4.6E-01	H									1.4E+09			0.1	Chloro-2-methylaniline HCl, 4-	3165-93-3	7.1E+02	1.7E+03		5.0E+02					
1.0E-01	P	7.7E-05	C	3.0E-03	X					1.4E+09			0.1	Chloro-2-methylaniline, 4-	95-69-2	3.3E+03	7.7E+03	2.2E+07	2.3E+03	3.5E+03	8.3E+03		2.5E+03	
2.7E-01	X									1.2E+04	1.6E+04	1	0.1	Chloroacetaldehyde, 2-	107-20-0	1.2E+03			1.2E+03					
										1.4E+09			0.1	Chloroacetic Acid	79-11-8									
										3.0E-05			0.1	Chloroacetophenone, 2-	532-27-4						1.8E+05	1.8E+05		
2.0E-01	P			4.0E-03	I					1.4E+09			0.1	Chloroaniline, p-	106-47-8	1.6E+03	3.9E+03		1.1E+03	4.7E+03	1.1E+04		3.3E+03	
1.0E-01	P	7.7E-05	C	3.0E-03	X					1.4E+09			0.1	Chlorobenzene	108-90-7					2.3E+04		1.4E+03	1.3E+03	
1.1E-01	C	3.1E-05	C	2.0E-02	I					1.4E+09			0.1	Chlorobenzilate	510-15-6	3.0E+03	7.0E+03	5.4E+07	2.1E+03	2.3E+04	5.5E+04		1.6E+04	
										3.0E-02			0.1	Chlorobenzoic Acid, p-	74-11-3					3.5E+04	8.3E+04		2.5E+04	
										3.0E-03			0.1	Chlorobenzotrifluoride, 4-	98-56-6					3.5E+03		8.9E+03	2.5E+03	
										4.0E-02			0.1	Chlorobutane, 1-	109-69-3					4.7E+04			4.7E+04	
										5.0E+01			0.1	Chlorodifluoromethane	75-45-6							2.1E+05	2.1E+05	
										2.0E-02			0.1	Chloroethanol, 2-	107-07-3					2.3E+04			2.3E+04	
3.1E-02	C	2.3E-05	I	1.0E-02	I	9.8E-02	A	V		2.5E+03	1.4E+09	2.6E+03	1	Chloroform	67-66-3	1.1E+04		1.4E+02	1.4E+02	1.2E+04		1.1E+03	1.0E+03	
										9.0E-02			0.1	Chloromethane	74-87-3							4.6E+02	4.6E+02	
2.4E+00	C	6.9E-04	C							9.3E+03	1.4E+09	5.3E+03	1	Chloromethyl Methyl Ether	107-30-2	1.4E+02		9.5E+00	8.9E+00					
3.0E-01	P			3.0E-03	P	1.0E-05	X			1.4E+09			0.1	Chloronitrobenzene, o-	88-73-3	1.1E+03	2.6E+03		7.7E+02	3.5E+03	8.3E+03	6.0E+04	2.4E+03	
6.0E-02	P			7.0E-04	P	2.0E-03	P			1.4E+09			0.1	Chloronitrobenzene, p-	100-00-5	5.5E+03	1.3E+04		3.8E+03	8.2E+02	1.9E+03	1.2E+07	5.7E+02	
										5.0E-03			0.1	Chlorophenol, 2-	95-57-8					5.8E+03			5.8E+03	
										4.0E-04			0.1	Chloropicrin	76-06-2							8.2E+00	8.2E+00	
3.1E-03	C	8.9E-07	C	1.5E-02	I					1.4E+09			0.1	Chlorothalonil	1897-45-6	1.1E+05	2.5E+05	1.9E+09	7.4E+04	1.8E+04	4.1E+04		1.2E+04	
										2.0E-02			0.1	Chlorotoluene, o-	95-49-8					2.3E+04			2.3E+04	
										2.0E-02			0.1	Chlorotoluene, p-	106-43-4					2.3E+04			2.3E+04	
2.4E+02	C	6.9E-02	C							1.4E+09			0.1	Chlorozotocin	54749-90-5	1.4E+00	3.2E+00	2.4E+04	9.6E-01					
										2.0E-01			0.1	Chlorpropham	101-21-3					2.3E+05	5.5E+05		1.6E+05	
										1.0E-03			0.1	Chlorpyrifos	2921-88-2					1.2E+03	2.8E+03		8.2E+02	
										1.0E-02			0.1	Chlorpyrifos Methyl	5598-13-0					1.2E+04	2.8E+04		8.2E+03	
										5.0E-02			0.1	Chlorsulfuron	64902-72-3					5.8E+04	1.4E+05		4.1E+04	
										1.0E-02			0.1	Chlorthal-dimethyl	1861-32-1					1.2E+04	2.8E+04		8.2E+03	
										8.0E-04			0.1	Chlorthiophos	60238-56-4					9.3E+02	2.2E+03		6.6E+02	
										1.5E+00			0.013	Chromium(III), Insoluble Salts	16065-83-1					1.8E+06			1.8E+06	
5.0E-01	J	8.4E-02	S	3.0E-03	I	1.0E-04	I	M		1.4E+09			0.025	Chromium(VI)	18540-29-9	6.5E+02		2.0E+04	6.3E+02	3.5E+03		6.0E+05	3.5E+03	
										1.4E+09			0.013	Chromium, Total	7440-7-3									
										1.3E-02			0.1	Clofentazine	74115-24-5					1.5E+04	3.6E+04		1.1E+04	
										3.0E-04			0.1	Cobalt	7440-48-4					3.5E+02		3.6E+04	3.5E+02	
										6.2E-04			0.1	Coke Oven Emissions	8007-5-2									
										4.0E-02			0.1	Copper	7440-50-8					4.7E+04			4.7E+04	
										5.0E-02			0.1	Cresol, m-	108-39-4					5.8E+04	1.4E+05	3.6E+09	4.1E+04	
										5.0E-02			0.1	Cresol, o-	95-48-7					5.8E+04	1.4E+05	3.6E+09	4.1E+04	
										1.0E-01			0.1	Cresol, p-	106-44-5				1.2E+05	2.8E+05	3.6E+09	8.2E+04		
										1.0E-01			0.1	Cresol, p-chloro-m-	59-50-7				1.2E+05	2.8E+05			8.2E+04	
										1.0E-01			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					1.7E+04	1.4E+09	1.9E+04	1	Crotonaldehyde, trans-	123-73-9	1.7E+02			1.7E+02	1.2E+03			1.2E+03	
										1.0E-01			0.1	Cumene	98-82-8					1.2E+05		1.1E+04	9.9E+03	
2.2E-01	C	6.3E-05	C							1.4E+09			0.1	Cupferron	135-20-6	1.5E+03	3.5E+03	2.6E+07	1.0E+03					
8.4E-01	H			2.0E-03	H					1.4E+09			0.1	Cyanazine	21725-46-2	3.9E+02	9.2E+02		2.7E+02	2.3E+03	5.5E+03			1.6E+03
										1.0E-03			0.1	Cyanides										
										5.0E-03			0.1	Calcium Cyanide	592-01-8					1.2E+03			1.2E+03	
										6.0E-04			0.1	Copper Cyanide	544-92-3					5.8E+03			5.8E+03	
										9.5E+05	5.3E+04	1	0.1	Cyanide (CN-)	57-12-5					7.0E+02		1.9E+02	1.5E+02	
										1.0E-03			0.1	Cyanogen	460-19-5					1.2E+03			1.2E+03	
										9.0E-02			0.1	Cyanogen Bromide	506-68-3					1.1E+05			1.1E+05	
										5.0E-02			0.1	Cyanogen Chloride	506-77-4					5.8E+04			5.8E+04	
										6.0E-04			0.1	Hydrogen Cyanide	74-90-8					7.0E+02		1.8E+02	1.5E+02	
										2.0E-03			0.1	Potassium Cyanide	151-50-8					2.3E+03			2.3E+03	
										5.0E-03			0.04	Potassium Silver Cyanide	506-61-6					5.8E+03			5.8E+03	
										1.0E-01			0.04	Silver Cyanide	506-64-9					1.2E+05			1.2E+05	
										1.0E-03			0.1	Sodium Cyanide	143-33-9					1.2E+03			1.2E+03	
										2.0E-04			0.1	Thiocyanates	NA					2.3E+02			2.3E+02	

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)															Toxicity and Chemical-specific Information				Contaminant				Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k _e (ug/m ³) ⁻¹	IUR (ug/m ³) ⁻¹	k _e (mg/kg-day)	RfD _o (mg/kg-day)	k _e (mg/m ³) ⁻¹	RfC _i (mg/m ³) ⁻¹	k _e (mg/m ³) ⁻¹	v _o (mg/m ³) ⁻¹	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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-04	X				V		1.4E+09				1	*Thiocyanic Acid	463-56-9					2.3E+02			2.3E+02						
			5.0E-02	I						1.4E+09				1	*Zinc Cyanide	557-21-1					5.8E+04			5.8E+04						
			6.0E+00	I	V					1.2E+02	1.4E+09	1.0E+03		1	Cyclohexane	110-82-7							2.7E+04	2.7E+04						
2.3E-02	H									1.4E+09				0.1	Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	1.4E+04	3.4E+04		1.0E+04										
			5.0E+00	I	7.0E-01	P	V			5.1E+03	1.4E+09	4.2E+04		1	Cyclohexanone	108-94-1					5.8E+06		1.3E+05	1.3E+05						
			2.0E-01	I				V		2.8E+02	1.4E+09	1.5E+03		1	Cyclohexene	110-83-8					5.8E+03		6.4E+03	3.1E+03						
			2.0E-01	I						2.9E+05	1.4E+09	7.5E+04		1	Cyclohexylamine	108-91-8					2.3E+05			2.3E+05						
			2.5E-02	I						1.4E+09				0.1	Cyfluthrin	68359-37-5					2.9E+04	6.9E+04		2.1E+04						
			5.0E-03	I						1.4E+09				0.1	Cyhalothrin	68085-85-8					5.8E+03	1.4E+04		4.1E+03						
			1.0E-02	I						1.4E+09				0.1	Cypermethrin	52315-07-8					1.2E+04	2.8E+04		8.2E+03						
			7.5E-03	I						1.4E+09				0.1	Cyromazine	66215-27-8					8.8E+03	2.1E+04		6.2E+03						
2.4E-01	I	6.9E-05	C							1.4E+09				0.1	DDD	72-54-8	1.4E+03	3.2E+03	2.4E+07	9.6E+02										
3.4E-01	I	9.7E-05	C					V		1.4E+09	2.1E+06			1	DDE, p,p'-	72-55-9	9.6E+02		2.7E+04	9.3E+02										
3.4E-01	I	9.7E-05	I	5.0E-04	I					1.4E+09				0.03	DDT	50-29-3	9.6E+02	7.6E+03	1.7E+07	8.5E+02	5.8E+02	4.6E+03		5.2E+02						
			3.0E-02	I						1.4E+09				1	Dalapon	75-99-0					3.5E+04	8.3E+04		2.5E+04						
1.8E-02	C	5.1E-06	C	1.5E-01	I					1.4E+09				0.1	Daminozide	1596-84-5	1.8E+04	4.3E+04	3.3E+08	1.3E+04	1.8E+05	4.1E+05		1.2E+05						
7.0E-04	I		7.0E-03	I						1.4E+09				0.1	Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'-(BDE-209)	1163-19-5	4.7E+05	1.1E+06		3.3E+05	8.2E+03	1.9E+04		5.7E+03						
			4.0E-05	I						1.4E+09				0.1	Demeton	8065-48-3					4.7E+01	1.1E+02		3.3E+01						
1.2E-03	I		6.0E-01	I						1.4E+09				0.1	Di(2-ethylhexyl)adipate	103-23-1	2.7E+05	6.4E+05		1.9E+05	7.0E+05	1.7E+06		4.9E+05						
6.1E-02	H		7.0E-04	A						1.4E+09				0.1	Diallate	2303-16-4	5.4E+03	1.3E+04		3.8E+03										
			1.0E-02	X				V		1.4E+09	5.2E+05			1	Dibenzothiophene	132-65-0					1.2E+04			1.2E+04						
8.0E-01	P	6.0E-03	P	2.0E-04	I	V	M			9.8E+02	1.4E+09	3.2E+04		1	Dibromo-3-chloropropane, 1,2-	96-12-8	4.1E+02		6.5E+00	6.4E+00	2.3E+02		2.8E+01	2.5E+01						
			4.0E-04	X				V		1.6E+02	1.4E+09	1.9E+04		1	Dibromobenzene, 1,3-	108-36-1					4.7E+02			4.7E+02						
			1.0E-02	I				V		1.4E+09	2.2E+04			1	Dibromobenzene, 1,4-	106-37-6					1.2E+04			1.2E+04						
8.4E-02	I		2.0E-02	I				V		8.0E+02	1.4E+09	8.0E+03		1	Dibromochloromethane	124-48-1	3.9E+03			3.9E+03	2.3E+04			2.3E+04						
2.0E+00	I	6.0E-04	I	9.0E-03	I	9.0E-03	I	V		1.3E+03	1.4E+09	8.6E+03		1	Dibromoethane, 1,2-	106-93-4	1.6E+02		1.8E+01	1.6E+01	1.1E+04		3.4E+02	3.3E+02						
			4.0E-03	X	V					2.8E+03	1.4E+09	5.6E+03		1	Dibromomethane (Methylene Bromide)	74-95-3							9.9E+01							
			3.0E-04	P						1.4E+09				0.1	Dibutyltin Compounds	NA					3.5E+02	8.3E+02		2.5E+02						
			3.0E-02	I						1.4E+09				0.1	Dicamba	1918-00-9					3.5E+04	8.3E+04		2.5E+04						
4.2E-03	P							V		5.5E+02	1.4E+09	3.2E+03		1	Dichloro-2-butene, 1,4-	764-41-0			9.4E-01	9.4E-01										
4.2E-03	P							V		5.2E+02	1.4E+09	1.1E+04		1	Dichloro-2-butene, cis-1,4-	1476-11-5			3.2E+00	3.2E+00										
4.2E-03	P							V		7.6E+02	1.4E+09	1.1E+04		1	Dichloro-2-butene, trans-1,4-	110-57-5			3.2E+00	3.2E+00										
5.0E-02	I		4.0E-03	I						1.4E+09				0.1	Dichloroacetic Acid	79-43-6	6.5E+03	1.5E+04		4.6E+03	4.7E+03	1.1E+04		3.3E+03						
			9.0E-02	I	2.0E-01	H	V			3.8E+02	1.4E+09	1.2E+04		1	Dichlorobenzene, 1,2-	95-50-1					1.1E+05		1.0E+04	9.3E+03						
5.4E-03	C	1.1E-05	C	7.0E-02	A	8.0E-01	I	V		1.4E+09	1.0E+04			1	Dichlorobenzene, 1,4-	106-46-7	6.1E+04		1.2E+03	1.1E+03	8.2E+04		3.7E+04	2.5E+04						
4.5E-01	I	3.4E-04	C							1.4E+09				0.1	Dichlorobenzidine, 3,3'-	91-94-1	7.3E+02	1.7E+03	4.9E+06	5.1E+02										
			9.0E-03	X						1.4E+09				0.1	Dichlorobenzophenone, 4,4'-	90-98-2					1.1E+04	2.5E+04		7.4E+03						
			2.0E-01	I	1.0E-01	X	V			8.5E+02	1.4E+09	8.4E+02		1	Dichlorodifluoromethane	75-71-8					2.3E+05		3.7E+02	3.7E+02						
5.7E-03	C	1.6E-06	C	2.0E-01	P			V		1.7E+03	1.4E+09	2.1E+03		1	Dichloroethane, 1,1-	75-34-3	5.7E+04		1.6E+03	1.6E+03	2.3E+05			2.3E+05						
9.1E-02	I	2.6E-05	I	6.0E-03	X	7.0E-03	P	V		3.0E+03	1.4E+09	4.6E+03		1	Dichloroethane, 1,2-	107-06-2	3.6E+03		2.2E+02	2.0E+02	7.0E+03		1.4E+02	1.4E+02						
			5.0E-02	I	2.0E-01	I	V			1.2E+03	1.4E+09	1.2E+03		1	Dichloroethylene, 1,1-	75-35-4					5.8E+04		1.0E+03	1.0E+03						
2.0E-03	I							V		2.4E+03	1.4E+09	2.5E+03		1	Dichloroethylene, 1,2-cis-	156-59-2					2.3E+03		2.3E+03							
2.0E-02	I							V		1.9E+03	1.4E+09	1.8E+03		1	Dichloroethylene, 1,2-trans-	156-60-5					2.3E+04		2.3E+04							
3.0E-03	I									1.4E+09				0.1	Dichlorophenol, 2,4-	120-83-2					3.5E+03	8.3E+03		2.5E+03						
1.0E-02	I									1.4E+09				0.05	Dichlorophenoxy Acetic Acid, 2,4-	94-75-7					1.2E+04	5.5E+04		9.6E+03						
8.0E-03	I									1.4E+09				0.1	Dichlorophenoxybutyric Acid, 4-(2,4-	94-82-6					9.3E+03	2.2E+04		6.6E+03						
3.6E-02	C	1.0E-05	C	9.0E-02	A	4.0E-03	I	V		1.4E+03	1.4E+09	3.8E+03		1	Dichloropropane, 1,2-	78-87-5	9.1E+03		4.6E+02	4.4E+02	1.1E+05		6.6E+01	6.6E+01						
			2.0E-02	P				V		1.5E+03	1.4E+09	6.8E+03		1	Dichloropropane, 1,3-	142-28-9					2.3E+04			2.3E+04						
			3.0E-03	I						1.4E+09				0.1	Dichloropropanol, 2,3-	616-23-9					3.5E+03	8.3E+03		2.5E+03						
1.0E-01	I	4.0E-06	I	3.0E-02	I	2.0E-02	I	V		1.6E+03	1.4E+09	3.6E+03		1	Dichloropropene, 1,3-	542-75-6	3.3E+03		1.1E+03	8.2E+02	3.5E+04		3.1E+02	3.1E+02						
2.9E-01	I	8.3E-05	C	5.0E-04	I	5.0E-04	I			1.4E+09				0.1	Dichlorvos	62-73-7	1.1E+03	2.7E+03	2.0E+07	7.9E+02	5.8E+02	1.4E+03	3.0E+06		4.1E+02					
			1.0E-04	I						1.4E+09				0.1	Dicrotophos	141-66-2					1.2E+02	2.8E+02		8.2E+01						
			8.0E-02	P	3.0E-04	X	V			2.6E+02	1.4E+09	4.1E+03		1	Dicyclopentadiene	77-73-6					9.3E+04		5.4E+00	5.4E+00						
1.6E+01	I	4.6E-03	I	5.0E-05	I					1.4E+09				0.1	Dieldrin	60-57-1	2.0E+01	4.8E+01	3.6E+05											

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)														Toxicity and Chemical-specific Information				Contaminant				Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³) ⁻¹	k _e (y)	Rfd _o (mg/kg-day)	k _e (y)	RfC _i (mg/m ³)	k _e (y)	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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)						
3.5E+02	C	1.0E-01	C	1.0E-03	P	V			1.1E+05	1.4E+09	1.4E+05	1	0.1	Diethylformamide	617-84-5					1.2E+03			1.2E+03						
										1.4E+09			0.1	Diethylstilbestrol	56-53-1	9.3E-01	2.2E+00	1.7E+04	6.6E-01										
				8.0E-02	I					1.4E+09			0.1	Difenzoquat	43222-48-6					9.3E+04	2.2E+05		6.6E+04						
				2.0E-02	I					1.4E+09			0.1	Diflubenuron	35367-38-5					2.3E+04	5.5E+04		1.6E+04						
4.4E-02	C	1.3E-05	C						4.0E+01	I	V			Diffluoroethane, 1,1-Dihydrosafrole	75-37-6 94-58-6	7.4E+03		1.2E+04	4.5E+03			2.0E+05	2.0E+05						
														Diisopropyl Ether	108-20-3							9.4E+03	9.4E+03						
				8.0E-02	I					5.3E+02	1.4E+09	3.8E+04	1	Diisopropyl Methylphosphonate	1445-75-6					9.3E+04			9.3E+04						
				2.0E-02	I					1.4E+09			0.1	Dimethipin	55290-64-7					2.3E+04	5.5E+04		1.6E+04						
				2.0E-04	I					1.4E+09			0.1	Dimethoate	60-51-5					2.3E+02	5.5E+02		1.6E+02						
1.6E+00	P									1.4E+09			0.1	Dimethoxybenzidine, 3,3'-	119-90-4	2.0E+02	4.8E+02		1.4E+02										
1.7E-03	P			6.0E-02	P					1.4E+09			0.1	Dimethyl methylphosphonate	756-79-6	1.9E+05	4.5E+05		1.4E+05	7.0E+04	1.7E+05		4.9E+04						
4.6E+00	C	1.3E-03	C							1.4E+09			0.1	Dimethylamino azobenzene [p-]	60-11-7	7.1E+01	1.7E+02	1.3E+06	5.0E+01										
5.8E-01	H									1.4E+09			0.1	Dimethylaniline HCl, 2,4-	21436-96-4	5.6E+02	1.3E+03		4.0E+02	2.3E+03	5.5E+03		1.6E+03						
2.0E-01	P			2.0E-03	X					1.4E+09			0.1	Dimethylaniline, 2,4-	95-68-1	1.6E+03	3.9E+03		1.1E+03	2.3E+03			2.3E+03						
				2.0E-03	I				8.3E+02	1.4E+09	3.1E+04	1		Dimethylaniline, N,N-Dimethylbenzidine, 3,3'-	121-69-7														
1.1E+01	P									1.4E+09			0.1	Dimethylhydrazine, 1,1-	119-93-7	3.0E+01	7.0E+01		2.1E+01	1.2E+05		1.7E+04	1.5E+04						
				1.0E-01	P	3.0E-02	I	V		1.1E+05	1.4E+09	1.3E+05	1	Dimethylformamide	68-12-2					1.2E+02		2.4E-01	2.4E-01						
				1.0E-04	X	2.0E-06	X	V		1.7E+05	1.4E+09	2.8E+04	1	Dimethylhydrazine, 1,2-	57-14-7														
5.5E+02	C	1.6E-01	C							1.9E+05	1.4E+09	1.7E+05	1	Dimethylphenol, 2,4-	540-73-8	5.9E-01		1.3E+00	4.1E-01	2.3E+04	5.5E+04		1.6E+04						
				2.0E-02	I					1.4E+09			0.1	Dimethylphenol, 2,6-	105-67-9					7.0E+02	1.7E+03		4.9E+02						
				6.0E-04	I					1.4E+09			0.1	Dimethylphenol, 3,4-	95-65-8					1.2E+03	2.8E+03		8.2E+02						
4.5E-02	C	1.3E-05	C							4.7E+02	1.4E+09	5.5E+03	1	Dimethylvinylchloride	513-37-1	7.3E+03		5.2E+02	4.8E+02	9.3E+01	2.2E+02		6.6E+01						
				8.0E-05	X					1.4E+09			0.1	Dinitro-o-cresol, 4,6-	534-52-1														
				2.0E-03	I					1.4E+09			0.1	Dinitro-o-cyclohexyl Phenol, 4,6-	131-89-5					2.3E+03	5.5E+03		1.6E+03						
				1.0E-04	P					1.4E+09			0.1	Dinitrobenzene, 1,2-	528-29-0					1.2E+02	2.8E+02		8.2E+01						
				1.0E-04	I					1.4E+09			0.1	Dinitrobenzene, 1,3-	99-65-0					1.2E+02	2.8E+02		8.2E+01						
6.8E-01	I									1.4E+09			0.1	Dinitrobenzene, 1,4-	100-25-4					1.2E+02	2.8E+02		8.2E+01						
				2.0E-03	I					1.4E+09			0.1	Dinitrophenol, 2,4-	51-28-5					2.3E+03	5.5E+03		1.6E+03						
										1.4E+09			0.1	Dinitrotoluene Mixture, 2,4/2,6-	NA	4.8E+02	1.1E+03		3.4E+02										
3.1E-01	C	8.9E-05	C	2.0E-03	I					1.4E+09			0.102	Dinitrotoluene, 2,4-	121-14-1	1.1E+03	2.4E+03	1.9E+07	7.4E+02	2.3E+03	5.4E+03		1.6E+03						
1.5E+00	P			3.0E-04	X					1.4E+09			0.099	Dinitrotoluene, 2,6-	606-20-1	2.2E+02	5.2E+02		1.5E+02	3.5E+02	8.4E+02		2.5E+02						
				2.0E-03	S					1.4E+09			0.006	Dinitrotoluene, 2-Amino-4,6-	35572-8-2					2.3E+03	9.2E+04		2.3E+03						
				2.0E-03	S					1.4E+09			0.009	Dinitrotoluene, 4-Amino-2,6-	19406-51-0					2.3E+03	6.1E+04		2.3E+03						
4.5E-01	X			9.0E-04	X					1.4E+09			0.1	Dinitrotoluene, Technical grade	25321-34-6	7.3E+02	1.7E+03		5.1E+02	1.1E+03	2.5E+03		7.4E+02						
				1.0E-03	I					1.4E+09			0.1	Dinoseb	89-85-7					1.2E+03	2.8E+03		8.2E+02						
1.0E-01	I	5.0E-06	I	3.0E-02	I	3.0E-02	I	V		1.2E+05	1.4E+09	4.0E+04	1	Dioxane, 1,4-	123-91-1	3.3E+03		9.7E+03	2.4E+03	3.5E+04		5.2E+03	4.5E+03						
6.2E+03	I	1.3E+00	I							1.4E+09			0.03	Dioxins	NA	5.3E-02	4.2E-01	1.3E+03	4.7E-02										
1.3E+05	C	3.8E+01	C	7.0E-10	I	4.0E-08	C	V		1.4E+09	2.0E+06	1	0.03	1,2,3,7,8-PeCDD, 2,3,7,8-	1746-01-6	2.5E-03	2.0E-02	6.3E-02	2.2E-03	8.2E-04	6.4E-03	3.4E-01	7.2E-04						
				3.0E-02	I					1.4E+09			0.1	Diphenamid	957-51-7					3.5E+04	8.3E+04		2.5E+04						
				8.0E-04	X					1.4E+09			0.1	Diphenyl Sulfone	127-63-9					9.3E+02	2.2E+03		6.6E+02						
				2.5E-02	I					1.4E+09			0.1	Diphenylamine	122-39-4					2.9E+04	6.9E+04		2.1E+04						
8.0E-01	I	2.2E-04	I							1.4E+09			0.1	Diphenylhydrazine, 1,2-	122-66-7	4.1E+02	9.7E+02	7.6E+06	2.9E+02	2.6E+03	6.1E+03		1.8E+03						
				2.2E-03	I					1.4E+09			0.1	Diquat	85-00-7														
7.1E+00	C	1.4E-01	C							1.4E+09			0.1	Direct Black 38	1937-37-7	4.6E+01	1.1E+02	1.2E+04	3.2E+01										
7.4E+00	C	1.4E-01	C							1.4E+09			0.1	Direct Blue 6	2602-46-2	4.4E+01	1.0E+02	1.2E+04	3.1E+01										
6.7E+00	C	1.4E-01	C							1.4E+09			0.1	Direct Brown 95	16071-86-6	4.9E+01	1.2E+02	1.2E+04	3.4E+01										
				4.0E-05	I					1.4E+09			0.1	Disulfoton	298-04-4					4.7E+01	1.1E+02		3.3E+01						
				1.0E-02	I					1.4E+09	4.5E+04	1		Dithiane, 1,4-	505-29-3					1.2E+04									
				2.0E-03	I					1.4E+09			0.1	Diuron	330-54-1					2.3E+03	5.5E+03		1.6E+03						
				4.0E-03	I					1.4E+09			0.1	Dodine	2439-10-3					4.7E+03	1.1E+04		3.3E+03						
				2.5E-02	I					1.4E+09	1.2E+05	1		EPTC	759-94-4					2.9E+04			2.9E+04						
				6.0E-03	I					1.4E+09	4.1E+05	1		Endosulfan	115-29-7					7.0E+03			7.0E+03						
				2.0E-02	I					1.4E+09			0.1	Endothall	145-73-3					2.3E+04	5.5E+04		1.6E+04						
				3.0E-04	I					1.4E+09			0.1	Endrin	72-20-8					3.5E+02	8.3E+02		2.5E+02						
9.9E-03	I	1.2E-06	I	6.0E-03	P	1.0E-03	I	V		1.1E+04	1.4E+09	1.9E+04	1	Epichlorohydrin	106-89-8	3.3E+04		1.9E+04	1.2E+04	7.0E+03		8.3E+01	8.2E+01						
				2.0E-02	I					1.5E+04	1.4E+09	7.7E+03	1	Epoxybutane, 1,2-	106-88-7							6.7E+02	6.7E+02						

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)																								
Toxicity and Chemical-specific Information															Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k e	IUR (ug/m ³) ⁻¹	k e	RfD _o (mg/kg- day)	k e	RfC _i (mg/m ³) ⁻¹	k e	v	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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)
				4.0E-02	P					1.4E+09		1	0.1		Ethanol, 2-(2-methoxyethoxy)-	111-77-3					4.7E+04	1.1E+05		3.3E+04
				5.0E-03	I					1.4E+09		1	0.1		Ethephon	16672-87-0					5.8E+03	1.4E+04		4.1E+03
				5.0E-04	I					1.4E+09		1	0.1		Ethion	563-12-2					5.8E+02	1.4E+03		4.1E+02
				1.0E-01	P	6.0E-02	P	V		2.4E+04	1.4E+09	6.2E+04	1		Ethoxyethanol Acetate, 2-	111-15-9					1.2E+05		1.6E+04	1.4E+04
				9.0E-02	P	2.0E-01	I	V		1.1E+05	1.4E+09	9.8E+04	1		Ethoxyethanol, 2-	110-80-5					1.1E+05		8.6E+04	4.7E+04
				9.0E-01	I	7.0E-02	P	V		1.1E+04	1.4E+09	8.6E+03	1		Ethyl Acetate	141-78-6					1.1E+06		2.6E+03	2.6E+03
				5.0E-03	P	8.0E-03	P	V		2.5E+03	1.4E+09	6.3E+03	1		Ethyl Acrylate	140-88-5					5.8E+03		2.2E+02	2.1E+02
						1.0E+01	I	V		2.1E+03	1.4E+09	1.3E+03	1		Ethyl Chloride (Chloroethane)	75-00-3							5.7E+04	5.7E+04
				2.0E-01	I			V		1.0E+04	1.4E+09	3.1E+03	1		Ethyl Ether	60-29-7					2.3E+05			2.3E+05
						3.0E-01	P	V		1.1E+03	1.4E+09	5.8E+03	1		Ethyl Methacrylate	97-63-2							7.6E+03	7.6E+03
				1.0E-05	I					1.4E+09			0.1		Ethyl-p-nitrophenyl Phosphonate	2104-64-5					1.2E+01	2.8E+01		8.2E+00
1.1E-02	C	2.5E-06	C	1.0E-01	I	1.0E+00	I	V		4.8E+02	1.4E+09	5.7E+03	1		Ethylbenzene	100-41-4	3.0E+04		2.8E+03	2.5E+03	1.2E+05		2.5E+04	2.0E+04
				7.0E-02	P					1.4E+09			0.1		Ethylene Cyanohydrin	109-78-4					8.2E+04	1.9E+05		5.7E+04
				9.0E-02	P			V		1.9E+05	1.4E+09	1.8E+05	1		Ethylene Diamine	107-15-3					1.1E+05			1.1E+05
				2.0E+00	I	4.0E-01	C			1.4E+09			0.1		Ethylene Glycol	107-21-1					2.3E+06	5.5E+06	2.4E+09	1.6E+06
				1.0E-01	I	1.6E+00	I			1.4E+09			0.1		Ethylene Glycol Monobutyl Ether	111-76-2					1.2E+05	2.8E+05	9.5E+09	8.2E+04
3.1E-01	C	8.8E-05	C			3.0E-02	C	V		1.2E+05	1.4E+09	6.1E+03	1		Ethylene Oxide	75-21-8	1.1E+03		8.5E+01	7.9E+01			8.0E+02	8.0E+02
4.5E-02	C	1.3E-05	C	8.0E-05	I					1.4E+09			0.1		Ethylene Thiourea	96-45-7	7.3E+03	1.7E+04	1.3E+08	5.1E+03	9.3E+01	2.2E+02		6.6E+01
6.5E+01	C	1.9E-02	C					V		1.5E+05	1.4E+09	2.4E+04	1		Ethyleneimine	151-56-4	5.0E+00		1.5E+00	1.2E+00				
				3.0E+00	I					1.4E+09			0.1		Ethylphthalyl Ethyl Glycolate	84-72-0					3.5E+06	8.3E+06		2.5E+06
				2.5E-04	I					1.4E+09			0.1		Fenamiphos	22224-92-6					2.9E+02	6.9E+02		2.1E+02
				2.5E-02	I					1.4E+09			0.1		Fenpropathrin	39515-41-8					2.9E+04	6.9E+04		2.1E+04
				2.5E-02	I					1.4E+09			0.1		Fenvalerate	51630-58-1					2.9E+04	6.9E+04		2.1E+04
				1.3E-02	I					1.4E+09			0.1		Fluometuron	2164-17-2					1.5E+04	3.6E+04		1.1E+04
				4.0E-02	C	1.3E-02	C			1.4E+09			1		Fluoride	16984-48-8					4.7E+04		7.7E+07	4.7E+04
				6.0E-02	I	1.3E-02	C			1.4E+09			1		Fluorine (Soluble Fluoride)	7782-41-4					7.0E+04		7.7E+07	7.0E+04
				8.0E-02	I					1.4E+09			0.1		Fluridone	59756-60-4					9.3E+04	2.2E+05		6.6E+04
				2.0E-02	I					1.4E+09			0.1		Flurprimidol	56425-91-3					2.3E+04	5.5E+04		1.6E+04
				7.0E-04	I					1.4E+09			0.1		Flusilazole	85509-19-9					8.2E+02	1.9E+03		5.7E+02
				6.0E-02	I					1.4E+09			0.1		Flutolanil	66332-96-5					7.0E+04	1.7E+05		4.9E+04
				1.0E-02	I					1.4E+09			0.1		Fluvalinate	69409-94-5					1.2E+04	2.8E+04		8.2E+03
3.5E-03	I			1.0E-01	I					1.4E+09			0.1		Folpet	133-07-3	9.3E+04	2.2E+05		6.6E+04	1.2E+05	2.8E+05		8.2E+04
1.9E-01	I									1.4E+09			0.1		Fomesafen	72178-02-0	1.7E+03	4.1E+03		1.2E+03	2.3E+03	5.5E+03		1.6E+03
				1.3E-05	I	2.0E-01	I	9.8E-03	A	V	4.2E+04	1.4E+09	7.8E+04	1		Fonofos	944-22-9				2.3E+03			1.6E+03
										1.4E+09			0.1		Formaldehyde	50-00-0			7.3E+03	7.3E+03	2.3E+05		3.3E+03	3.3E+03
				9.0E-01	P	3.0E-04	X	V		1.1E+05	1.4E+09	9.3E+04	1		Formic Acid	64-18-6					1.1E+06		1.2E+02	1.2E+02
				3.0E+00	I					1.4E+09			0.1		Fosetyl-AL	39148-24-8					3.5E+06	8.3E+06		2.5E+06
										1.4E+09			0.03		Furans									
				1.0E-03	X			V		1.4E+09	1.6E+05	1	0.03		~Dibenzofuran	132-64-9					1.2E+03	9.2E+03		1.0E+03
				1.0E-03	I			V		6.2E+03	1.4E+09	2.6E+03	1	0.03	~Furan	110-00-9					1.2E+03	9.2E+03		1.0E+03
				9.0E-01	I	2.0E+00	I	V		1.7E+05	1.4E+09	1.2E+04	1	0.03	~Tetrahydrofuran	109-99-9					1.1E+06	8.3E+06	1.0E+05	9.4E+04
3.8E+00	H									1.4E+09			0.1		Furazolidone	67-45-8	8.6E+01	2.0E+02		6.0E+01	3.5E+03		1.1E+04	2.6E+03
				3.0E-03	I	5.0E-02	H	V		1.0E+04	1.4E+09	4.9E+04	1		Furfural	98-01-1								
1.5E+00	C	4.3E-04	C							1.4E+09			0.1		Furium	531-82-8	2.2E+02	5.2E+02	3.9E+06	1.5E+02				
3.0E-02	I	8.6E-06	C							1.4E+09			0.1		Furmecycloz	60568-05-0	1.1E+04	2.6E+04	1.9E+08	7.7E+03				
				4.0E-04	I					1.4E+09			0.1		Glufosinate, Ammonium	77182-82-2					4.7E+02	1.1E+03		3.3E+02
						8.0E-05	C			1.4E+09			0.1		Glutaraldehyde	111-30-8							4.8E+05	4.8E+05
				4.0E-04	I	1.0E-03	H	V		1.1E+05	1.4E+09	8.4E+04	1		Glycidyl	765-34-4					4.7E+02		3.7E+02	2.1E+02
				1.0E-01	I					1.4E+09			0.1		Glyphosate	1071-83-6					1.2E+05	2.8E+05		8.2E+04
				1.0E-02	X			V		1.4E+09	1.5E+05	1			Guanidine	113-00-8					1.2E+04			1.2E+04
				2.0E-02	P					1.4E+09			0.1		Guanidine Chloride	50-01-1					2.3E+04	5.5E+04		1.6E+04
				5.0E-05	I					1.4E+09			0.1		Haloxypol, Methyl	69806-40-2					5.8E+01	1.4E+02		4.1E+01
4.5E+00	I	1.3E-03	I	5.0E-04	I			V		1.4E+09	4.8E+05	1			Heptachlor	76-44-8	7.3E+01		4.5E+02	6.3E+01	5.8E+02			5.8E+02
9.1E+00	I	2.6E-03	I	1.3E-05	I			V		1.4E+09	8.4E+05	1			Heptachlor Epoxide	1024-57-3	3.6E+01		4.0E+02	3.3E+01	1.5E+01			1.5E+01
				2.0E-03	I			V		1.4E+09	3.8E+05	1			Hexabromobenzene	87-82-1					2.3E+03			2.3E+03
				2.0E-04	I					1.4E+09			0.1		Hexabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-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		1.4E+09	6.8E+0													

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)																							
Toxicity and Chemical-specific Information										Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1							
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³) ⁻¹	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _i (mg/m ³)	k _e (y)	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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)
1.8E+00	I	5.3E-04	I	3.0E-04	I					1.4E+09		1	0.1	Hexachlorocyclohexane, Beta-	319-85-7	1.8E+02	4.3E+02	3.1E+06	1.3E+02	3.5E+02	2.1E+03		3.0E+02
1.1E+00	C	3.1E-04	C	3.0E-04	I					1.4E+09		1	0.04	Hexachlorocyclohexane, Gamma- (Lindane)	58-89-9	3.0E+02	1.8E+03	5.4E+06	2.5E+02				
1.8E+00	I	5.1E-04	I							1.4E+09		1	0.1	Hexachlorocyclohexane, Technical	608-73-1	1.8E+02	4.3E+02	3.3E+06	1.3E+02				
4.0E-02	I	1.1E-05	C	6.0E-03	I	2.0E-04	I	V	1.6E+01	1.4E+09	8.5E+03	1		Hexachlorocyclopentadiene	77-47-4					7.0E+03		7.5E+00	7.5E+00
				7.0E-04	I	3.0E-02	I	V		1.4E+09	8.0E+03	1		Hexachloroethane	67-72-1	8.2E+03		8.9E+02	8.0E+02	8.2E+02		1.1E+03	4.6E+02
				3.0E-04	I					1.4E+09		1	0.1	Hexachlorophene	70-30-4					3.5E+02	8.3E+02		2.5E+02
1.1E-01	I			3.0E-03	I					1.4E+09		1	0.015	Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	3.0E+03	4.7E+04		2.8E+03	3.5E+03	5.5E+04		3.3E+03
				1.0E-05	I	V			3.4E+03	1.4E+09	3.0E+05	1		Hexamethylene Diisocyanate, 1,6-	822-06-0							1.3E+01	1.3E+01
				4.0E-04	P					1.4E+09		1	0.1	Hexamethylphosphoramide	680-31-9					4.7E+02	1.1E+03		3.3E+02
				7.0E-01	I	V			1.4E+02	1.4E+09	8.3E+02	1		Hexane, n-	110-54-3							2.5E+03	2.5E+03
				2.0E+00	P					1.4E+09		1	0.1	Hexanedioic Acid	124-04-9					2.3E+06	5.5E+06		1.6E+06
				5.0E-03	I	3.0E-02	I	V	3.3E+03	1.4E+09	1.3E+04	1		Hexanone, 2-	591-78-6					5.8E+03		1.7E+03	1.3E+03
				3.3E-02	I					1.4E+09		1	0.1	Hexazinone	51235-04-2					3.9E+04	9.1E+04		2.7E+04
				2.5E-02	I					1.4E+09		1	0.1	Hexythiazox	78587-05-0					2.9E+04	6.9E+04		2.1E+04
				3.0E-04	I					1.4E+09		1	0.1	Hydramethylnon	67485-29-4					3.5E+02	8.3E+02		2.5E+02
3.0E+00	I	4.9E-03	I			3.0E-05	P	V		1.4E+09		1		Hydrazine	302-01-2	1.1E+02		3.4E+05	1.1E+02			1.8E+05	1.8E+05
3.0E+00	I	4.9E-03	I							1.4E+09		1		Hydrazine Sulfate	10034-93-2	1.1E+02		3.4E+05	1.1E+02				
				2.0E-02	I	V				1.4E+09		1		Hydrogen Chloride	7647-01-0							1.2E+08	1.2E+08
				4.0E-02	C	1.4E-02	C	V		1.4E+09		1		Hydrogen Fluoride	7664-39-3					4.7E+04		8.3E+07	4.7E+04
6.0E-02	P			4.0E-02	P				2.0E-03	1.4E+09		1		Hydrogen Sulfide	7783-06-4							1.2E+07	1.2E+07
				2.5E-01	I					1.4E+09		1	0.1	Hydroquinone	123-31-9	5.5E+03	1.3E+04		3.8E+03	4.7E+04	1.1E+05		3.3E+04
				1.3E-02	I					1.4E+09		1	0.1	Imazalil	35554-44-0					1.5E+04	3.6E+04		1.1E+04
				2.5E-01	I					1.4E+09		1	0.1	Imazaquin	81335-37-7					2.9E+05	6.9E+05		2.1E+05
				2.5E-01	I					1.4E+09		1	0.1	Imazethapyr	81335-77-5					2.9E+05	6.9E+05		2.1E+05
				1.0E-02	A					1.4E+09		1		Iodine	7553-56-2							1.2E+04	1.2E+04
				4.0E-02	I					1.4E+09		1	0.1	Iprodione	36734-19-7					4.7E+04	1.1E+05		3.3E+04
				7.0E-01	P					1.4E+09		1		Iron	7439-89-6					8.2E+05			8.2E+05
9.5E-04	I			3.0E-01	I		V		1.0E+04	1.4E+09	2.8E+04	1		Isobutyl Alcohol	78-83-1				2.4E+05	3.5E+05			3.5E+05
				2.0E-01	I	2.0E+00	C			1.4E+09		1	0.1	Isophorone	78-59-1	3.4E+05	8.1E+05			2.3E+05	5.5E+05	1.2E+10	1.6E+05
				1.5E-02	I		V			1.4E+09	4.2E+05	1		Isopropalin	33820-53-0					1.8E+04			1.8E+04
				2.0E+00	P	2.0E-01	P	V	1.1E+05	1.4E+09	2.8E+04	1		Isopropanol	67-63-0					2.3E+06		2.4E+04	2.4E+04
				1.0E-01	I					1.4E+09		1	0.1	Isopropyl Methyl Phosphonic Acid	1832-54-8					1.2E+05	2.8E+05		8.2E+04
				5.0E-02	I					1.4E+09		1	0.1	Isoxaben	82558-50-7					5.8E+04	1.4E+05		4.1E+04
				3.0E-01	A	V				1.4E+09		1		JP-7	NA							1.8E+09	1.8E+09
				2.0E-03	I					1.4E+09		1	0.1	Lactofen	77501-53-4					2.3E+03	5.5E+03		1.6E+03
														Lead Compounds									
5.0E-01	C	1.5E-01	C	2.0E-02	C	2.0E-04	C		M	1.4E+09	0.025			*Lead Chromate	7758-97-6	6.5E+02		1.1E+04	6.2E+02	2.3E+04		1.2E+06	2.3E+04
8.5E-03	C	1.2E-05	C							1.4E+09		1		*Lead Phosphate	7446-27-7	3.8E+04		1.4E+08	3.8E+04				
8.5E-03	C	1.2E-05	C							1.4E+09		1	0.1	*Lead acetate	301-04-2	3.8E+04	9.1E+04	1.4E+08	2.7E+04				
										1.4E+09		1		*Lead and Compounds	7439-92-1								8.0E+02
8.5E-03	C	1.2E-05	C							1.4E+09		1	0.1	*Lead subacetate	1335-32-6	3.8E+04	9.1E+04	1.4E+08	2.7E+04				
				1.0E-07	I		V		2.4E+00	1.4E+09	1.9E+03	1		*Tetraethyl Lead	78-00-2					1.2E-01			1.2E-01
				5.0E-06	P		V		3.8E+02	1.4E+09	2.6E+04	1		Lewisite	541-25-3					5.8E+00			5.8E+00
				2.0E-03	I					1.4E+09		1	0.1	Linuron	330-55-2					2.3E+03	5.5E+03		1.6E+03
				2.0E-03	P					1.4E+09		1		Lithium	7439-93-2					2.3E+03			2.3E+03
				5.0E-04	I					1.4E+09		1	0.1	MCPA	94-74-6					5.8E+02	1.4E+03		4.1E+02
				1.0E-02	I					1.4E+09		1	0.1	MCPB	94-81-5					1.2E+04	2.8E+04		8.2E+03
				1.0E-03	I					1.4E+09		1	0.1	MCPD	93-65-2					1.2E+03	2.8E+03		8.2E+02
				2.0E-02	I					1.4E+09		1	0.1	Malathion	121-75-5					2.3E+04	5.5E+04		1.6E+04
				1.0E-01	I	7.0E-04	C			1.4E+09		1	0.1	Maleic Anhydride	108-31-6					1.2E+05	2.8E+05	4.2E+06	8.0E+04
				5.0E-01	I					1.4E+09		1	0.1	Maleic Hydrazide	123-33-1					5.8E+05	1.4E+06		4.1E+05
				1.0E-04	P					1.4E+09		1	0.1	Malononitrile	109-77-3					1.2E+02	2.8E+02		8.2E+01
				3.0E-02	H					1.4E+09		1	0.1	Mancozeb	8018-01-7					3.5E+04	8.3E+04		2.5E+04
				5.0E-03	I					1.4E+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		Manganese (Diet)	7439-96-5					2.8E+04		3.0E+05	2.6E+04
				2.4E-02	S	5.0E-05	I			1.4E+09	0.04			Manganese (Non-diet)	7439-96-5					1.1E+02	2.5E+02		7.4E+01
				9.0E-05	H					1.4E+09		1	0.1	Mephsolan	950-10-7								
				3.0E-02	I					1.4E+09		1	0.1	Mepiquat Chloride	24307-26-4					3.5E+04	8.3E+04		2.5E+04
														Mercury Compounds									

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)																									
Toxicity and Chemical-specific Information															Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1				
SFO (mg/kg-day) ⁻¹	k e	IUR (ug/m ³) ⁻¹	k e	RfD _o (mg/kg-day)	k e	RfC _i (mg/m ³)	k e	v	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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	I	3.0E-04	S	1.4E+09	0.07										~Mercuric Chloride (and other Mercury salts)	7487-94-7					3.5E+02		1.8E+06	3.5E+02	
		3.0E-04	I	V	3.1E+00	1.4E+09	3.5E+04	1							~Mercury (elemental)	7439-97-6						4.6E+01		4.6E+01	
		1.0E-04	I			1.4E+09		1							~Methyl Mercury	22967-92-6					1.2E+02			1.2E+02	
		8.0E-05	I			1.4E+09		1	0.1						~Phenylmercuric Acetate	62-38-4					9.3E+01	2.2E+02		6.6E+01	
		3.0E-05	I	V		1.4E+09	1.9E+06	1							Merphos	150-50-5					3.5E+01			3.5E+01	
		3.0E-05	I			1.4E+09		1	0.1						Merphos Oxide	78-48-8					3.5E+01	8.3E+01		2.5E+01	
		6.0E-02	I			1.4E+09		1	0.1						Metalaxyl	57837-19-1					7.0E+04	1.7E+05		4.9E+04	
		1.0E-04	I	3.0E-02	P	V	4.6E+03	1.4E+09	6.8E+03	1					Methacrylonitrile	126-98-7					1.2E+02		8.9E+02	1.0E+02	
		5.0E-05	I			1.4E+09		1	0.1						Methamidophos	10265-92-6					5.8E+01	1.4E+02		4.1E+01	
		2.0E+00	I	2.0E+01	I	V	1.1E+05	1.4E+09	2.9E+04	1					Methanol	67-56-1					2.3E+06		2.5E+06	1.2E+06	
		1.0E-03	I			1.4E+09		1	0.1						Methidathion	950-37-8					1.2E+03	2.8E+03		8.2E+02	
4.9E-02	C	1.4E-05	C			1.4E+09		1	0.1						Methomyl	16752-77-5					2.9E+04	6.9E+04		2.1E+04	
		5.0E-03	I			1.4E+09		1	0.1						Methoxychlor	72-43-5					5.8E+03	1.4E+04		4.1E+03	
		8.0E-03	P	1.0E-03	P	V	1.2E+05	1.4E+09	1.2E+05	1					Methoxyethanol Acetate, 2-	110-49-6					9.3E+03		5.4E+02	5.1E+02	
		5.0E-03	P	2.0E-02	I	V	1.1E+05	1.4E+09	1.0E+05	1					Methoxyethanol, 2-	109-86-4					5.8E+03		8.8E+03	3.5E+03	
		1.0E+00	X		V	2.9E+04	1.4E+09	8.1E+03	1						Methyl Acetate	79-20-9					1.2E+06			1.2E+06	
		2.0E-02	P	V	6.8E+03	1.4E+09	7.0E+03	1							Methyl Acrylate	96-33-3							6.1E+02	6.1E+02	
		6.0E-01	I	5.0E+00	I	V	2.8E+04	1.4E+09	1.2E+04	1					Methyl Ethyl Ketone (2-Butanone)	78-93-3					7.0E+05		2.7E+05	1.9E+05	
		1.0E-03	X	1.0E-03	P	2.0E-05	X	V	1.8E+05	1.4E+09	5.0E+04	1			Methyl Hydrazine	60-34-4			6.2E+01	6.2E+01	1.2E+03		4.4E+00	4.4E+00	
		3.0E+00	I	V	3.4E+03	1.4E+09	1.1E+04	1							Methyl Isobutyl Ketone (4-methyl-2-pentanone)	108-10-1							1.4E+05	1.4E+05	
		1.0E-03	C	V	1.0E+04	1.4E+09	4.4E+03	1							Methyl Isocyanate	624-83-9							1.9E+01	1.9E+01	
		1.4E+00	I	7.0E-01	I	V	2.4E+03	1.4E+09	6.3E+03	1					Methyl Methacrylate	80-62-6					1.6E+06		1.9E+04	1.9E+04	
		2.5E-04	I			1.4E+09		1	0.1						Methyl Parathion	298-00-0					2.9E+02	6.9E+02		2.1E+02	
		6.0E-02	X			1.4E+09		1	0.1						Methyl Phosphonic Acid	993-13-5					7.0E+04	1.7E+05		4.9E+04	
		6.0E-03	H	4.0E-02	H	V	3.9E+02	1.4E+09	2.4E+04	1					Methyl Styrene (Mixed Isomers)	25013-15-4					7.0E+03		4.3E+03	2.6E+03	
9.9E-02	C	2.8E-05	C			1.4E+09		1	0.1						Methyl methanesulfonate	66-27-3	3.3E+03	7.8E+03	6.0E+07	2.3E+03					
1.8E-03	C	2.6E-07	C		3.0E+00	I	V	8.9E+03	1.4E+09	4.9E+03	1				Methyl tert-Butyl Ether (MTBE)	1634-04-4	1.8E+05		2.3E+04	2.1E+04			6.4E+04	6.4E+04	
		3.0E-04	X			1.4E+09		1	0.1						Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2					3.5E+02	8.3E+02		2.5E+02	
		9.0E-03	P	2.0E-02	X		1.4E+09	1	0.1						Methyl-5-Nitroaniline, 2-	99-55-8	3.6E+04	8.6E+04		2.6E+04	2.3E+04	5.5E+04		1.6E+04	
8.3E+00	C	2.4E-03	C			1.4E+09		1	0.1						Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	3.9E+01	9.3E+01	6.9E+05	2.8E+01					
1.3E-01	C	3.7E-05	C			1.4E+09		1	0.1						Methylaniline Hydrochloride, 2-	636-21-5	2.5E+03	5.9E+03	4.5E+07	1.8E+03					
		1.0E-02	A			1.4E+09		1	0.1						Methylarsonic Acid	121-58-3					1.2E+04	2.8E+04		8.2E+03	
		2.0E-04	X			1.4E+09		1	0.1						Methylbenzene, 1,4-diamine hydrochloride, 2-	74612-12-7					2.3E+02	5.5E+02		1.6E+02	
1.0E-01	X	3.0E-04	X			1.4E+09		1	0.1						Methylbenzene-1,4-diamine sulfate, 2-	615-50-9	3.3E+03	7.7E+03		2.3E+03	3.5E+02	8.3E+02		2.5E+02	
2.2E+01	C	6.3E-03	C			1.4E+09		1	0.1						Methylcholanthrene, 3-	56-99-1	1.5E+01	3.5E+01	2.6E+05	1.0E+01					
2.0E-03	I	1.0E-08	I	6.0E-03	I	6.0E-01	I	V	M	3.3E+03	1.4E+09	2.2E+03	1		Methylene Chloride	75-09-1	1.6E+05		2.7E+05	1.0E+05	7.0E+03		5.8E+03	3.2E+03	
1.0E-01	P	4.3E-04	C	2.0E-03	P				M	1.4E+09			1	0.1	Methylene-bis(2-chloroaniline), 4,4'-	101-14-4	3.3E+03	7.7E+03	3.9E+06	2.3E+03	2.3E+03	5.5E+03		1.6E+03	
4.6E-02	I	1.3E-05	C			1.4E+09		1	0.1						Methylene-bis(N,N-dimethyl) Aniline, 4,4'-	101-61-1	7.1E+03	1.7E+04	1.3E+08	5.0E+03					
1.6E+00	C	4.6E-04	C		2.0E-02	C				1.4E+09			1	0.1	Methylenebisbenzenamine, 4,4'-	101-77-9	2.0E+02	4.8E+02	3.6E+06	1.4E+02			1.2E+08	1.2E+08	
				6.0E-04	I			1.4E+09		1			1	0.1	Methylenediphenyl Diisocyanate	101-68-8							3.6E+06	3.6E+06	
		7.0E-02	H		V	5.0E+02	1.4E+09	1.3E+04	1						Methylstyrene, Alpha-	98-83-9					8.2E+04			8.2E+04	
		1.5E-01	I			1.4E+09		1	0.1						Metolachlor	51218-45-2					1.8E+05	4.1E+05		1.2E+05	
		2.5E-02	I			1.4E+09		1	0.1						Metribuzin	21087-64-9					2.9E+04	6.9E+04		2.1E+04	
		2.5E-01	I			1.4E+09		1	0.1						Metsulfuron-methyl	74223-64-6					2.9E+05	6.9E+05		2.1E+05	
		3.0E+00	P		V	3.4E-01	1.4E+09	1.4E+03	1						Mineral oils	8012-95-1					3.5E+06			3.5E+06	
1.8E+01	C	5.1E-03	C	2.0E-04	I	V	1.4E+09	8.6E+05	1						Mirex	2385-85-5	1.8E+01		2.1E+02	1.7E+01	2.3E+02				2.3E+02
		2.0E-03	I			1.4E+09		1	0.1						Molinate	2212-67-1					2.3E+03	5.5E+03		1.6E+03	
		5.0E-03	I			1.4E+09		1							Molybdenum	7439-98-7					5.8E+03			5.8E+03	
		1.0E-01	I			1.4E+09		1							Monochloramine	10599-90-3					1.2E+05			1.2E+05	
		2.0E-03	P			1.4E+09		1	0.1						Monomethylaniline	100-61-8					2.3E+03	5.5E+03		1.6E+03	
		2.5E-02	I			1.4E+09		1	0.1						Myclobutanil	88671-89-0					2.9E+04	6.9E+04		2.1E+04	
		3.0E-04	X			1.4E+09		1	0.1						N,N'-Diphenyl-1,4-benzenediamine	74-31-7					3.5E+02	8.3E+02		2.5E+02	
		2.0E-03	I		V	1.4E+09	5.7E+04	1							Naled	300-76-5					2.3E+03			2.3E+03	
		3.0E-02	X	1.0E-01	P	V	1.4E+09		1						Naphtha, High Flash Aromatic (HFAN)	64742-95-6					3.5E+04		6.0E+08	3.5E+04	
1.8E+00	C	0.0E+00	C			1.4E+09		1	0.1						Naphthylamine, 2-	91-59-8	1.8E+02	4.3E+02		1.3E+02					
		1.0E-01	I			1.4E+09		1	0.1						Napropamide	15299-99-7					1.				

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³) ⁻¹	k _e (y)	Rfd _o (mg/kg-day)	k _e (y)	RfC _i (mg/m ³)	k _e (y)	o _v	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GIABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (mg/kg)	THQ=1 (mg/kg)	Dermal SL THQ=1 (mg/kg)	Inhalation SL THQ=1 (mg/kg)	Noncarcinogenic SL THI=1 (mg/kg)
2.6E-04	C	1.1E-02	C	1.4E-05	C	V				1.4E+09					Nickel Carbonyl	13463-39-3			6.4E+06	6.4E+06	1.3E+04		8.3E+04	1.1E+04
2.6E-04	C	1.1E-02	C	1.4E-05	C					1.4E+09					Nickel Hydroxide	12054-48-7			6.4E+06	6.4E+06	1.3E+04		8.3E+04	1.1E+04
2.6E-04	C	1.1E-02	C	2.0E-05	C					1.4E+09	0.04				Nickel Oxide	1313-99-1			6.4E+06	6.4E+06	1.3E+04		1.2E+05	1.2E+04
2.4E-04	I	1.1E-02	C	1.4E-05	C					1.4E+09	0.04				Nickel Refinery Dust	NA			6.9E+06	6.9E+06	1.3E+04		8.3E+04	1.1E+04
2.6E-04	C	2.0E-02	I	9.0E-05	A					1.4E+09	0.04				Nickel Soluble Salts	7440-02-0			6.4E+06	6.4E+06	2.3E+04		5.4E+05	2.2E+04
1.7E+00	C	4.8E-04	I	1.1E-02	C	1.4E-05	C			1.4E+09	0.04				Nickel Subsulfide	12035-72-2	1.9E+02		3.5E+06	1.9E+02	1.3E+04		8.3E+04	1.1E+04
2.6E-04	C	1.1E-02	C	1.4E-05	C					1.4E+09	1	0.1			Nickelocene	1271-28-9			6.4E+06	6.4E+06	1.3E+04	3.0E+04	8.3E+04	8.1E+03
		1.6E+00	I							1.4E+09	1				Nitrate	14797-55-8					1.9E+06			1.9E+06
										1.4E+09	1				Nitrate + Nitrite (as N)	NA								
		1.0E-01	I							1.4E+09	1				Nitrite	14797-65-0								
		1.0E-02	X	5.0E-05	X					1.4E+09	1	0.1			Nitroaniline, 2-	88-74-4					1.2E+05			1.2E+05
2.0E-02	P			4.0E-03	P	6.0E-03	P			1.4E+09	1	0.1			Nitroaniline, 4-	100-01-6	1.6E+04	3.9E+04		1.1E+04	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	1		Nitrobenzene	98-95-3			2.2E+03	2.2E+03	2.3E+03		2.9E+03	1.3E+03
		3.0E+03	P							1.4E+09	1	0.1			Nitrocellulose	9004-70-0					3.5E+09	8.3E+09		2.5E+09
1.3E+00	C	3.7E-04	C							1.4E+09	1	0.1			Nitrofurantoin	67-20-9					8.2E+04	1.9E+05		5.7E+04
1.7E-02	P		1.0E-04	P						1.4E+09	1	0.1			Nitrofurazone	59-87-0	2.5E+02	5.9E+02	4.5E+06	1.8E+02				
		2.7E-03	H							1.4E+09	1	0.1			Nitroglycerin	55-63-0	1.9E+04	4.5E+04		1.4E+04	1.2E+02	2.8E+02		8.2E+01
		8.8E-06	P		1.0E-01	I				1.4E+09	1	0.1			Nitroguanidine	556-88-7					1.2E+05	2.8E+05		8.2E+04
		2.7E-03	H		5.0E-03	P	V			1.8E+04	1.4E+09	1.7E+04	1		Nitromethane	75-52-5			2.4E+03	2.4E+03			3.7E+02	3.7E+02
		2.7E-03	H		2.0E-02	I	V			4.9E+03	1.4E+09	1.3E+04	1		Nitropropane, 2-	79-46-9			6.0E+00	6.0E+00	1.2E+03		1.2E+03	1.2E+03
2.7E+01	C	7.7E-03	C						M	1.4E+09	1	0.1			Nitroso-N-ethylurea, N-	759-73-9	1.2E+01	2.9E+01	2.2E+05	8.5E+00				
1.2E+02	C	3.4E-02	C						M	1.4E+09	1	0.1			Nitroso-N-methylurea, N-	684-93-5	2.7E+00	6.4E+00	4.9E+04	1.9E+00				
5.4E+00	I	1.6E-03	I						V	1.4E+09	2.4E+05	1			Nitroso-di-N-butylamine, N-	924-16-3	6.1E+01		1.9E+02	4.6E+01				
7.0E+00	I	2.0E-03	C							1.4E+09	1	0.1			Nitroso-di-N-propylamine, N-	621-64-7	4.7E+01	1.1E+02	8.3E+05	3.3E+01				
2.8E+00	I	8.0E-04	C							1.4E+09	1	0.1			Nitrosodiethanolamine, N-	1116-54-7	1.2E+02	2.8E+02	2.1E+06	8.2E+01				
1.5E+02	I	4.3E-02	I						M	1.4E+09	1	0.1			Nitrosodiethylamine, N-	55-18-5	2.2E+00	5.2E+00	3.9E+04	1.5E+00				
5.1E+01	I	1.4E-02	I	8.0E-06	P	4.0E-05	X	V	M	2.4E+05	1.4E+09	8.2E+04	1		Nitrosodimethylamine, N-	62-75-9	6.4E+00		7.2E+00	3.4E+00	9.3E+00		1.4E+01	5.7E+00
4.9E-03	I	2.6E-06	C							1.4E+09	1	0.1			Nitrosodiphenylamine, N-	86-30-6	6.7E+04	1.6E+05	6.4E+08	4.7E+04				
2.2E+01	I	6.3E-03	C						V	1.1E+05	1.4E+09	1.2E+05	1		Nitrosomethylethylamine, N-	10595-95-6	1.5E+01		2.4E+01	9.1E+00				
6.7E+00	C	1.9E-03	C							1.4E+09	1	0.1			Nitrosomorpholine [N-]	59-89-2	4.9E+01	1.2E+02	8.8E+05	3.4E+01				
9.4E+00	C	2.7E-03	C							1.4E+09	1	0.1			Nitrosopiperidine [N-]	100-75-4	3.5E+01	8.2E+01	6.2E+05	2.4E+01				
2.1E+00	I	6.1E-04	I							1.4E+09	1	0.1			Nitrosopyrrolidine, N-	930-55-2	1.6E+02	3.7E+02	2.7E+06	1.1E+02				
2.2E-01	P		1.0E-04	X						1.4E+09	1	0.1			Nitrotoluene, m-	99-08-1					1.2E+02	2.8E+02		8.2E+01
1.6E-02	P		9.0E-04	P					V	1.5E+03	1.4E+09	1.4E+05	1		Nitrotoluene, o-	88-72-7	1.5E+03			1.5E+03	1.1E+03		1.1E+03	
			4.0E-03	P						1.4E+09	1	0.1			Nitrotoluene, p-	99-99-1	2.0E+04	4.8E+04		1.4E+04	4.7E+03	1.1E+04		3.3E+03
		3.0E-04	X	2.0E-02	P	V				6.9E+00	1.4E+09	1.0E+03	1		Nonane, n-	111-84-2					3.5E+02		9.1E+01	7.2E+01
		4.0E-02	I							1.4E+09	1	0.1			Norflurazon	27314-13-2					4.7E+04	1.1E+05		3.3E+04
		3.0E-03	I							1.4E+09	1	0.1			Octabromodiphenyl Ether	32536-52-0					3.5E+03	8.3E+03		2.5E+03
		5.0E-02	I							1.4E+09	1	0.006			Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0					5.8E+04	2.3E+06		5.7E+04
		2.0E-03	H							1.4E+09	1	0.1			Octamethylpyrophosphoramide	152-16-9					2.3E+03	5.5E+03		1.6E+03
		5.0E-02	I							1.4E+09	1	0.1			Oryzalin	19044-88-3					5.8E+04	1.4E+05		4.1E+04
		5.0E-03	I							1.4E+09	1	0.1			Oxadiazon	19666-30-9					5.8E+03	1.4E+04		4.1E+03
		2.5E-02	I							1.4E+09	1	0.1			Oxamyl	23135-22-0					2.9E+04	6.9E+04		2.1E+04
		3.0E-03	I							1.4E+09	1	0.1			Oxyfluorfen	42874-03-3					3.5E+03	8.3E+03		2.5E+03
		1.3E-02	I							1.4E+09	1	0.1			Paclitaxel	76738-62-0					1.5E+04	3.6E+04		1.1E+04
		4.5E-03	I							1.4E+09	1	0.1			Paraquat Dichloride	1910-42-5					5.3E+03	1.2E+04		3.7E+03
		6.0E-03	H							1.4E+09	1	0.1			Parathion	56-38-2					7.0E+03	1.7E+04		4.9E+03
		5.0E-02	H							1.4E+09	4.5E+04	1			Perbulate	1114-71-2					5.8E+04			5.8E+04
		4.0E-02	I							1.4E+09	1	0.1			Pendimethalin	40487-42-1					4.7E+04	1.1E+05		3.3E+04
		2.0E-03	I						V	3.1E-01	1.4E+09	5.1E+05	1		Pentabromodiphenyl Ether	32534-81-9					2.3E+03			2.3E+03
		1.0E-04	I							1.4E+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
9.0E-02	P			8.0E-04	I					1.4E+09	8.1E+04	1			Pentachlorobenzene	608-93-5					9.3E+02			9.3E+02
										4.6E+02	1.4E+09	9.7E+03	1		Pentachloroethane	76-01-7	3.6E+03		3.6E+03					
2.6E-01	H			3.0E-03	I					1.4E+09	4.3E+05	1			Pentachloronitrobenzene	82-68-8	1.3E+03		1.3E+03		3.5E+03			3.5E+03
4.0E-01	I	5.1E-06	C	5.0E-03	I					1.4E+09	1	0.25			Pentachlorophenol	87-86-5	8.2E+02	7.7E+02	3.3E+08	4.0E+02	5.8E+03	5.5E		

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice) ; 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 Guide); s = Concentration may exceed Csat (See User Guide)

Toxicity and Chemical-specific Information													Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k e	IUR (ug/m ³) ⁻¹	k e	RfD _o (mg/kg-day)	k e	RfC _i (mg/m ³) ⁻¹	k e	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-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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)
				7.0E-04	I							1.4E+09				*Lithium Perchlorate	7791-03-9					8.2E+02			8.2E+02
				7.0E-04	I							1.4E+09				*Perchlorate and Perchlorate Salts	14797-73-0					8.2E+02			8.2E+02
				7.0E-04	I							1.4E+09				*Potassium Perchlorate	7778-74-7					8.2E+02			8.2E+02
				7.0E-04	I							1.4E+09				*Sodium Perchlorate	7601-89-0					8.2E+02			8.2E+02
				2.0E-02	P					V		1.4E+09	1.3E+05			Perfluorobutane Sulfonate	375-73-5					2.3E+04			2.3E+04
				5.0E-02	I							1.4E+09		0.1		Permethrin	52645-53-1					5.8E+04	1.4E+05		4.1E+04
2.2E-03	C	6.3E-07	C									1.4E+09			0.1	Phenacetin	62-44-2	1.5E+05	3.5E+05	2.6E+09	1.0E+05				
				2.5E-01	I							1.4E+09			0.1	Phenmedipham	13684-63-4					2.9E+05	6.9E+05		2.1E+05
				3.0E-01	I	2.0E-01	C					1.4E+09			0.1	Phenol	108-95-2					3.5E+05	8.3E+05	1.2E+09	2.5E+05
				4.0E-03	I							1.4E+09			0.1	Phenol, 2-(1-methylethoxy)-, methylcarbamate	114-26-1					4.7E+03	1.1E+04		3.3E+03
				5.0E-04	X							1.4E+09			0.1	Phenothiazine	92-84-2					5.8E+02	1.4E+03		4.1E+02
				6.0E-03	I							1.4E+09			0.1	Phenylenediamine, m-	108-45-2					7.0E+03	1.7E+04		4.9E+03
4.7E-02	H											1.4E+09			0.1	Phenylenediamine, o-	95-54-5	7.0E+03	1.6E+04		4.9E+03				
				1.9E-01	H							1.4E+09			0.1	Phenylenediamine, p-	106-50-3					2.2E+05	5.2E+05		1.6E+05
1.9E-03	H											1.4E+09			0.1	Phenylphenol, 2-	90-43-7				1.2E+05				
				2.0E-04	H							1.4E+09			0.1	Phorate	298-02-2					2.3E+02	5.5E+02		1.6E+02
						3.0E-04	I	V			1.6E+03	1.4E+09	9.8E+02		1	Phosgene	75-44-5							1.3E+00	1.3E+00
				2.0E-02	I							1.4E+09			0.1	Phosmet	732-11-6					2.3E+04	5.5E+04		1.6E+04
																Phosphates, Inorganic									
				4.9E+01	P							1.4E+09			1	*Aluminum metaphosphate	13776-88-0					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Ammonium polyphosphate	68333-79-9					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Calcium pyrophosphate	7790-76-3					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Diammonium phosphate	7783-28-0					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Dicalcium phosphate	7757-93-9					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Dimagnesium phosphate	7782-75-4					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Dipotassium phosphate	7758-11-4					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Disodium phosphate	7558-79-4					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Monoaluminum phosphate	13530-50-2					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Monoammonium phosphate	7722-76-1					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Monocalcium phosphate	7758-23-8					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Monomagnesium phosphate	7757-86-0					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Monopotassium phosphate	7778-77-0					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Monosodium phosphate	7558-80-7					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Polyphosphoric acid	8017-15-1					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Potassium tripolyphosphate	13845-36-8					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Sodium acid pyrophosphate	7758-15-9					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Sodium aluminum phosphate (acidic)	7785-83-8					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Sodium aluminum phosphate (anhydrous)	10279-59-1					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Sodium aluminum phosphate (tetrahydrate)	10305-76-7					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Sodium hexametaphosphate	10124-56-8					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Sodium polyphosphate	68915-31-1					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Sodium trimetaphosphate	7785-84-4					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Sodium tripolyphosphate	7758-29-4					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Tetrapotassium phosphate	7320-34-5					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Tetrasodium pyrophosphate	7722-88-5					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Trialuminum sodium tetra decahydrogenooctaoorthophosphate (dihydrate)	15136-87-5					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Tricalcium phosphate	7758-87-4					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Trimagnesium phosphate	7757-87-1					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Tripotassium phosphate	7778-53-2					5.7E+07			5.7E+07
				4.9E+01	P							1.4E+09			1	*Trisodium phosphate	7601-54-9					5.7E+07			5.7E+07
				3.0E-04	I	3.0E-04	I	V				1.4E+09			1	Phosphine	7803-51-2					3.5E+02		1.8E+06	3.5E+02
				4.9E+01	P	1.0E-02	I					1.4E+09			1	Phosphoric Acid	7664-38-2					5.7E+07		6.0E+07	2.9E+07
				2.0E-05	I							1.4E+09	6.9E+03		1	Phosphorus, White	7723-14-0					2.3E+01			2.3E+01
																Phthalates									
1.4E-02	I	2.4E-06	C	2.0E-02	I							1.4E+09			0.1	*Bis(2-ethylhexyl)phthalate	117-81-7	2.3E+04	5.5E+04	6.9E+08	1.6E+04	2.3E+04	5.5E+04		1.6E+04
1.9E-03	P			2.0E-01	I							1.4E+09			0.1	*Butyl Benzyl Phthalate	85-68-7	1.7E+05	4.1E+05		1.2E+05	2.3E+05	5.5E+05		1.6E+05
				1.0E+00	I							1.4E+09			0.1	*Butylphthalyl Butylglycolate	85-70-1					1.2E+06	2.8E+06		8.2E+05
				1.0E-01	I							1.4E+09			0.1	*Dibutyl Phthalate	84-74-2					1.2E+05	2.8E+05		8.2E+04
				8.0E-01	I							1.4E+09			0.1	*Diethyl Phthalate	84-66-2					9.3E+05	2.2E+06		6.6E+05

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1						
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³) ⁻¹	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _o (mg/m ³)	k _e (y)	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GI/ABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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-01	I					V		1.4E+09	2.1E+04	1		**Dimethylterephthalate	120-61-6								1.2E+05		1.2E+05	
		1.0E-02	P							1.4E+09		1	0.1	**Octyl Phthalate, di-N-	117-84-0								1.2E+04	2.8E+04	8.2E+03	
		1.0E+00	H							1.4E+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					1.4E+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							1.4E+09		1	0.1	Picloram	1918-02-1								8.2E+04	1.9E+05	5.7E+04	
		1.0E-04	X							1.4E+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							1.4E+09		1	0.1	Picric Acid (2,4,6-Trinitrophenol)	88-89-1								1.1E+03	2.5E+03	7.4E+02	
3.0E+01	C	8.6E-03	C	7.0E-06	H					1.4E+09		1	0.1	Pirimiphos, Methyl	29232-93-7								1.2E+04	2.8E+04	8.2E+03	
										1.4E+09		1	0.1	Polybrominated Biphenyls Polychlorinated Biphenyls (PCBs)	59536-65-1	1.1E+01	2.6E+01	1.9E+05	7.7E+00				8.2E+00	1.9E+01	5.7E+00	
7.0E-02	S	2.0E-05	S	7.0E-05	I					1.4E+09	7.1E+05	1	0.14	**Aroclor 1016	12674-11-2	4.7E+03	7.9E+03	4.4E+04	2.7E+03				8.2E+01	1.4E+02	5.1E+01	
2.0E+00	S	5.7E-04	S							1.4E+09	2.0E+05	1	0.14	**Aroclor 1221	11104-28-2	1.6E+02	2.8E+02	4.4E+02	8.3E+01							
2.0E+00	S	5.7E-04	S							1.4E+09	1.1E+05	1	0.14	**Aroclor 1232	11141-16-5	1.6E+02	2.8E+02	2.4E+02	7.2E+01							
2.0E+00	S	5.7E-04	S							1.4E+09	5.9E+05	1	0.14	**Aroclor 1242	53469-21-9	1.6E+02	2.8E+02	1.3E+03	9.5E+01							
2.0E+00	S	5.7E-04	S							1.4E+09	6.3E+05	1	0.14	**Aroclor 1248	12672-29-6	1.6E+02	2.8E+02	1.3E+03	9.5E+01							
2.0E+00	S	5.7E-04	S	2.0E-05	I					1.4E+09	8.4E+05	1	0.14	**Aroclor 1254	11097-69-1	1.6E+02	2.8E+02	1.8E+03	9.7E+01				2.3E+01	3.9E+01	1.5E+01	
2.0E+00	S	5.7E-04	S							1.4E+09	1.3E+06	1	0.14	**Aroclor 1260	11096-82-5	1.6E+02	2.8E+02	2.8E+03	9.9E+01							
3.9E+00	E	1.1E-03	E	6.0E-04	X					1.4E+09	9.6E+05	1	0.14	**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		1.4E+09	2.4E+06	1	0.14	**Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189)	39635-31-9	8.4E+01	1.4E+02	2.6E+03	5.2E+01				2.7E+01	4.6E+01	1.4E+04	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	1.6E+06	1	0.14	**Hexachlorobiphenyl, 2,3,4,4',5,5'- (PCB 167)	52663-72-6	8.4E+01	1.4E+02	1.7E+03	5.1E+01				2.7E+01	4.6E+01	9.2E+03	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	1.0E+06	1	0.14	**Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157)	69782-90-7	8.4E+01	1.4E+02	1.1E+03	5.0E+01				2.7E+01	4.6E+01	6.1E+03	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	1.1E+06	1	0.14	**Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 156)	38380-08-4	8.4E+01	1.4E+02	1.2E+03	5.0E+01				2.7E+01	4.6E+01	6.5E+03	1.7E+01
3.9E+03	E	1.1E+00	E	2.3E-08	E	1.3E-06	E	V		1.4E+09	1.6E+06	1	0.14	**Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169)	32774-16-6	8.4E-02	1.4E-01	1.7E+00	5.1E-02				2.7E-02	4.6E-02	9.2E+00	1.7E-02
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	7.3E+05	1	0.14	**Pentachlorobiphenyl, 2',3,4,4',5'- (PCB 123)	65510-44-3	8.4E+01	1.4E+02	7.9E+02	4.9E+01				2.7E+01	4.6E+01	4.3E+03	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	5.9E+05	1	0.14	**Pentachlorobiphenyl, 2,3',4,4',5'- (PCB 118)	31508-00-6	8.4E+01	1.4E+02	6.3E+02	4.9E+01				2.7E+01	4.6E+01	3.4E+03	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	6.0E+05	1	0.14	**Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105)	32598-14-4	8.4E+01	1.4E+02	6.5E+02	4.9E+01				2.7E+01	4.6E+01	3.5E+03	1.7E+01
3.9E+00	E	1.1E-03	E	2.3E-05	E	1.3E-03	E	V		1.4E+09	1.1E+06	1	0.14	**Tetrachlorobiphenyl, 2,3,4,4',5'- (PCB 114)	74472-37-0	8.4E+01	1.4E+02	1.1E+03	5.0E+01				2.7E+01	4.6E+01	6.1E+03	1.7E+01
1.3E+04	E	3.8E+00	E	7.0E-09	E	4.0E-07	E	V		1.4E+09	7.3E+05	1	0.14	**Pentachlorobiphenyl, 3,3',4,4',5'- (PCB 126)	57465-28-8	2.5E-02	4.2E-02	2.3E-01	1.5E-02				8.2E-03	1.4E-02	1.3E+00	5.1E-03
2.0E+00	I	5.7E-04	I							1.4E+09	5.3E+05	1	0.14	**Polychlorinated Biphenyls (high risk)	1336-36-3	1.6E+02	2.8E+02	1.1E+03	9.4E+01							
4.0E-01	I	1.0E-04	I							1.4E+09		1	0.14	**Polychlorinated Biphenyls (low risk)	1336-36-3											
7.0E-02	I	2.0E-05	I							1.4E+09		1	0.14	**Polychlorinated Biphenyls (lowest risk)	1336-36-3											
1.3E+01	E	3.8E-03	E	7.0E-06	E	4.0E-04	E	V		1.4E+09		1	0.14	**Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77)	32598-13-3	2.5E+01	4.2E+01	4.4E+05	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		1.4E+09	5.1E+05	1	0.14	**Tetrachlorobiphenyl, 3,4,4',5'- (PCB 81)	70362-50-4	8.4E+00	1.4E+01	5.5E+01	4.8E+00				2.7E+00	4.6E+00	3.0E+02	1.7E+00
				6.0E-04	I					1.4E+09		1	0.1	Polymeric Methylenediphenyl Disocyanate (PMDI)	9016-37-9									3.6E+06	3.6E+06	
		6.0E-02	I							1.4E+09	1.4E+05	1	0.13	Polynuclear Aromatic Hydrocarbons (PAHs) **Acenaphthene	83-32-9								7.0E+04	1.3E+05	4.5E+04	
		3.0E-01	I							1.4E+09	5.2E+05	1	0.13	**Anthracene	120-21-7								3.5E+05	6.4E+05	2.3E+05	
7.3E-01	E	1.1E-04	C							1.4E+09	4.4E+06	1	0.13	**Benz[a]anthracene	56-55-3	4.5E+02	8.1E+02	4.9E+04	2.9E+02							
1.2E+00	C	1.1E-04	C							1.4E+09		1	0.13	**Benzo[j]fluoranthene	205-82-3	2.7E+02	5.0E+02	1.5E+07	1.8E+02							
7.3E+00	I	1.1E-03	C							1.4E+09		1	0.13	**Benzo[a]pyrene	50-32-8	4.5E+01	8.1E+01	1.5E+06	2.9E+01							
7.3E-01	E	1.1E-04	C							1.4E+09		1	0.13	**Benzo[b]fluoranthene	205-99-2	4.5E+02	8.1E+02	1.5E+07	2.9E+02							
7.3E-02	E	1.1E-04	C							1.4E+09		1	0.13	**Benzo[k]fluoranthene	207-08-9	4.5E+03	8.1E+03	1.5E+07	2.9E+03							
				8.0E-02	I					1.4E+09	8.0E+04	1	0.13	**Chloronaphthalene, Beta-	91-58-7								9.3E+04	1.7E+05	6.0E+04	
7.3E-03	E	1.1E-05	C							1.4E+09		1	0.13	**Chrysene	218-01-9	4.5E+04	8.1E+04	1.5E+08	2.9E+04							
7.3E+00	E	1.2E-03	C							1.4E+09		1	0.13	**Dibenz[a,h]anthracene	53-70-3	4.5E+01	8.1E+01	1.4E+06	2.9E+01							
1.2E+01	C	1.1E-03	C							1.4E+09		1	0.13	**Dibenzo[a,e]pyrene	192-65-4	2.7E+01	5.0E+01	1.5E+06	1.8E+01							
2.5E+02	C	7.1E-02	C							1.4E+09		1	0.13	**Dimethylbenz[a]anthracene, 7,12-	57-97-6	1.3E+00	2.4E+00	2.3E+04	8.4E-01							
		4.0E-02	I							1.4E+09		1	0.13	**Fluoranthene	206-44-0								4.7E+04	8.5E+04	3.0E+04	
		4.0E-02	I							1.4E+09	2.8E+05	1	0.13	**Fluorene	86-73-7								4.7E+04	8.5E+04	3.0E+04	
7.3E-01	E	1.1E-04	C							1.4E+09		1	0.13	**indeno[1,2,3-cd]pyrene	193-39-5	4.5E+02	8.1E+02	1.5E+07	2.9E+02							
2.9E-02	P			7.0E-02	A					3.9E+02	1.4E+09	5.9E+04	1	0.13	**Methylnaphthalene, 1-	90-12-0	1.1E+04	2.0E+04	1.5E+07	7.3E+03						

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRTV; A = ATSDR; C = Cal EPA; X = APPENDIX PPRTV SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)																								
Toxicity and Chemical-specific Information															Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k _e (y)	IUR (ug/m ³) ⁻¹	k _e (y)	RfD _o (mg/kg-day)	k _e (y)	RfC _i (mg/m ³) ⁻¹	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-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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)	
				4.0E-03	I					1.4E+09			0.1	Prometryn	7287-19-6					4.7E+03	1.1E+04		3.3E+03	
				1.3E-02	I					1.4E+09			0.1	Propachlor	1918-16-7					1.5E+04	3.6E+04		1.1E+04	
				5.0E-03	I					1.4E+09			0.1	Propanil	709-98-8					5.8E+03	1.4E+04		4.1E+03	
				2.0E-02	I					1.4E+09			0.1	Propargite	2312-35-8					2.3E+04	5.5E+04		1.6E+04	
				2.0E-03	I			V	1.1E+05	1.4E+09	6.3E+04		1	Propargyl Alcohol	107-19-7					2.3E+03			2.3E+03	
				2.0E-02	I					1.4E+09			0.1	Propazine	139-40-2					2.3E+04	5.5E+04		1.6E+04	
				2.0E-02	I					1.4E+09			0.1	Propham	122-42-9					2.3E+04	5.5E+04		1.6E+04	
				1.3E-02	I					1.4E+09			0.1	Propiconazole	60207-90-1					1.5E+04	3.6E+04		1.1E+04	
				8.0E-03	I	V			3.3E+04	1.4E+09	8.9E+03		1	Propionaldehyde	123-38-6							3.1E+02	3.1E+02	
				1.0E-01	X	1.0E+00	X	V	2.6E+02	1.4E+09	7.0E+03		1	Propyl benzene	103-65-1					1.2E+05		3.1E+04	2.4E+04	
								C	3.0E+00	C	3.5E+02	1.4E+09	7.0E+02	1	Propylene	115-07-1						9.3E+03	9.3E+03	
				2.0E+01	P					1.4E+09			0.1	Propylene Glycol	57-55-6					2.3E+07	5.5E+07		1.6E+07	
				2.7E-04	A					1.4E+09			0.1	Propylene Glycol Dinitrate	6423-43-4							1.6E+06	1.6E+06	
				7.0E-01	H	2.0E+00	I	V	1.1E+05	1.4E+09	7.8E+04		1	Propylene Glycol Monomethyl Ether	107-98-2					8.2E+05		6.9E+05	3.7E+05	
				2.4E-01	I	3.7E-06	I			3.0E-02	I	V	7.8E+04	1.4E+09	1.0E+04	1			1.4E+03		3.4E+03	9.7E+02	1.4E+03	
				7.5E-02	I					1.4E+09			0.1	Propylene Oxide	75-56-9	1.4E+03		3.4E+03	9.7E+02				1.4E+03	
				1.0E-03	I			V	5.3E+05	1.4E+09	5.5E+04		1	Propylamide	23950-58-5					8.8E+04	2.1E+05		6.2E+04	
										1.4E+09			0.1	Pyridine	110-86-1					1.2E+03			1.2E+03	
				3.0E+00	I					1.4E+09			0.1	Quinalphos	13593-03-8					5.8E+02	1.4E+03		4.1E+02	
				9.0E-03	I					1.4E+09			0.1	Quinoline	91-22-5	1.1E+02	2.6E+02		7.7E+01				1.4E+03	
										1.4E+09			0.1	Quizalofop-ethyl	76578-14-8					1.1E+04	2.5E+04		7.4E+03	
										1.4E+09			1	Refractory Ceramic Fibers	NA							1.8E+08	1.8E+08	
				3.0E-02	I					1.4E+09			0.1	Resmethrin	10453-86-8					3.5E+04	8.3E+04		2.5E+04	
				5.0E-02	H			V	1.4E+09	4.7E+05			1	Ronnel	299-84-3					5.8E+04			5.8E+04	
				4.0E-03	I					1.4E+09			0.1	Rotenone	83-79-4					4.7E+03	1.1E+04		3.3E+03	
				2.2E-01	C	6.3E-05	C			1.4E+09			0.1	Safrole	94-59-7	1.5E+03	3.5E+03	2.6E+07	1.0E+03				5.8E+03	
				5.0E-03	I	2.0E-02	C			1.4E+09			1	Selenious Acid	7783-00-8					5.8E+03			5.8E+03	
				5.0E-03	C	2.0E-02	C			1.4E+09			1	Selenium	7782-49-2					5.8E+03		1.2E+08	5.8E+03	
				9.0E-02	I					1.4E+09			0.1	Selenium Sulfide	7446-34-6					1.1E+05	2.5E+05		7.4E+04	
										1.4E+09			1	Sethoxydim	74051-80-2								1.8E+07	
				3.0E-03	C					1.4E+09			1	Silica (crystalline, respirable)	7631-86-9								1.8E+07	
				5.0E-03	I					1.4E+09	0.04			Silver	7440-22-4					5.8E+03			5.8E+03	
				5.0E-03	I					1.4E+09			0.1	Simazine	1972-34-9	2.7E+03	6.4E+03		1.9E+03			1.4E+04		4.1E+03
				1.3E-02	I					1.4E+09			0.1	Sodium Adifluorfen	62176-59-9					1.5E+04	3.6E+04		1.1E+04	
				4.0E-03	I					1.4E+09			0.1	Sodium Azide	26628-22-8					4.7E+03			4.7E+03	
				2.0E-02	C	2.0E-04	C	M		1.4E+09	0.025			Sodium Dichromate	10588-01-9	6.5E+02		1.1E+04	6.2E+02			1.2E+06	2.3E+04	
				3.0E-02	I					1.4E+09			0.1	Sodium Diethyldithiocarbamate	148-11-5	1.2E+03	2.9E+03		8.5E+02				2.5E+04	
				5.0E-02	A	1.3E-02	C			1.4E+09			1	Sodium Fluoride	7681-49-4					5.8E+04			5.8E+04	
				2.0E-05	I					1.4E+09			0.1	Sodium Fluoroacetate	62-74-8					2.3E+01	5.5E+01		1.6E+01	
				1.0E-03	H					1.4E+09			1	Sodium Metavanadate	13718-26-8					1.2E+03			1.2E+03	
				8.0E-04	P					1.4E+09			1	Sodium Tungstate	13472-45-2					9.3E+02			9.3E+02	
				8.0E-04	P					1.4E+09			1	Sodium Tungstate Dihydrate	10213-10-2					9.3E+02			9.3E+02	
				3.0E-02	I					1.4E+09			0.1	Stirofos (Tetrachlorovinphos)	961-11-5	1.4E+04	3.2E+04		9.6E+03			8.3E+04	2.5E+04	
				2.0E-02	C	2.0E-04	C	M		1.4E+09	0.025			Strontium Chromate	7789-06-2	6.5E+02		1.1E+04	6.2E+02			1.2E+06	2.3E+04	
				6.0E-01	I					1.4E+09			1	Strontium, Stable	7440-24-6					7.0E+05			7.0E+05	
				3.0E-04	I					1.4E+09			0.1	Strychnine	57-24-9					3.5E+02	8.3E+02		2.5E+02	
				2.0E-01	I	1.0E+00	I	V	8.7E+02	1.4E+09	9.4E+03		1	Styrene	100-42-5					2.3E+05		4.1E+04	3.5E+04	
				3.0E-03	P					1.4E+09			0.1	Styrene-Acrylonitrile (SAN) Trimer	NA					3.5E+03	8.3E+03		2.5E+03	
				1.0E-03	P	2.0E-03	X			1.4E+09			0.1	Sulfolane	126-33-0					1.2E+03	2.8E+03	1.2E+07	8.2E+02	
				8.0E-04	P					1.4E+09			0.1	Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9					9.3E+02	2.2E+03		6.6E+02	
										1.4E+09			1	Sulfur Trioxide	7446-11-9							6.0E+06	6.0E+06	
				1.0E-03	C					1.4E+09			1	Sulfuric Acid	7664-93-9							6.0E+06	6.0E+06	
				7.1E-06	I	5.0E-02	H			1.4E+09			0.1	Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethylphenoxy)-1-methylethyl ester	140-57-8	1.3E+04	3.1E+04	2.3E+08	9.2E+03				4.1E+04	
				3.0E-02	H					1.4E+09			0.1	TCMTB	21564-17-0					3.5E+04	8.3E+04		2.5E+04	
				7.0E-02	I					1.4E+09			0.1	Tebuthiuron	34014-18-1					8.2E+04	1.9E+05		5.7E+04	
				2.0E-02	H					1.4E+09			0.1	Temephos	3383-96-8					2.3E+04	5.5E+04		1.6E+04	
				1.3E-02	I					1.4E+09			0.1	Terbacil	5902-51-2					1.5E+04	3.6E+04		1.1E+04	
				2.5E-05	H			V	3.1E+01	1.4E+09	2.6E+05		1	Terbufos	13071-79-9					2.9E+01			2.9E+01	
				1.0E-03	I					1.4E+09			0.1	Terbutryn	886-50-0					1.2E+03	2.8E+03		8.2E+02	
				1.0E-04	I					1.4E+09			0.1	Tetrabromodiphenyl ether, 2,2',4,4'-(BDE-47)	5436-43-1					1.2E+02	2.8E+02		8.2E+01	

Regional Removal Management Level (RML) Composite Worker Soil Table (TR=1E-04, HQ=1) May 2016

Key: I = IRIS; P = PPRVT; A = ATSDR; C = Cal EPA; X = APPENDIX PPRVT SCREEN (See FAQ #27); H = HEAST; F = See FAQ; J = New Jersey; O = EPA Office of Water; E = see user guide Section 2.3.5; L = see user guide on lead; M = mutagen; S = see user guide Section 5; V = volatile; R = RBA applied (See User Guide for Arsenic notice); 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 Guide); s = Concentration may exceed Csat (See User Guide)																										
Toxicity and Chemical-specific Information															Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1					
SFO (mg/kg-day) ⁻¹	k e	IUR (ug/m ³) ⁻¹	k e	RfD _o (mg/kg- day)	k e	RfC _i (mg/m ³) ⁻¹	k e	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-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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)	
2.6E-02	I	7.4E-06	I	3.0E-04	I			V				1.4E+09	5.1E+04	1		Tetrachlorobenzene, 1,2,4,5-	95-94-3					3.5E+02				3.5E+02
2.0E-01	I	5.8E-05	C	2.0E-02	I			V				6.8E+02	1.4E+09	5.7E+03	1	Tetrachloroethane, 1,1,1,2-	630-20-6	1.3E+04		9.4E+02	8.8E+02	3.5E+04			3.5E+04	
2.1E-03	I	2.6E-07	I	6.0E-03	I	4.0E-02	I	V				1.7E+02	1.4E+09	2.4E+03	1	Tetrachloroethylene	127-18-4	1.6E+05		1.1E+04	1.0E+04	7.0E+03		4.1E+02	3.9E+02	
2.0E+01	H			3.0E-02	I			V				1.4E+09	1.1E+05	1	0.1	Tetrachlorophenol, 2,3,4,6-	58-90-2					3.5E+04	8.3E+04		4.1E+02	2.5E+04
				5.0E-04	I			V				1.4E+09		1	0.1	Tetraethyl Dithiopyrophosphate	3689-24-5					5.8E+02	1.4E+03			4.1E+02
				2.0E-03	P			V			2.1E+03	1.4E+09	1.2E+03	1	0.0007	Tetrafluoroethane, 1,1,1,2-	811-97-2					2.3E+03	8.5E+05	4.3E+05		4.3E+05
				2.0E-05	S			V				1.4E+09		1		Tetryl (Trinitrophenylmethylnitramine)	479-45-8									2.3E+03
				1.0E-05	X			V				1.4E+09		1		Thallic Oxide	1314-32-5					2.3E+01				2.3E+01
				1.0E-05	X			V				1.4E+09		1		Thallium (I) Nitrate	10102-45-1					1.2E+01				1.2E+01
				1.0E-05	X			V				1.4E+09		1		Thallium (Soluble Salts)	7440-28-0					1.2E+01				1.2E+01
				1.0E-05	X			V				1.4E+09		1		Thallium Acetate	563-68-8					1.2E+01				1.2E+01
				2.0E-05	X			V				1.4E+09		1		Thallium Carbonate	6533-73-9					2.3E+01				2.3E+01
				1.0E-05	X			V				1.4E+09		1		Thallium Chloride	7791-12-0					1.2E+01				1.2E+01
				1.0E-05	S			V				1.4E+09		1		Thallium Selenite	12039-52-0					1.2E+01				1.2E+01
				2.0E-05	X			V				1.4E+09		1		Thallium Sulfate	7446-18-6					2.3E+01				2.3E+01
				1.3E-02	I			V				1.4E+09		1	0.1	Thifensulfuron-methyl	79277-27-3				1.5E+04	3.6E+04				1.1E+04
				1.0E-02	I			V				1.4E+09		1	0.1	Thiobencarb	28249-77-6					1.2E+04	2.8E+04			8.2E+03
				7.0E-02	X			V				1.4E+09		1	0.0075	Thiodiglycol	111-48-8					8.2E+04	2.6E+06			7.9E+04
				3.0E-04	H			V				1.4E+09		1	0.1	Thiofanox	39196-18-4					3.5E+02	8.3E+02			2.5E+02
				8.0E-02	I			V				1.4E+09		1	0.1	Thiophanate, Methyl	23564-05-8					9.3E+04	2.2E+05			6.6E+04
				5.0E-03	I			V				1.4E+09		1	0.1	Thiram	137-26-8					5.8E+03	1.4E+04			4.1E+03
				6.0E-01	H			V				1.4E+09		1		Tin	7440-31-5					7.0E+05				7.0E+05
				1.0E-04	A	V		V				1.4E+09		1		Titanium Tetrachloride	7550-45-0					9.3E+04		6.0E+05		6.0E+05
				5.0E+00	I	V		V			8.2E+02	1.4E+09	4.3E+03	1		Toluene	108-88-3					9.4E+04				9.4E+04
				8.0E-06	C	V		V				1.4E+09	7.6E+05	1		Toluene-2,4-diisocyanate	584-84-9			8.5E+04	8.5E+04			2.7E+01		2.7E+01
1.8E-01	X			2.0E-04	X			V				1.4E+09		1	0.1	Toluene-2,5-diamine	95-70-5	1.8E+03	4.3E+03			2.3E+02	5.5E+02			1.6E+02
				1.1E-05	C			V				1.4E+09	6.3E+05	1		Toluene-2,6-diisocyanate	91-08-7			7.0E+04	7.0E+04	2.3E+02		2.2E+01		2.2E+01
1.6E-02	P	5.1E-05	C					V				1.4E+09		1	0.1	Toluidine, o- (Methylaniline, 2-)	95-53-4	2.0E+04	4.8E+04	3.3E+07	1.4E+04					
3.0E-02	P			4.0E-03	X			V				1.4E+09		1	0.1	Toluidine, p-	106-49-0	1.1E+04	2.6E+04		7.7E+03	4.7E+03	1.1E+04			3.3E+03
				3.0E+00	P			V			3.4E-01	1.4E+09	1.1E+03	1		Total Petroleum Hydrocarbons (Aliphatic High)	NA					3.5E+06				3.5E+06
				6.0E-01	P	V		V			1.4E+02	1.4E+09	8.3E+02	1		Total Petroleum Hydrocarbons (Aliphatic Low)	NA							2.2E+03	2.2E+03	
				1.0E-02	X	1.0E-01	P	V			6.9E+00	1.4E+09	1.0E+03	1		Total Petroleum Hydrocarbons (Aliphatic Medium)	NA					1.2E+04		4.6E+02	4.4E+02	
				4.0E-02	P			V				1.4E+09		1	0.1	Total Petroleum Hydrocarbons (Aromatic High)	NA					4.7E+04	1.1E+05			3.3E+04
				4.0E-03	P	3.0E-02	P	V			1.8E+03	1.4E+09	3.5E+03	1		Total Petroleum Hydrocarbons (Aromatic Low)	NA					4.7E+03		4.6E+02	4.2E+02	
1.1E+00	I	3.2E-04	I	4.0E-03	P	3.0E-03	P	V				1.4E+09	5.2E+04	1		Total Petroleum Hydrocarbons (Aromatic Medium)	NA					4.7E+03		6.9E+02	6.0E+02	
				7.5E-03	I			V				1.4E+09		1	0.1	Toxaphene	8001-35-2	3.0E+02	7.0E+02	5.2E+06	2.1E+02					6.2E+03
				3.0E-04	A			V				1.4E+09	3.4E+03	1		Tralometrin	66841-25-6					8.8E+03	2.1E+04			6.2E+03
				8.0E+01	X			V				1.4E+09		1	0.1	Tri-n-butyltin	688-73-3					3.5E+02				3.5E+02
				3.0E-02	I			V				1.4E+09		1	0.1	Triacetin	102-76-1					9.3E+07	2.2E+08			6.6E+07
				1.3E-02	I			V				1.4E+09	3.6E+05	1		Triadimefon	43121-43-3					3.5E+04	8.3E+04			2.5E+04
				1.0E-02	I			V				1.4E+09		1	0.1	Triallate	2303-17-5					1.5E+04				1.5E+04
				8.0E-03	I			V				1.4E+09		1	0.1	Triasulfuron	82097-50-5					1.2E+04	2.8E+04			8.2E+03
				5.0E-03	I			V				1.4E+09	4.8E+04	1		Tribenuron-methyl	101200-48-0					9.3E+03	2.2E+04			6.6E+03
9.0E-03	P			1.0E-02	P			V				1.4E+09		1	0.1	Tribromobenzene, 1,2,4-	615-54-3					5.8E+03				5.8E+03
				3.0E-04	P			V				1.4E+09		1	0.1	Tributyl Phosphate	126-73-8	3.6E+04	8.6E+04		2.6E+04	1.2E+04	2.8E+04			8.2E+03
				3.0E-04	P			V				1.4E+09		1	0.1	Tributyltin Compounds	NA					3.5E+02	8.3E+02			2.5E+02
				3.0E-04	I			V				1.4E+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	3.0E+01	H	V			9.1E+02	1.4E+09	1.3E+03	1		Trichloro-1,2,2-trifluoroethane, 1,1,2-	76-13-1					3.5E+07		1.7E+05	1.7E+05	
				2.0E-02	I			V				1.4E+09		1	0.1	Trichloroacetic Acid	76-03-9	4.7E+03	1.1E+04		3.3E+03	2.3E+04	5.5E+04			1.6E+04
2.9E-02	H							V				1.4E+09		1	0.1	Trichloroaniline HCl, 2,4,6-	33663-50-2	1.1E+04	2.7E+04		7.9E+03					
7.0E-03	X			3.0E-05	X			V				1.4E+09		1	0.1	Trichloroaniline, 2,4,6-	634-93-5	4.7E+04	1.1E+05		3.3E+04	3.5E+01	8.3E+01			2.5E+01
				8.0E-04	X			V				1.4E+09	3.2E+04	1		Trichlorobenzene, 1,2,3-	87-61-6					9.3E+02				9.3E+02
2.9E-02	P			1.0E-02	I	2.0E-03	P	V			4.0E+02	1.4E+09	3.0E+04	1		Trichlorobenzene, 1,2,4-	120-82-1	1.1E+04			1.1E+04	1.2E+04		2.6E+02	2.6E+02	
				2.0E+00	I	5.0E+00	I	V			6.4E+02	1.4E+09	1.7E+03	1		Trichloroethane, 1,1,1-	71-55-6					2.3E+06				

Toxicity and Chemical-specific Information														Contaminant		Carcinogenic Target Risk (TR) = 1E-04				Noncancer Hazard Index (HI) = 1			
SFO (mg/kg-day) ⁻¹	k e	IUR (ug/m ³) ⁻¹	k e	RfD _o (mg/kg-day)	k e	RfC _i (mg/m ³)	k e	muta- gen	C _{sat} (mg/kg)	PEF (m ³ /kg)	VF (m ³ /kg)	GI/ABS	ABS	Analyte	CAS No.	Ingestion SL TR=1E-04 (mg/kg)	Dermal SL TR=1E-04 (mg/kg)	Inhalation SL TR=1E-04 (mg/kg)	Carcinogenic SL TR=1E-04 (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)
1.1E-02	I	3.1E-06	I	1.0E-01	I				1.4E+09			1	0.1	Trichlorophenol, 2,4,5-	95-95-4					1.2E+05	2.8E+05		8.2E+04
				1.0E-03	P				1.4E+09			1	0.1	Trichlorophenol, 2,4,6-	88-06-2	3.0E+04	7.0E+04	5.4E+08	2.1E+04	1.2E+03	2.8E+03		8.2E+02
				1.0E-02	I				1.4E+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				1.4E+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		1.3E+03	1.4E+09	1.5E+04	1		Trichloropropane, 1,1,2-	598-77-6					5.8E+03			5.8E+03
				4.0E-03	I	3.0E-04	I	V	M	1.4E+03	1.4E+09	1.6E+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		3.1E+02	1.4E+09	2.3E+03	1	Trichloropropene, 1,2,3-	96-19-5					3.5E+03		3.1E+00	3.1E+00
				2.0E-02	A				1.4E+09			1	0.1	Tricresyl Phosphate (TCP)	1330-78-5					2.3E+04	5.5E+04		1.6E+04
				3.0E-03	I				1.4E+09			1	0.1	Tri-diphenyl	58138-08-2					3.5E+03	8.3E+03		2.5E+03
						7.0E-03	I	V		2.8E+04	1.4E+09	1.6E+04	1	Triethylamine	121-44-8							4.8E+02	4.8E+02
				2.0E+00	P				1.4E+09			1	0.1	Triethylene Glycol	112-27-6					2.3E+06	5.5E+06		1.6E+06
7.7E-03	I			2.0E+01	P	V			4.8E+03	1.4E+09	7.1E+02	1		Trifluoroethane, 1,1,1-	420-46-2							6.2E+04	6.2E+04
				7.5E-03	I		V		1.4E+09	5.1E+05		1		Trifluralin	1582-09-8	4.2E+04			4.2E+04	8.8E+03			8.8E+03
2.0E-02	P			1.0E-02	P				1.4E+09			1	0.1	Trimethyl Phosphate	512-56-1	1.6E+04	3.9E+04		1.1E+04	1.2E+04	2.8E+04		8.2E+03
				5.0E-03	P	V			2.9E+02	1.4E+09	9.4E+03	1		Trimethylbenzene, 1,2,3-	526-73-8							2.1E+02	2.1E+02
				7.0E-03	P	V			2.2E+02	1.4E+09	7.9E+03	1		Trimethylbenzene, 1,2,4-	95-63-6							2.4E+02	2.4E+02
				1.0E-02	X		V		1.8E+02	1.4E+09	6.6E+03	1		Trimethylbenzene, 1,3,5-	108-67-8					1.2E+04			1.2E+04
				1.0E-02	X		V		3.0E+01	1.4E+09	1.0E+03	1		Trimethylpentene, 2,4,4-	25167-70-8					1.2E+04			1.2E+04
				3.0E-02	I				1.4E+09			1	0.019	Trinitrobenzene, 1,3,5-	99-35-4					3.5E+04	4.4E+05		3.2E+04
3.0E-02	I			5.0E-04	I				1.4E+09			1	0.032	Trinitrotoluene, 2,4,6-	118-96-7	1.1E+04	8.0E+04		9.6E+03	5.8E+02	4.3E+03		5.1E+02
				2.0E-02	P				1.4E+09			1	0.1	Triphenylphosphine Oxide	791-28-6					2.3E+04	5.5E+04		1.6E+04
				2.0E-02	A				1.4E+09			1	0.1	Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8					2.3E+04	5.5E+04		1.6E+04
2.3E+00	C	6.6E-04	C	1.0E-02	X				1.4E+09			1	0.1	Tris(1-chloro-2-propyl)phosphate	13674-84-5					1.2E+04	2.8E+04		8.2E+03
2.0E-02	P			7.0E-03	P				1.4E+09			1	0.1	Tris(2,3-dibromopropyl)phosphate	12672-7	1.4E+02		1.7E+03	1.3E+02				
									1.4E+09			1	0.1	Tris(2-chloroethyl)phosphate	11596-8	1.6E+04	3.9E+04		1.1E+04	8.2E+03	1.9E+04		5.7E+03
3.2E-03	P			1.0E-01	P				1.4E+09			1	0.1	Tris(2-ethylhexyl)phosphate	78-42-2	1.0E+05	2.4E+05		7.2E+04	1.2E+05	2.8E+05		8.2E+04
				8.0E-04	P				1.4E+09			1		Tungsten	7440-33-7					9.3E+02			9.3E+02
				3.0E-03	I	4.0E-05	A		1.4E+09			1		Uranium (Soluble Salts)	NA					3.5E+03		2.4E+05	3.5E+03
1.0E+00	C	2.9E-04	C					M	1.4E+09			1	0.1	Urethane	51-96	3.3E+02	7.7E+02	5.7E+06	2.3E+02				
		8.3E-03	P	9.0E-03	I	7.0E-06	P		1.4E+09		0.026	1		Vanadium Pentoxide	1314-62-1			2.0E+05	2.0E+05	1.1E+04		4.2E+04	8.4E+03
				5.0E-03	S	1.0E-04	A		1.4E+09		0.026	1		Vanadium and Compounds	7440-62-2					5.9E+03		6.0E+05	5.8E+03
				1.0E-03	I		V		1.4E+09	1.2E+05		1		Vernolate	1929-77-7					1.2E+03			1.2E+03
				2.5E-02	I				1.4E+09			1	0.1	Vinclozolin	50471-44-8					2.9E+04	6.9E+04		2.1E+04
				1.0E+00	H	2.0E-01	I	V		2.8E+03	1.4E+09	4.4E+03	1	Vinyl Acetate	108-05-4					1.2E+06			3.8E+03
		3.2E-05	H			3.0E-03	I	V		2.5E+03	1.4E+09	1.4E+03	1	Vinyl Bromide	593-60-2			5.2E+01	5.2E+01			1.8E+01	1.8E+01
7.2E-01	I	4.4E-06	I	3.0E-03	I	1.0E-01	I	V	M	3.9E+03	1.4E+09	9.6E+02	1	Vinyl Chloride	75-01-4	4.5E+02		2.7E+02	1.7E+02	3.5E+03		4.2E+02	3.7E+02
				3.0E-04	I				1.4E+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	V		3.9E+02	1.4E+09	5.6E+03	1	Xylene, p-	106-42-3					2.3E+05		2.4E+03	2.4E+03
				2.0E-01	S	1.0E-01	S	V		3.9E+02	1.4E+09	5.5E+03	1	Xylene, m-	108-38-3					2.3E+05		2.4E+03	2.4E+03
				2.0E-01	S	1.0E-01	S	V		4.3E+02	1.4E+09	6.5E+03	1	Xylene, o-	95-47-6					2.3E+05		2.8E+03	2.8E+03
				2.0E-01	I	1.0E-01	I	V		2.6E+02	1.4E+09	5.7E+03	1	Xylenes	1330-20-7					2.3E+05		2.5E+03	2.5E+03
				3.0E-04	I				1.4E+09			1		Zinc Phosphide	1314-84-7					3.5E+02			3.5E+02
				3.0E-01	I				1.4E+09			1		Zinc and Compounds	7440-66-6					3.5E+05			3.5E+05
				5.0E-02	I				1.4E+09			1	0.1	Zineb	12122-67-7					5.8E+04	1.4E+05		4.1E+04
				8.0E-05	X				1.4E+09			1		Zirconium	7440-67-7					9.3E+01			9.3E+01

Regional Removal Management Level (RML) Chemical-specific Parameters Supporting Table May 2016

Contaminant		Molecular Weight		Volatility Parameters				Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients				Water Solubility		Tapwater Dermal Parameters											
Analyte	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m ³ /mole)		H' and HLC Ref		VP	VP Ref	MP	MP Ref	Density (g/cm ³)	Density Ref	Dia (cm ² /s)	Dw (cm ² /s)	D _a and D _w Ref	K _{oc} (L/kg)	K _{oc} Ref	K _{oc}	K _{oc} Ref	log K _{ow} (unitless)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/event)	t* (hr)	K _p (cm ² /hr)	K Ref			
					H' (unitless)	H' and HLC Ref																											
Acetate	30560-19-1	1.8E+02	PHYSPROP	2.0E-11	5.0E-13	EPI	1.7E-06	PHYSPROP	8.8E+01	PHYSPROP	1.4E+00	CRCB9	3.7E-02	8.0E-06	WATER9	1.0E+01	EPI	8.5E-01	PHYSPROP	8.2E+05	PHYSPROP	2.1E-04	1.1E+00	2.7E+00	4.0E-05	EPI	1.3E-03	1.9E-01	4.5E-01	5.0E-04	EPI		
Acetaldehyde	75-07-0	4.4E+01	PHYSPROP	2.7E-03	6.7E-05	PHYSPROP	9.0E+02	PHYSPROP	1.2E+02	PHYSPROP	7.8E-01	CRCB9	1.3E-01	1.4E-05	WATER9	1.0E+00	EPI	3.4E-01	PHYSPROP	1.0E+06	PHYSPROP	1.0E-03	1.0E+00	2.2E+02	PHYSPROP	3.1E-02	3.4E+00	8.2E+01	5.0E-03	EPI			
Acetochlor	34256-82-1	2.7E+02	PHYSPROP	9.1E-07	2.2E-08	PHYSPROP	2.8E-05	PHYSPROP	1.1E+01	PubChem	1.1E+00	PubChem	2.2E-02	5.6E-06	WATER9	3.0E+02	EPI	3.0E+00	PHYSPROP	2.2E+02	PHYSPROP	1.0E+06	PHYSPROP	1.5E-03	2.2E-01	5.3E-01	5.0E-04	EPI	1.3E-03	1.9E-01	4.5E-01	5.0E-04	EPI
Acetone	67-64-1	5.8E+01	PHYSPROP	1.4E-03	3.5E-05	PHYSPROP	2.3E+02	PHYSPROP	9.5E+01	PHYSPROP	7.8E-01	CRCB9	1.1E-01	1.2E-05	WATER9	2.4E+00	EPI	2.4E-01	PHYSPROP	1.0E+06	PHYSPROP	1.0E-03	1.0E+00	2.2E+02	PHYSPROP	3.1E-02	3.4E+00	8.2E+01	5.0E-03	EPI			
Acetone Cyanohydrin	75-86-5	8.5E+01	PHYSPROP	8.1E-08	2.0E-09	PHYSPROP	3.4E-01	PHYSPROP	1.3E+02	PHYSPROP	9.4E-01	CRCB9	8.6E-02	1.0E-05	WATER9	1.0E+00	EPI	3.0E-02	PHYSPROP	1.0E+06	PHYSPROP	1.0E-03	1.0E+00	2.2E+02	PHYSPROP	3.1E-02	3.4E+00	8.2E+01	5.0E-03	EPI			
Acetone	75-07-0	4.4E+01	PHYSPROP	1.4E-03	3.5E-05	PHYSPROP	8.9E+01	PHYSPROP	4.4E+01	PHYSPROP	7.9E-01	CRCB9	1.3E-01	1.4E-05	WATER9	4.7E+00	EPI	3.4E-01	PHYSPROP	1.0E+06	PHYSPROP	1.0E-03	1.0E+00	2.2E+02	PHYSPROP	3.1E-02	3.4E+00	8.2E+01	5.0E-03	EPI			
Acetophenone	98-86-2	1.2E+02	PHYSPROP	4.3E-04	1.0E-05	PHYSPROP	4.0E-01	PHYSPROP	2.0E+01	PHYSPROP	1.0E+00	CRCB9	6.5E-02	8.7E-06	WATER9	5.2E+01	EPI	1.6E+00	PHYSPROP	1.0E+06	PHYSPROP	1.0E-03	1.0E+00	2.2E+02	PHYSPROP	3.1E-02	3.4E+00	8.2E+01	5.0E-03	EPI			
Acetylaminofluorene, 2-	53-96-3	2.2E+02	PHYSPROP	7.8E-09	1.9E-10	PHYSPROP	9.4E-08	PHYSPROP	1.9E+02	PHYSPROP	1.1E+00	CRCB9	5.2E-02	6.0E-06	WATER9	2.2E+03	EPI	3.1E+00	PHYSPROP	5.5E+00	PHYSPROP	7.2E-02	1.9E+01	4.5E+00	1.2E-02	RAGSE	2.2E-03	2.2E-01	5.2E-01	7.5E-04	EPI		
Acrolein	107-02-8	5.6E+01	PHYSPROP	5.0E-03	1.2E-04	PHYSPROP	2.7E+02	PHYSPROP	8.8E-01	PHYSPROP	8.4E-01	CRCB9	1.1E-01	1.2E-05	WATER9	1.0E+00	EPI	1.0E-02	PHYSPROP	2.1E+05	PHYSPROP	2.2E-03	2.2E-01	5.2E-01	7.5E-04	EPI	1.3E-03	1.9E-01	4.5E-01	5.0E-04	EPI		
Acrylamide	79-06-1	7.1E+01	PHYSPROP	7.0E-08	1.7E-09	EPI	7.0E-03	PHYSPROP	8.5E+01	PHYSPROP	1.2E+00	LANGE	1.1E-01	1.3E-05	WATER9	5.7E+00	EPI	6.7E-01	PHYSPROP	3.9E+05	PHYSPROP	7.3E-04	2.6E-01	6.3E-01	2.2E-04	EPI	3.4E-03	2.7E-01	6.4E-01	1.1E-03	EPI		
Acrylic Acid	79-10-7	7.2E+01	PHYSPROP	1.5E-05	3.7E-07	EPI	4.0E+00	PHYSPROP	1.3E+01	PHYSPROP	1.1E+00	CRCB9	1.0E-01	1.2E-05	WATER9	1.4E+00	EPI	3.5E-01	PHYSPROP	1.0E+06	PHYSPROP	3.3E-03	2.1E-01	5.0E-01	1.2E-03	EPI	3.4E-03	2.7E-01	6.4E-01	1.1E-03	EPI		
Acrylonitrile	107-13-1	5.3E+01	PHYSPROP	5.6E-03	1.4E-04	PHYSPROP	1.1E+02	PHYSPROP	8.4E-01	PHYSPROP	8.0E-01	CRCB9	1.1E-01	1.2E-05	WATER9	8.5E+00	EPI	2.5E-01	PHYSPROP	7.0E+04	PHYSPROP	3.3E-03	2.1E-01	5.0E-01	1.2E-03	EPI	3.4E-03	2.7E-01	6.4E-01	1.1E-03	EPI		
Adiponitrile	111-69-3	1.1E+02	PHYSPROP	4.9E-08	1.2E-09	EPI	6.8E-04	PHYSPROP	1.0E+00	PHYSPROP	9.7E-01	CRCB9	7.1E-02	9.0E-06	WATER9	2.0E+01	EPI	3.2E-01	PHYSPROP	8.0E+04	PHYSPROP	9.5E-04	4.2E-01	1.0E+00	2.4E-04	EPI	3.4E-03	2.7E-01	6.4E-01	1.1E-03	EPI		
Alachlor	15972-60-8	2.7E+02	PHYSPROP	3.4E-07	8.3E-09	PHYSPROP	2.2E-05	PHYSPROP	4.0E+01	PHYSPROP	1.1E+00	CRCB9	2.3E-02	5.7E-06	WATER9	3.1E+02	EPI	3.5E+00	PHYSPROP	2.4E+02	PHYSPROP	6.6E-02	3.4E+00	8.2E+00	1.1E-02	EPI	4.0E-03	1.2E+00	2.9E+00	7.6E-04	EPI		
Aldicarb	116-06-3	1.9E+02	PHYSPROP	5.9E-08	1.4E-09	EPI	3.5E-05	PHYSPROP	9.9E+01	PHYSPROP	1.2E+00	CRCB9	3.2E-02	7.2E-06	WATER9	2.5E+01	EPI	1.1E+00	PHYSPROP	6.0E+03	PHYSPROP	4.0E-03	1.2E+00	2.9E+00	7.6E-04	EPI	4.0E-03	1.2E+00	2.9E+00	7.6E-04	EPI		
Aldicarb Sulfone	1646-88-4	2.2E+02	PHYSPROP	1.4E-07	3.4E-09	EPI	9.0E-05	PHYSPROP	1.4E+02	PHYSPROP	1.6E+00	PubChem	5.2E-02	6.1E-06	WATER9	1.0E+01	EPI	5.7E-01	PHYSPROP	1.0E+04	PHYSPROP	2.1E-04	1.8E+00	4.4E+00	3.7E-05	EPI	1.8E-04	1.5E+00	3.6E+00	3.3E-05	EPI		
Aldicarb sulfoxide	1646-87-3	2.1E+02	PHYSPROP	4.0E-08	9.7E-10	EPI	1.0E-04	PHYSPROP	7.8E+01	EPI	1.6E+00	PubChem	5.4E-02	6.4E-06	WATER9	1.0E+01	EPI	7.8E-01	PHYSPROP	2.8E+04	PHYSPROP	2.2E+00	1.2E+01	4.8E+01	2.9E-04	EPI	1.8E-04	1.5E+00	3.6E+00	3.3E-05	EPI		
Aldrin	309-00-2	3.6E+02	PHYSPROP	1.8E-03	4.4E-05	PHYSPROP	1.2E-04	PHYSPROP	1.0E+02	PHYSPROP	1.6E+00	PubChem	2.3E-02	5.8E-06	WATER9	8.2E+04	EPI	6.5E+00	PHYSPROP	1.7E-02	PHYSPROP	2.2E+00	1.2E+01	4.8E+01	2.9E-04	EPI	1.8E-04	1.5E+00	3.6E+00	3.3E-05	EPI		
Allyl Alcohol	107-18-6	5.8E+01	PHYSPROP	2.1E-04	5.0E-06	PHYSPROP	2.6E+01	PHYSPROP	1.3E+02	PHYSPROP	8.5E-01	CRCB9	1.1E-01	1.2E-05	WATER9	1.9E+00	EPI	1.7E-01	PHYSPROP	1.0E+06	PHYSPROP	2.8E-03	2.2E-01	5.3E-01	5.0E-04	EPI	1.3E-03	1.9E-01	4.5E-01	5.0E-04	EPI		
Allyl Chloride	107-05-1	7.7E+01	PHYSPROP	4.5E-01	1.1E-02	EPI	3.7E+02	PHYSPROP	9.4E+01	PHYSPROP	9.4E-01	CRCB9	9.4E-02	1.1E-05	WATER9	4.0E+01	EPI	1.9E+00	PHYSPROP	3.4E+03	PHYSPROP	2.0E-03	1.5E-01	8.6E-01	1.0E-03	RAGSE	2.0E-03	1.5E-01	8.6E-01	1.0E-03	RAGSE		
Aluminum	7429-90-5	2.7E+01	CRCB9				0.0E+00	NIOSH	6.6E+02	CRCB9	2.7E+00	CRCB9						1.5E+03	BAES			2.9E-03	2.2E-01	5.3E-01	1.0E-03	RAGSE	4.6E-02	2.0E+00	4.7E+00	7.9E-03	EPI		
Aluminum Phosphide	20859-73-8	5.8E+01	PHYSPROP	1.6E-03	3.8E-05	CRCB9	2.6E+03	CRCB9	2.4E+00	CRCB9	2.4E+00	CRCB9	2.4E+00	CRCB9									2.9E-03	2.2E-01	5.3E-01	1.0E-03	RAGSE	4.6E-02	2.0E+00	4.7E+00	7.9E-03	EPI	
Ameryn	834-12-8	2.3E+02	PHYSPROP	9.9E-08	2.4E-09	EPI	2.7E-06	PHYSPROP	8.8E+01	PHYSPROP	1.2E+00	PubChem	5.1E-02	6.0E-06	WATER9	4.3E+02	EPI	3.0E+00	PHYSPROP	2.1E+02	PHYSPROP	4.6E-02	2.0E+00	4.7E+00	7.9E-03	EPI	4.6E-02	2.0E+00	4.7E+00	7.9E-03	EPI		
Aminobiphenyl, 4-	92-67-1	1.7E+02	PHYSPROP	6.0E-06	1.5E-07	PHYSPROP	1.2E-04	PHYSPROP	5.4E+01	PHYSPROP	1.2E+00	PubChem	6.2E-02	7.3E-06	WATER9	2.5E+03	EPI	2.9E+00	PHYSPROP	2.0E+02	PHYSPROP	7.0E-02	9.3E-01	2.2E+00	1.4E-02	EPI	4.6E-02	2.0E+00	4.7E+00	7.9E-03	EPI		
Aminophenol, m-	591-27-5	1.1E+02	PHYSPROP	8.1E-09	2.0E-10	PHYSPROP	9.6E-03	PHYSPROP	1.2E+02	PHYSPROP	1.2E+00	PubChem	8.3E-02	9.7E-06	WATER9	9.0E+01	EPI	2.1E-01	PHYSPROP	2.7E+04	PHYSPROP	2.1E-03	4.3E-01	1.0E+00	5.3E-04	EPI	1.6E-03	4.3E-01	1.0E+00	4.1E-04	EPI		
Aminophenol, p-	123-30-8	1.1E+02	PHYSPROP	1.5E-08	3.6E-10	EPI	4.0E-05	EPI	1.9E+02	PHYSPROP	1.2E+00	PubChem	8.3E-02	9.7E-06	WATER9	9.0E+01	EPI	2.1E-01	PHYSPROP	2.7E+04	PHYSPROP	2.1E-03	4.3E-01	1.0E+00	5.3E-04	EPI	1.6E-03	4.3E-01	1.0E+00	4.1E-04	EPI		
Amiratz	33089-61-1	2.9E+02	PHYSPROP	4.0E-04	9.9E-06	PHYSPROP	2.0E-06	PHYSPROP	8.6E+01	PHYSPROP	1.1E+00	CRCB9	2.2E-02	5.4E-06	WATER9	2.6E+05	EPI	5.5E+00	PHYSPROP	1.0E+00	PHYSPROP	1.1E+00	4.6E+00	1.8E+01	1.6E-01	EPI	1.1E+00	4.6E+00	1.8E+01	1.6E-01	EPI		
Ammonia	7664-41-7	1.7E+01	PHYSPROP	6.6E-04	1.6E-05	PHYSPROP	7.5E+03	PHYSPROP	7.8E+01	PHYSPROP	7.0E-01	CRCB9	2.3E-01	2.2E-05	WATER9	1.0E+00	OTHER	4.8E+05	PHYSPROP	1.6E-03	1.3E-01	3.1E-01	1.0E-03	RAGSE	1.6E-03	1.3E-01	3.1E-01	1.0E-03	RAGSE				
Ammonium Sulfamate	7773-06-0	1.1E+02	CRCB9				0.0E+00	NIOSH	1.3E+02	CRCB9	1.8E+00	PubChem											4.1E+01	SSL	4.5E-03	3.5E-01	8.4E-01	1.9E-03	EPI				
Amyl Alcohol, tert-	75-85-4	8.8E+01	PHYSPROP	5.6E-04	1.4E-05	PHYSPROP	1.7E+01	PHYSPROP	9.1E+00	PHYSPROP	8.1E-01	CRCB9	7.9E-02	9.1E-06	WATER9	4.1E+00	EPI	8.9E-01	PHYSPROP	1.1E+06	PHYSPROP	7.1E-03	4.6E-01	1.1E+00	1.0E-03	RAGSE	4.1E-03	3.3E-01	7.9E-01	2.0E-03	EPI		
Aniline	62-53-3	9.3E+01	PHYSPROP	8.3E-05	2.0E-06	PHYSPROP	6.7E-01	PHYSPROP	4.0E+00	PHYSPROP	1.0E+00	CRCB9	8.3E-02	1.0E-05	WATER9	7.0E+01	EPI	9.0E-01	PHYSPROP	3.6E+04	PHYSPROP	6.9E-03	3.5E-01	8.4E-01	1.9E-03	EPI	1.1E-01	1.5E+00	3.7E+00	1.9E-02	EPI		
Anthraquinone, 9,10-	84-65-1	2.1E+02	PHYSPROP	9.6E-07	2.4E-08	EPI	1.2E-07	PHYSPROP	2.9E+02	PHYSPROP	1.2E+00	CRCB9	5.4E-02	6.3E-06	WATER9	5.0E+03	EPI	3.4E+00	PHYSPROP	1.4E+00	PHYSPROP	4.3E-03	5.3E-01	1.3E+00	1.0E-03	RAGSE	4.3E-03	5.3E-01	1.3E+00	1.0E-03	RAGSE		
Antimony (metallic)	7440-36-0	1.2E+02	PHYSPROP				0.0E+00	NIOSH																									

Regional Removal Management Level (RML) Chemical-specific Parameters Supporting Table May 2016

Contaminant	Molecular Weight	Volatility Parameters						Melting Point		Density		Diffusivity in Air and Water			Partition Coefficients				Water Solubility		Tapwater Dermal Parameters							
		CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m ³ /mole)	H' and HLC Ref	VP	VP Ref	MP	MP Ref	Density (g/cm ³)	Density Ref	Dia (cm ² /s)	Diw (cm ² /s)	D ₁₀ and D ₁₀₀ Ref	K _{oc} (L/kg)	K _{oc} Ref	K _{oc}	K _{oc} Ref	log K _{ow} (unitless)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr)	t* (hr)	K _p (cm/hr)
Butylated hydroxyanisole	25013-16-5	3.6E+02	PHYSPROP	4.8E-05	1.2E-06	PHYSPROP	2.5E-03	PHYSPROP	5.1E+01	PHYSPROP			3.8E-02	4.4E-06	WATER9	8.4E+02	EPI	3.5E+00	PHYSPROP	2.1E+02	PHYSPROP	2.0E+01	PHYSPROP	2.4E-01	1.1E+01	2.6E+01	3.3E-02	EPI
Butylated hydroxytoluene	128-37-0	2.2E+02	PHYSPROP	1.7E-04	4.1E-06	PHYSPROP	5.2E-03	EPI	7.1E+01	PHYSPROP	8.9E-01	CRCB9	2.3E-02	5.6E-06	WATER9	1.5E+04	EPI	5.1E+00	PHYSPROP	6.0E+01	PHYSPROP	1.3E+00	PHYSPROP	1.3E+00	1.8E+00	7.1E+00	2.2E-01	EPI
Butylbenzene, n-	104-51-8	1.3E+02	PHYSPROP	6.5E-01	1.6E-02	EPI	1.1E+00	PHYSPROP	-8.8E+01	PHYSPROP	8.6E-01	CRCB9	5.3E-02	7.3E-06	WATER9	1.5E+03	EPI	4.4E+00	PHYSPROP	1.2E+01	PHYSPROP	1.0E+00	PHYSPROP	1.0E+00	5.9E-01	2.3E+00	2.3E-01	EPI
n-Butylbenzene, sec-	135-98-8	1.3E+02	PHYSPROP	7.2E-01	1.8E-02	EPI	1.8E+00	PHYSPROP	-8.3E+01	PHYSPROP	8.6E-01	LANGE	5.3E-02	7.3E-06	WATER9	1.3E+03	EPI	4.6E+00	PHYSPROP	1.3E+01	PHYSPROP	1.3E+00	PHYSPROP	1.3E+00	5.9E-01	2.3E+00	3.0E-01	EPI
Butylbenzene, tert-	98-06-6	1.3E+02	PHYSPROP	5.4E-01	1.3E-02	EPI	2.2E+00	PHYSPROP	-5.8E+01	PHYSPROP	8.7E-01	CRCB9	5.3E-02	7.4E-06	WATER9	1.0E+03	EPI	4.1E+00	PHYSPROP	3.0E+01	PHYSPROP	3.0E+01	PHYSPROP	6.6E-01	5.9E-01	2.3E+00	1.5E-01	EPI
Cacodylic Acid	75-60-5	3.4E+02	PHYSPROP	1.1E+02	1.8E-14	PHYSPROP	0.0E+00	NIOSH	2.2E+02	PHYSPROP	8.7E+00	CRCB9	7.1E-02	8.3E-06	WATER9	4.4E+01	EPI	3.6E-01	PHYSPROP	2.0E+06	PHYSPROP	3.0E+06	PHYSPROP	2.1E-03	6.2E-01	3.5E+00	4.6E-04	EPI
Cadmium (Salt)	7440-43-9	1.1E+02	PHYSPROP	0.0E+00			0.0E+00	NIOSH	3.2E+02	PHYSPROP	8.7E+00	CRCB9				7.5E+01	SSL							4.1E-03	4.5E-01	1.1E+00	1.0E-03	RAGSE
Cadmium (Water)	7440-43-9	1.1E+02	PHYSPROP	0.0E+00			0.0E+00	NIOSH	3.2E+02	PHYSPROP	8.7E+00	CRCB9				7.5E+01	SSL							4.1E-03	4.5E-01	1.1E+00	1.0E-03	RAGSE
Calcium Chromate	13765-19-0	1.6E+02	CRCB9	1.0E-06	2.5E-08	PHYSPROP	1.6E-03	EPI	6.9E+01	PHYSPROP	1.0E+00	LANGE	6.9E-02	9.0E-06	WATER9	2.5E+01	EPI	-1.9E-01	YAWS	7.7E+05	PHYSPROP	4.8E-03	7.9E-01	1.9E+00	1.0E-03	RAGSE		
Caprolactam	105-60-2	1.1E+02	PHYSPROP	1.0E-06	2.5E-08	PHYSPROP	1.6E-03	EPI	6.9E+01	PHYSPROP	1.0E+00	LANGE	6.9E-02	9.0E-06	WATER9	2.5E+01	EPI	-1.9E-01	YAWS	7.7E+05	PHYSPROP	4.1E-03	4.5E-01	1.1E+00	1.0E-03	RAGSE		
Captafol	2425-06-1	3.5E+02	PHYSPROP	2.0E-07	4.9E-09	EPI	1.5E-08	EPI	1.6E+02	PHYSPROP	3.8E-02	4.5E-06	WATER9	7.8E+02	EPI	3.8E+00	PHYSPROP	1.4E+00	PHYSPROP	1.4E+00	PHYSPROP	4.1E-02	9.5E+00	2.3E+01	5.8E-03	EPI		
Captan	133-06-2	3.0E+02	PHYSPROP	2.9E-07	7.0E-09	EPI	9.0E-08	PHYSPROP	1.8E+02	PHYSPROP	1.7E+00	CRCB9	2.6E-02	6.9E-06	WATER9	2.5E+02	EPI	2.8E+00	PHYSPROP	5.1E+00	PHYSPROP	1.6E-02	5.1E+00	1.2E+01	2.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E-07	3.3E-09	EPI	1.4E-06	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRCB9	2.7E-02	7.1E-06	WATER9	3.5E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	2.4E-02	1.4E+00	3.4E+00	4.3E-03	EPI		
Carbaryl	63-25-2	2.0E+02	PHYSPROP	1.3E																								

Regional Removal Management Level (RML) Chemical-specific Parameters Supporting Table May 2016

Contaminant	Molecular Weight		Volatility Parameters				Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients				Water Solubility		Tapwater Dermal Parameters										
	Analyte	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m ³ /mole)	H' and HLC Ref	VP	VP Ref	MP	MP Ref	Density (g/cm ³)	Density Ref	Dia (cm ² /s)	Diw (hr/cm/s)	D _a and D _w Ref	K _{oc} (L/kg)	K _{oc} Ref	K _{oc}	K _{oc} Ref	log K _{ow} (unitless)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/event)	t [*] (hr)	K _p (cm/hr)	K Ref		
~Potassium Cyanide	151-50-8	6.5E+01	PHYSPROP				0.0E+00	NIOSH	6.3E+02	PHYSPROP	1.6E+00	CRCB9											7.2E+05	PHYSPROP	6.2E-03	2.4E-01	5.8E-01	2.0E-03	RAGSE		
~Potassium Silver Cyanide	506-61-6	2.0E+02	PHYSPROP																				1.1E+02	1.4E+00	3.3E+00	2.0E-03	RAGSE				
~Silver Cyanide	506-64-9	1.3E+02	PHYSPROP																				2.3E+01	PHYSPROP	4.5E-03	5.9E-01	1.4E+00	1.0E-03	RAGSE		
~Sodium Cyanide	143-33-9	4.9E+01	PHYSPROP				0.0E+00	NIOSH	5.6E+02	PHYSPROP	1.6E+00	CRCB9											5.8E+05	CRCB9	2.7E-03	2.0E-01	4.7E-01	1.0E-03	RAGSE		
~Thiocyanates	NA																														
~Thiocyanic Acid	463-56-0	5.9E+01	PHYSPROP				4.7E+00	PPRTV	5.0E+00	PPRTV	1.1E+00	PPRTV	1.2E-01	1.4E-05	WATER9							5.8E-01	OTHER								
~Zinc Cyanide	557-21-1	1.2E+02	PHYSPROP																				4.7E+00	CRCB9	3.5E-03	4.8E-01	1.1E+00	6.0E-04	RAGSE		
Cyclohexane	110-82-7	8.4E+01	PHYSPROP	6.1E+00	1.5E-01	PHYSPROP	9.7E+01	PHYSPROP	6.6E+00	2.0E+02	CRCB9	7.7E-01	CRCB9	8.0E-02	9.1E-06	WATER9	1.5E+02	EPI	3.4E+00	PHYSPROP	5.5E+01	PHYSPROP	3.6E-01	3.1E-01	7.5E-01	1.0E-01	EPI				
Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro-	87-84-3	5.1E+02	PHYSPROP	3.9E+05	9.6E-07	PHYSPROP	3.5E+06	PHYSPROP	2.0E+02	CRCB9	3.0E+02	3.2E+06	WATER9	2.8E+03	EPI	4.7E+00	PHYSPROP	5.5E+01	PHYSPROP	5.5E+01	PHYSPROP	2.5E+02	PHYSPROP	2.5E+02	7.9E+01	1.9E+02	2.8E-03	EPI			
Cyclohexanone	108-94-1	9.8E+01	PHYSPROP	3.7E-04	9.0E-06	PHYSPROP	4.3E+00	PHYSPROP	-3.1E-01	PHYSPROP	9.5E-01	CRCB9	7.7E-02	9.4E-06	WATER9	1.7E+01	EPI	8.1E-01	PHYSPROP	2.5E+04	PHYSPROP	5.8E-03	3.7E-01	8.9E-01	1.5E-03	EPI					
Cyclohexene	110-83-8	8.2E+01	PHYSPROP	1.9E+00	4.6E-02	PHYSPROP	8.9E+01	PHYSPROP	-1.0E+02	PHYSPROP	8.1E-01	NIOSH	8.3E-02	9.5E-06	WATER9	1.5E+02	EPI	2.9E+00	PHYSPROP	2.1E+02	PHYSPROP	1.5E-01	3.0E-01	7.3E-01	4.3E-02	EPI					
Cyclohexylamine	108-91-8	9.9E+01	PHYSPROP	1.7E-04	4.2E-06	PHYSPROP	1.0E+01	PHYSPROP	-1.8E-01	PHYSPROP	8.2E-01	CRCB9	7.1E-02	8.5E-06	WATER9	3.2E+01	EPI	1.5E+00	PHYSPROP	1.0E+06	PHYSPROP	1.6E-02	3.8E-01	9.1E-01	4.3E-03	EPI					
Cyfluthrin	68359-37-5	4.3E+02	PHYSPROP	1.2E-06	2.9E-08	EPI	1.5E-10	PHYSPROP	6.0E+01	PHYSPROP	3.3E-02	3.9E-06	WATER9	1.3E+05	EPI	6.0E+00	PHYSPROP	3.0E+03	PHYSPROP	3.0E+03	PHYSPROP	4.1E-01	2.8E+01	6.8E+01	5.2E-02	EPI					
Cyhalothrin	68085-85-8	4.5E+02	PHYSPROP	6.1E-05	1.5E-06	EPI	1.5E-09	PHYSPROP	4.9E+01	PHYSPROP	3.2E-02	3.8E-06	WATER9	3.4E+05	EPI	6.9E+00	PHYSPROP	5.0E+03	PHYSPROP	5.0E+03	PHYSPROP	1.7E+00	3.5E+01	1.4E+02	2.1E-01	EPI					
Cypermethrin	52315-07-8	4.2E+02	PHYSPROP	1.7E-05	4.2E-07	EPI	3.1E-09	PHYSPROP	8.1E+01	PHYSPROP	1.3E+00	CRCB9	1.9E-02	4.7E-06	WATER9	8.0E+04	EPI	6.6E+00	PHYSPROP	4.0E+03	PHYSPROP	6.0E-01	3.2E+01	9.1E+01	7.7E-02	EPI					
Cyromazine	66215-07-8	1.7E+02	PHYSPROP	2.3E-12	5.7E-14	EPI	3.4E-09	PHYSPROP	2.2E+02	PHYSPROP	6.3E-02	7.3E-06	WATER9	2.9E+01	EPI	6.1E-02	PHYSPROP	1.3E+04	PHYSPROP	1.3E+04	PHYSPROP	4.0E-03	9.0E-01	2.2E+00	8.0E-04	EPI					
DD	72-54-8	3.2E+02	PHYSPROP	2.7E-04	6.6E-06	PHYSPROP	1.4E-06	PHYSPROP	1.1E+02	PHYSPROP	4.1E-02	4.7E-06	WATER9	1.2E+05	EPI	6.0E+00	PHYSPROP	9.0E+02	PHYSPROP	9.0E+02	PHYSPROP	1.7E+00	6.5E+00	2.6E+01	2.5E-01	EPI					
DDE, p,p'	72-55-9	3.2E+02	PHYSPROP	1.7E-03	4.2E-05	PHYSPROP	6.0E-06	EPI	8.9E+01	PHYSPROP	1.4E+00	LookChem	2.3E-02	5.9E-06	WATER9	1.2E+05	EPI	6.5E+00	PHYSPROP	4.0E-02	PHYSPROP	3.7E+00	6.4E+00	2.7E+01	5.5E-01	EPI					
DDT	50-29-3	3.5E+02	PHYSPROP	3.4E-04	8.3E-06	PHYSPROP	1.6E-07	PHYSPROP	1.1E+02	PHYSPROP	3.8E-02	4.4E-06	WATER9	1.7E+05	EPI	6.9E+00	PHYSPROP	5.5E+03	PHYSPROP	5.5E+03	PHYSPROP	4.5E+00	1.0E+01	4.4E+01	6.3E-01	EPI					
Dalapon	75-99-0	1.4E+02	PHYSPROP	2.3E-06	5.7E-08	EPI	1.5E-04	EPI	-5.0E+00	PHYSPROP	1.4E+00	CRCB9	6.0E-02	9.4E-06	WATER9	3.2E+00	EPI	7.8E-01	PHYSPROP	5.0E+05	PHYSPROP	3.7E-03	6.6E-01	1.6E+00	8.2E-04	EPI					
Diazinon	1596-94-5	1.6E+02	PHYSPROP	1.7E-08	4.2E-10	EPI	2.0E-04	PHYSPROP	1.5E+02	PHYSPROP	6.4E-02	7.5E-06	WATER9	1.0E+01	EPI	1.5E+00	PHYSPROP	1.0E+05	PHYSPROP	9.7E-05	2.3E-01	2.0E+00	2.0E-05	EPI							
Decabromodiphenyl ether, 2,2',3',4',4',5',5',6',6'-(BDE-209)	1163-19-5	9.6E+02	PHYSPROP	4.9E-07	1.2E-08	PHYSPROP	4.7E-12	PHYSPROP	3.1E+02	PHYSPROP	3.0E+00	IRIS Profile	1.9E-02	4.8E-06	WATER9	2.8E+05	EPI	1.2E+01	PHYSPROP	1.0E+04	PHYSPROP	8.6E+00	2.7E+01	1.1E+05	7.3E-01	EPI					
Demeton	8065-48-3	3.2E+02	PHYSPROP	1.6E-04	3.8E-06	PHYSPROP	3.4E-04	PHYSPROP	8.5E+01	PHYSPROP	1.1E+00	PubChem	1.1E-02	3.8E-06	WATER9	6.7E+02	EPI	3.2E+00	PHYSPROP	6.7E+02	PHYSPROP	6.6E-02	8.2E+01	2.0E+02	7.6E-03	RAGSE					
Di(2-ethylhexyl)adipate	103-23-1	5.7E+02	PHYSPROP	1.8E-05	4.3E-07	PHYSPROP	8.4E-07	PHYSPROP	-6.8E+01	PHYSPROP	9.2E-01	CRCB9	1.7E-02	4.2E-06	WATER9	3.6E+04	EPI	6.1E+00	PHYSPROP	7.8E+01	PHYSPROP	2.4E+01	1.3E+01	5.8E+01	3.2E+00	EPI					
Diallate	2303-16-4	2.7E+02	PHYSPROP	1.6E-04	3.8E-06	EPI	1.5E-04	PHYSPROP	2.5E+01	PHYSPROP	9.2E-01	CRCB9	4.5E-02	5.3E-06	WATER9	6.4E+02	EPI	4.5E+00	PHYSPROP	1.4E+01	PHYSPROP	2.9E+01	3.4E+00	8.2E+00	4.6E-02	EPI					
Diazinon	333-41-5	3.0E+02	PHYSPROP	4.6E-06	1.1E-07	PHYSPROP	9.0E-05	PHYSPROP	8.8E+01	EPI	1.1E+00	CRCB9	2.1E-02	5.2E-06	WATER9	3.0E+03	EPI	3.8E+00	PHYSPROP	4.0E+01	PHYSPROP	7.0E-02	5.3E+00	1.3E+01	1.0E-02	EPI					
Dibenzothiazophene	132-65-0	1.8E+02	PHYSPROP	1.4E-03	3.4E-05	EPI	2.1E-04	EPI	9.7E+01	PHYSPROP	1.3E+00	ChemNet	3.6E-02	7.6E-06	WATER9	9.2E+03	EPI	4.4E+00	PHYSPROP	1.5E+00	PHYSPROP	6.2E-01	1.1E+00	4.5E+00	1.2E-01	EPI					
Dibromo-3-chloropropane, 1,2-	96-12-8	2.4E+02	PHYSPROP	6.0E-03	1.5E-04	EPI	5.8E-01	PHYSPROP	6.0E+00	PHYSPROP	2.1E+00	CRCB9	3.0E+00	2.8E-06	WATER9	1.2E+02	EPI	3.0E+00	PHYSPROP	1.2E+03	PHYSPROP	4.1E-02	2.2E+00	5.3E+00	6.9E-03	EPI					
Dibromobenzene, 1,4-	108-36-1	2.4E+02	PHYSPROP	5.1E-02	1.2E-03	EPI	2.7E-01	PHYSPROP	-7.0E+00	PHYSPROP	2.0E+00	CRCB9	3.1E-02	8.5E-06	WATER9	3.8E+02	EPI	3.8E+00	PHYSPROP	6.8E+01	PHYSPROP	1.4E-01	2.2E+00	5.3E+00	2.3E-02	EPI					
Dibromobenzene, 1,3-	106-37-6	2.4E+02	PHYSPROP	3.7E-02	8.9E-04	EPI	5.8E-02	PHYSPROP	8.7E+01	PHYSPROP	2.3E+00	CRCB9	3.3E-02	9.3E-06	WATER9	3.8E+02	EPI	3.8E+00	PHYSPROP	2.0E+01	PHYSPROP	1.4E-01	2.2E+00	5.3E+00	2.5E-02	EPI					
Dibromochloromethane	124-48-1	2.1E+02	PHYSPROP	3.2E-02	7.8E-04	PHYSPROP	5.5E+00	PHYSPROP	-2.0E+01	PHYSPROP	2.5E+00	CRCB9	3.7E-02	1.1E-05	WATER9	3.2E+01	EPI	2.2E+00	PHYSPROP	2.7E+03	PHYSPROP	1.6E-02	1.5E+00	3.7E+00	2.9E-03	EPI					
Dibromomethane	106-93-4	1.9E+02	PHYSPROP	2.7E-02	6.5E-04	PHYSPROP	1.1E+01	PHYSPROP	9.9E+00	PHYSPROP	2.2E+00	CRCB9	4.8E-02	1.0E-05	WATER9	4.0E+01	EPI	2.0E+00	PHYSPROP	3.9E+03	PHYSPROP	1.5E-02	1.2E+00	2.8E+00	2.8E-03	EPI					
Dibromomethane (Methylene Bromide)	74-95-3	1.7E+02	PHYSPROP	3.4E-02	8.2E-04	PHYSPROP	4.4E+01	PHYSPROP	-5.3E+01	PHYSPROP	2.5E+00	CRCB9	5.5E-02	1.2E-05	WATER9	2.2E+01	EPI	1.7E+00	PHYSPROP	1.2E+04	PHYSPROP	1.1E-02	9.9E-01	2.4E+00	2.2E-03	EPI					
Dibutyltin Compounds	NA																														
Dicamba	1918-00-9	2.2E+02	PHYSPROP	8.9E-08	2.2E-09	EPI	1.3E-05	PHYSPROP	1.2E+02	PHYSPROP	1.6E+00	CRCB9	2.9E-02	7.8E-06	WATER9	2.9E+01	EPI	2.2E+00	PHYSPROP	8.3E+03	PHYSPROP	1.5E-02	1.8E+00	4.4E+00	2.7E-03	EPI					
Dichloro-2-butene, 1,4-	764-01-0	1.3E+02	PHYSPROP	3.5E-01	8.5E-03	PHYSPROP	3.0E+00	EPI	3.5E+00	PHYSPROP	1.2E+00	LANGE	6.7E-02	9.3E-06	WATER9	1.3E+02	EPI	2.6E+00	PHYSPROP	5.8E+02	PHYSPROP	7.1E-02	5.3E-01	1.3E+00	1.7E-02	EPI					
Dichloro-2-butene, cis-1,4-	1476-11-5	1.3E+02	PHYSPROP	2.7E-02	6.6E-04	EPI	4.1E+00	PHYSPROP	4.8E+01	PHYSPROP	1.2E+00	CRCB9	6.7E-02	9.3E-06	WATER9	1.3E+02	EPI	2.6E+00	PHYSPROP	5.8E+02	PHYSPROP	7.1E-02	5.3E-01	1.3E+00	1.7E-02	EPI					
Dichloro-2-butene, trans-1,4-	110-57-6	1.3E+02	PHYSPROP	2.7E-02	6.6E-04	EPI	3.4E+00	PHYSPROP	2.0E+00	PHYSPROP	1.2E+00	CRCB9	6.6E-02	9.3E-06	WATER9	1.3E+02	EPI	2.6E+00	PHYSPROP	5.5E+02	PHYSPROP	7.1E-02	5.3E-01	1.3E+00	1.7E-02	EPI					
Dichloroacetic Acid	79-43-6	1.3E+02	PHYSPROP	3.4E-07	8.4E-09	PHYSPROP	1.8E-01	PHYSPROP	1.4E+01	PHYSPROP	1.6E+00	CRCB9	7.2E-02	1.1E-05	WATER9	2.3E+00	EPI	9.2E-01	PHYSPROP	1.0E+06	PHYSPROP	5.3E-03	5.5E-01	1.3E+00	1.2E-03	EPI					
Dichlorobenzene, 1,2-	95-50-1	1.5E+02	PHYSPROP	7.8E-02	1.9E-03	PHYSPROP	1.4E+00	PHYSPROP	-1.7E+01	PHYSPROP	1.3E+00	CRCB9	5.6E-02	8.9E-06	WATER9	3.8E+02	EPI	3.4E+00	PHYSPROP	1.6E+02	PHYSPROP	2.1									

Regional Removal Management Level (RML) Chemical-specific Parameters Supporting Table May 2016

Contaminant	Molecular Weight		Volatility Parameters				Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients				Water Solubility		Tapwater Dermal Parameters									
	Analyte	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m ³ /mole)	H' and HLC Ref	VP	VP Ref	MP	MP Ref	Density (g/cm ³)	Density Ref	Dia (cm ² /s)	Diw (cm ² /s)	D _a and D _w Ref	K _{oc} (L/kg)	K _{oc} Ref	K _{oc}	K _{oc} Ref	log K _{ow} (unitless)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/event)	t [*] (hr)	K _p (cm ² /hr)	K Ref	
Dimethylphenol, 3,4-	95-65-8	1.2E+02	PHYSPROP	1.7E-05	4.2E-07	PHYSPROP	3.6E-02	EPI	6.1E+01	PHYSPROP	9.8E-01	CRC89	6.3E-02	8.4E-06	WATER9	4.9E+02	EPI	2.2E+00	PHYSPROP	4.8E+03	PHYSPROP	4.8E+03	PHYSPROP	4.0E+03	PHYSPROP	4.2E-02	5.1E-01	1.2E+00	9.8E-03	EPI
Dithienylvinylchloride	513-37-1	9.1E+01	PHYSPROP	4.8E-02	1.2E-03	CRC89	2.1E+02	EPI	9.2E+01	EPI	9.2E-01	CRC89	8.1E-02	9.7E-06	WATER9	6.1E+01	EPI	2.6E+00	PHYSPROP	4.8E+03	PHYSPROP	4.8E+03	PHYSPROP	4.0E+03	PHYSPROP	9.3E-02	3.4E-01	8.1E-01	2.5E-02	EPI
Dinitro-o-cresol, 4,6-	534-52-1	2.0E+02	PHYSPROP	5.7E-05	1.4E-06	PHYSPROP	1.2E-04	PHYSPROP	8.7E+01	PHYSPROP	1.3E+00	CRC89	5.6E-02	6.5E-06	WATER9	7.5E+02	EPI	2.1E+00	PHYSPROP	2.0E+02	PHYSPROP	2.0E+02	PHYSPROP	2.0E+02	PHYSPROP	1.7E-02	1.4E+00	3.2E+00	3.2E-03	EPI
Dinitro-o-cyclohexyl Phenol, 4-	131-89-5	2.7E+02	PHYSPROP	4.2E-08	5.5E-08	PHYSPROP	4.2E-08	PHYSPROP	1.1E+02	PHYSPROP	1.3E+00	CRC89	4.6E-02	5.4E-06	WATER9	1.7E+04	EPI	4.1E+00	PHYSPROP	1.5E+01	PHYSPROP	1.5E+01	PHYSPROP	1.5E+01	PHYSPROP	1.7E-01	3.3E+00	7.8E+00	2.8E-02	EPI
Dinitrobenzene, 1,2-	528-29-0	1.7E+02	PHYSPROP	2.2E-06	5.3E-08	EPI	4.6E-05	EPI	1.2E+02	PHYSPROP	1.3E+00	CRC89	4.5E-02	8.3E-06	WATER9	3.6E+02	EPI	1.7E+00	PHYSPROP	3.1E+02	PHYSPROP	3.1E+02	PHYSPROP	3.1E+02	PHYSPROP	4.2E-02	9.2E-01	2.2E+00	2.4E-03	EPI
Dinitrobenzene, 1,3-	99-65-0	1.7E+02	PHYSPROP	2.0E-06	4.9E-08	PHYSPROP	9.0E-04	EPI	9.0E+01	PHYSPROP	1.6E+00	CRC89	4.8E-02	9.2E-06	WATER9	3.5E+02	EPI	1.5E+00	PHYSPROP	5.3E+02	PHYSPROP	5.3E+02	PHYSPROP	5.3E+02	PHYSPROP	6.7E-03	9.2E-01	2.2E+00	1.7E-03	EPI
Dinitrobenzene, 1,4-	100-25-4	1.7E+02	PHYSPROP	3.4E-05	8.4E-08	PHYSPROP	3.4E-05	PHYSPROP	1.7E+02	PHYSPROP	1.6E+00	CRC89	4.9E-02	9.4E-06	WATER9	3.5E+02	EPI	1.5E+00	PHYSPROP	6.9E+01	PHYSPROP	6.9E+01	PHYSPROP	6.9E+01	PHYSPROP	8.3E-03	9.2E-01	2.2E+00	1.7E-03	EPI
Dinitrophenol, 2,4-	51-28-5	1.8E+02	PHYSPROP	3.5E-06	8.6E-08	PHYSPROP	3.9E-04	PHYSPROP	1.1E+02	PHYSPROP	1.7E+00	CRC89	4.1E-02	9.1E-06	WATER9	4.6E+02	EPI	1.7E+00	PHYSPROP	2.8E+03	PHYSPROP	2.8E+03	PHYSPROP	2.8E+03	PHYSPROP	9.8E-03	1.1E+00	2.7E+00	1.9E-03	EPI
Dinitrotoluene Mixture, 2,4/2,6	NA	1.8E+02	EPI	1.6E-05	4.0E-07	EPI	2.2E-03	EPI	6.0E+01	EPI	1.3E+00	CRC89	5.9E-02	6.9E-06	WATER9	2.8E+02	EPI	2.2E+00	EPI	2.7E+02	PHYSPROP	2.7E+02	PHYSPROP	2.7E+02	PHYSPROP	2.2E-02	1.1E+00	2.6E+00	4.2E-03	EPI
Dinitrotoluene, 2,4-	121-14-2	1.8E+02	PHYSPROP	2.2E-06	5.4E-08	PHYSPROP	1.5E-04	PHYSPROP	7.1E+01	PHYSPROP	1.3E+00	CRC89	3.8E-02	7.9E-06	WATER9	5.8E+02	EPI	2.0E+00	PHYSPROP	2.0E+02	PHYSPROP	2.0E+02	PHYSPROP	2.0E+02	PHYSPROP	1.6E-02	1.1E+00	2.6E+00	3.1E-03	EPI
Dinitrotoluene, 2,6-	606-20-2	1.8E+02	PHYSPROP	3.1E-05	7.5E-07	EPI	5.7E-04	PHYSPROP	6.6E+01	PHYSPROP	1.3E+00	CRC89	3.7E-02	7.8E-06	WATER9	5.9E+02	EPI	2.1E+00	PHYSPROP	1.8E+02	PHYSPROP	1.8E+02	PHYSPROP	1.8E+02	PHYSPROP	1.9E-02	1.1E+00	2.6E+00	3.7E-03	EPI
Dinitrotoluene, 2-Amino-4,6-	35572-78-2	2.0E+02	PHYSPROP	1.3E-09	3.3E-11	PHYSPROP	1.1E-05	PHYSPROP	1.7E+02	PHYSPROP	1.2E+00	CRC89	5.6E-02	6.6E-06	WATER9	2.8E+02	EPI	1.8E+00	PHYSPROP	1.2E+03	PHYSPROP	1.2E+03	PHYSPROP	1.2E+03	PHYSPROP	1.1E-02	1.3E+00	3.2E+00	2.0E-03	EPI
Dinitrotoluene, 4-Amino-2,6-	19406-51-0	2.0E+02	PHYSPROP	1.3E-09	3.3E-11	PHYSPROP	1.1E-05	PHYSPROP	1.7E+02	PHYSPROP	1.2E+00	CRC89	5.6E-02	6.6E-06	WATER9	2.8E+02	EPI	1.8E+00	PHYSPROP	1.2E+03	PHYSPROP	1.2E+03	PHYSPROP	1.2E+03	PHYSPROP	1.1E-02	1.3E+00	3.2E+00	2.0E-03	EPI
Dinitrotoluene, Technical grade	25321-14-6	5.5E+02	PHYSPROP	3.8E-06	9.3E-08	PHYSPROP	4.0E-04	PHYSPROP	6.0E+01	EPI	1.3E+00	CRC89	2.8E-02	6.3E-06	WATER9	5.9E+02	EPI	2.2E+00	PHYSPROP	2.7E+02	PHYSPROP	2.7E+02	PHYSPROP	2.7E+02	PHYSPROP	3.7E-02	1.2E+00	2.9E+02	4.2E-03	EPI
Dinoseb	88-85-7	2.4E+02	PHYSPROP	1.9E-05	4.6E-07	EPI	7.5E-05	PHYSPROP	4.0E+01	PHYSPROP	1.3E+00	CRC89	2.5E-02	3.5E-06	WATER9	4.3E+03	EPI	3.6E+00	PHYSPROP	5.2E+01	PHYSPROP	5.2E+01	PHYSPROP	5.2E+01	PHYSPROP	9.7E-02	2.3E+00	6.6E+00	1.6E-02	EPI
Dioxane, 1,4-	123-91-1	8.8E+01	PHYSPROP	2.0E-04	4.8E-06	PHYSPROP	3.8E+01	PHYSPROP	1.2E+01	PHYSPROP	1.0E+00	CRC89	8.7E-02	1.1E-05	WATER9	2.6E+00	EPI	-2.7E-01	PHYSPROP	1.0E+06	PHYSPROP	1.0E+06	PHYSPROP	1.0E+06	PHYSPROP	1.2E-03	3.3E-01	7.9E-01	3.3E-04	EPI
**Hexachlorodibenzo-p-dioxin, Mixture	NA	3.9E+02	EPI	2.3E-04	5.7E-06	EPI	4.4E-11	EPI	2.5E+02	EPI	1.0E+00	CRC89	4.3E-02	4.2E-06	WATER9	7.0E+05	EPI	8.2E+00	EPI	4.0E-06	EPI	4.0E-06	EPI	4.0E-06	EPI	2.2E+01	1.6E+01	7.5E-01	2.9E+00	EPI
**TCDD, 2,3,7,8-	1746-01-6	3.2E+02	PHYSPROP	2.0E-03	5.0E-05	EPI	1.5E-09	PHYSPROP	3.1E+02	PHYSPROP	1.8E+00	PubChem	4.7E-02	6.8E-06	WATER9	2.5E+05	EPI	6.8E+00	PHYSPROP	2.0E-04	PHYSPROP	2.0E-04	PHYSPROP	2.0E-04	PHYSPROP	5.6E+00	6.7E+00	2.9E+01	8.1E-01	EPI
Diphenamid	957-51-7	2.4E+02	PHYSPROP	1.5E-09	3.6E-11	EPI	3.0E-08	PHYSPROP	1.4E+02	PHYSPROP	1.2E+00	CRC89	2.4E-02	6.2E-06	WATER9	4.8E+03	EPI	2.2E+00	PHYSPROP	2.6E+02	PHYSPROP	2.6E+02	PHYSPROP	2.6E+02	PHYSPROP	3.3E-02	2.3E+00	5.5E+00	5.6E-03	EPI
Diphenyl Sulfone	127-63-9	2.2E+02	PHYSPROP	1.0E-05	2.5E-07	PHYSPROP	1.5E-05	PHYSPROP	1.3E+02	PHYSPROP	1.3E+00	CRC89	2.7E-02	6.9E-06	WATER9	1.1E+03	EPI	2.4E+00	PHYSPROP	3.1E+02	PHYSPROP	3.1E+02	PHYSPROP	3.1E+02	PHYSPROP	2.1E-02	1.8E+00	4.2E+00	3.7E-03	EPI
Diphenylamine	12239-4	1.7E+02	PHYSPROP	1.1E-04	2.7E-06	EPI	6.7E-04	PHYSPROP	5.3E+01	PHYSPROP	1.2E+00	CRC89	4.2E-02	7.6E-06	WATER9	8.3E+02	EPI	3.5E+00	PHYSPROP	5.3E+01	PHYSPROP	5.3E+01	PHYSPROP	5.3E+01	PHYSPROP	1.9E-01	9.3E-01	3.2E+00	3.7E-02	EPI
Diphenylhydrazine, 1,2-	122-66-7	1.8E+02	PHYSPROP	2.0E-05	4.8E-07	EPI	4.4E-04	EPI	1.3E+02	PHYSPROP	1.2E+00	CRC89	3.4E-02	7.2E-06	WATER9	1.5E+03	EPI	2.9E+00	PHYSPROP	2.2E+02	PHYSPROP	2.2E+02	PHYSPROP	2.2E+02	PHYSPROP	6.8E-02	1.1E+00	2.7E+00	1.3E-02	EPI
Diquat	85-00-7	3.4E+02	PHYSPROP	1.8E-12	1.4E-13	PHYSPROP	1.8E-06	PHYSPROP	3.4E+02	PHYSPROP	1.2E+00	CRC89	2.1E-02	5.2E-06	WATER9	9.3E+03	EPI	-4.6E+00	PHYSPROP	7.1E+05	PHYSPROP	7.1E+05	PHYSPROP	7.1E+05	PHYSPROP	1.7E-06	8.9E+00	2.1E+01	2.4E-07	EPI
Direct Black 38	1937-37-7	7.8E+02	PHYSPROP	3.4E-38	8.2E-40	PHYSPROP	1.5E-36	PHYSPROP	3.5E+02	EPI	1.1E+00	CRC89	2.2E-02	2.6E-06	WATER9	2.4E+08	EPI	4.9E+00	PHYSPROP	3.0E+03	PHYSPROP	3.0E+03	PHYSPROP	3.0E+03	PHYSPROP	2.2E-03	2.4E+03	5.9E+03	2.1E-04	EPI
Direct Blue 6	2602-46-2	9.3E+02	PHYSPROP	3.7E-42	9.1E-44	PHYSPROP	9.5E-39	PHYSPROP	3.5E+02	EPI	1.1E+00	CRC89	2.0E-02	2.3E-06	WATER9	7.9E+08	EPI	2.6E+00	PHYSPROP	1.4E-04	PHYSPROP	1.4E-04	PHYSPROP	1.4E-04	PHYSPROP	2.0E-08	1.8E+04	4.2E+04	1.7E-09	EPI
Direct Brown 95	16071-86-6	7.6E+02	PHYSPROP	1.4E-41	PHYSPROP	PHYSPROP	1.4E-41	PHYSPROP	3.5E+02	EPI	1.1E+00	CRC89	2.3E-02	2.7E-06	WATER9	7.0E+06	EPI	-6.5E+00	PHYSPROP	1.0E+06	PHYSPROP	1.0E+06	PHYSPROP	1.0E+06	PHYSPROP	4.1E-11	1.9E+03	4.6E+03	3.9E-12	EPI
Disulfoton	298-04-4	2.7E+02	PHYSPROP	8.8E-05	2.2E-06	EPI	9.8E-05	PHYSPROP	1.2E+02	PHYSPROP	1.1E+00	CRC89	2.3E-02	5.7E-06	WATER9	8.4E+02	EPI	4.0E+00	PHYSPROP	1.6E+01	PHYSPROP	1.6E+01	PHYSPROP	1.6E+01	PHYSPROP	1.4E-01	3.6E+00	8.7E+00	2.1E-02	EPI
Dithiane, 1,4-	505-29-3	1.2E+02	PHYSPROP	1.7E-03	4.2E-05	EPI	8.0E-02	PHYSPROP	1.1E+02	PHYSPROP	1.1E+00	ChemNet	6.8E-02	9.3E-06	WATER9	1.5E+02	EPI	7.7E-01	PHYSPROP	3.0E+03	PHYSPROP	3.0E+03	PHYSPROP	3.0E+03	PHYSPROP	4.6E-03	5.0E-01	1.2E+00	1.1E-03	EPI
Diuron	330-54-1	2.3E+02	PHYSPROP	2.1E-08	5.0E-10	EPI	6.9E-08	PHYSPROP	1.6E+02	PHYSPROP	1.1E+00	CRC89	5.0E-02	5.9E-06	WATER9	1.1E+02	EPI	2.7E+00	PHYSPROP	4.2E+01	PHYSPROP	4.2E+01	PHYSPROP	4.2E+01	PHYSPROP	2.7E-02	2.1E+00	5.1E+00	4.7E-03	EPI
Dodine	2439-10-3	2.9E+02	PHYSPROP	3.7E-09	9.0E-11	EPI	1.5E-07	PHYSPROP	1.4E+02	PHYSPROP	1.1E+00	CRC89	4.4E-02	5.1E-06	WATER9	2.5E+03	EPI	1.2E+00	PHYSPROP	6.3E+02	PHYSPROP	6.3E+02	PHYSPROP	6.3E+02	PHYSPROP	1.4E-03	4.3E+00	1.0E+01	2.2E-04	EPI
EPTC	759-94-4	1.9E+02	PHYSPROP	6.5E-04	1.6E-05	EPI	2.4E-02	PHYSPROP	6.1E+01	EPI	9.5E-01	CRC89	2.9E-02	6.4E-06	WATER9	1.6E+02	EPI	3.2E+00	PHYSPROP	3.8E+02	PHYSPROP	3.8E+02	PHYSPROP	3.8E+02	PHYSPROP	9.7E-02	1.2E+00	2.6E+00	1.8E-02	EPI
Endosulfan	115-29-7	4.1E+02	PHYSPROP	2.7E-03	6.5E-05	PHYSPROP	1.7E-07	PHYSPROP	1.1E+02	PHYSPROP	1.7E+00	CRC89	2.2E-02	5.8E-06	WATER9	6.8E+03	EPI	3.8E+00	PHYSPROP	3.3E-01	PHYSPROP	3.3E-01	PHYSPROP	3.3E-01	PHYSPROP	2.2E-02	2.0E+01	4.8E+01	2.9E-03	EPI
Endothal	145-73-3	1.9E+02	PHYSPROP	1.6E-14	3.9E-16	EPI	1.6E-10	PHYSPROP	1.4E+02	PHYSPROP	1.4E+00	CRC89	3.7E-02	8.2E-06	WATER9	1.9E+01	EPI	1.9E+00	PHYSPROP	1.0E+05	PHYSPROP	1.0E+05	PHYSPROP	1.0E+05	PHYSPROP	1.4E-02	1.2E+00	2.8E+00	2.6E-03	EPI
Endrin	72-20-8	3.8E+02	PHYSPROP</																											

Regional Removal Management Level (RML) Chemical-specific Parameters Supporting Table May 2016

Contaminant	Molecular Weight		Volatility Parameters				Melting Point		Density		Diffusivity in Air and Water			Partition Coefficients				Water Solubility		Tapwater Dermal Parameters										
	Analyle	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m ³ /mole)	H' and HLC Ref	VP	VP Ref	MP	MP Ref	Density (g/cm ³)	Density Ref	Dia (cm ² /s)	Diw (cm ² /s)	D _a and D _w Ref	K _{oc} (L/kg)	K _{oc} Ref	K _{oc}	K _{oc} Ref	log K _{ow} (unitless)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/event)	t _h (hr)	K _p (cm ² /hr)	K Ref	
Guanine	113-00-8	5.9E+01	PHYSPROP	9.6E-10	2.3E-11	PHYSPROP	2.2E+00	PHYSPROP	5.0E+01	PHYSPROP	1.6E+00	GuideChem	1.4E-01	1.7E-05	WATER9	1.2E+01	1.8E+01	1.6E+00	PHYSPROP	1.8E+03	PHYSPROP	1.0E+06	PHYSPROP	1.0E+03	PHYSPROP	1.8E-04	2.3E-01	5.4E-01	6.0E-05	EPI
Guandine Chloride	50-01-1	9.6E+01	PHYSPROP	9.8E-17	2.2E-18	PHYSPROP	1.8E-06	PHYSPROP	1.8E+02	PHYSPROP	1.4E+00	CRC89	9.2E-02	1.2E-05	WATER9	1.4E+00	1.8E+01	1.6E+00	PHYSPROP	1.0E+06	PHYSPROP	1.0E+06	PHYSPROP	1.0E+03	PHYSPROP	1.5E-07	3.6E-01	8.7E-01	3.9E-08	EPI
Haloxafop, Methyl	69806-40-2	3.8E+02	PHYSPROP	1.3E-05	3.2E-07	EPI	6.0E-06	PHYSPROP	5.6E+01	PHYSPROP	1.6E+00	CRC89	3.6E-02	4.3E-06	WATER9	1.6E+00	1.8E+01	1.6E+00	PHYSPROP	5.5E+03	EPI	4.1E+00	PHYSPROP	9.9E+00	PHYSPROP	4.5E-02	1.3E+01	3.2E+01	6.0E-03	EPI
Heptachlor	76-44-8	3.7E+02	PHYSPROP	1.2E-02	2.9E-04	PHYSPROP	4.0E-04	PHYSPROP	9.6E+01	PHYSPROP	1.6E+00	CRC89	2.2E-02	5.7E-06	WATER9	1.6E+00	1.8E+01	1.6E+00	PHYSPROP	4.1E+04	EPI	6.1E+00	PHYSPROP	9.8E-01	PHYSPROP	1.1E+00	1.3E+01	5.0E+01	1.4E-01	EPI
Heptachlor Epoxide	1024-57-3	3.9E+02	PHYSPROP	8.6E-04	2.1E-05	PHYSPROP	2.0E-05	PHYSPROP	1.6E+02	PHYSPROP	1.9E+00	LookChem	2.4E-02	6.2E-06	WATER9	1.9E+00	2.4E+01	1.6E+00	PHYSPROP	1.0E+04	EPI	5.0E+00	PHYSPROP	2.0E-01	PHYSPROP	1.6E-01	1.6E+01	3.8E+01	2.1E-02	EPI
Hexabromobenzene	87-82-1	5.5E+02	PHYSPROP	1.6E-08	2.8E-05	PHYSPROP	1.6E-08	PHYSPROP	1.3E+02	PHYSPROP	3.0E+00	LookChem	2.5E-02	6.6E-06	WATER9	3.0E+00	3.6E+01	1.6E+00	PHYSPROP	1.6E-04	PHYSPROP	1.6E-04	PHYSPROP	1.2E-01	PHYSPROP	1.3E+02	3.1E+02	1.1E-02	1.4E-02	EPI
Hexabromodiphenyl ether, 2,2',4,4',5,5'-(BDE-153)	68631-49-2	6.4E+02	OTHER				5.8E-06	IRIS Profile							WATER9															
Hexachlorobenzene	118-74-1	2.8E+02	PHYSPROP	7.0E-02	1.7E-03	PHYSPROP	1.8E-05	PHYSPROP	3.3E+02	PHYSPROP	2.0E+00	CRC89	2.9E-02	7.8E-06	WATER9	2.0E+00	2.4E+01	1.6E+00	PHYSPROP	6.2E+03	EPI	5.7E+00	PHYSPROP	6.2E-03	PHYSPROP	1.6E+00	4.1E+00	1.7E+01	2.5E-01	EPI
Hexachlorobutadiene	87-68-3	2.6E+02	PHYSPROP	4.2E-01	1.0E-02	PHYSPROP	2.2E-01	PHYSPROP	3.2E+01	PHYSPROP	1.6E+00	CRC89	2.7E-02	7.0E-06	WATER9	1.6E+00	2.0E+01	1.6E+00	PHYSPROP	8.5E+02	EPI	4.8E+00	PHYSPROP	3.2E+00	PHYSPROP	5.0E-01	3.0E+00	7.3E+00	8.1E-02	EPI
Hexachlorocyclohexane, Alpha-	319-84-6	2.9E+02	PHYSPROP	2.7E-04	6.7E-06	PHYSPROP	3.5E-05	EPI	1.6E+02	PHYSPROP	1.9E+00	CRC89	4.3E-02	5.1E-06	WATER9	1.9E+00	2.4E+01	1.6E+00	PHYSPROP	2.8E+03	EPI	3.8E+00	PHYSPROP	2.0E+00	PHYSPROP	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclohexane, Beta-	319-85-7	2.9E+02	PHYSPROP	1.8E-05	4.4E-07	PHYSPROP	3.6E-07	PHYSPROP	3.1E+02	PHYSPROP	1.9E+00	CRC89	2.8E-02	7.4E-06	WATER9	1.9E+00	2.4E+01	1.6E+00	PHYSPROP	2.8E+03	EPI	3.8E+00	PHYSPROP	2.4E-01	PHYSPROP	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclohexane, Gamma-(Lindane)	58-89-9	2.9E+02	PHYSPROP	2.1E-04	5.1E-06	PHYSPROP	4.2E-05	PHYSPROP	1.1E+02	PHYSPROP	1.7E+00	CRC89	4.3E-02	5.1E-06	WATER9	1.7E+00	2.1E+01	1.6E+00	PHYSPROP	2.8E+03	EPI	3.7E+00	PHYSPROP	7.3E+00	PHYSPROP	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclohexane, Technical	608-73-1	2.9E+02	PHYSPROP	2.1E-04	5.1E-06	EPI	3.5E-05	EPI	1.1E+02	EPI	1.7E+00	CRC89	4.3E-02	5.1E-06	WATER9	1.7E+00	2.1E+01	1.6E+00	PHYSPROP	2.8E+03	EPI	4.1E+00	EPI	8.0E+00	PHYSPROP	1.4E-01	4.5E+00	1.1E+01	2.1E-02	EPI
Hexachlorocyclopentadiene	77-47-4	2.7E+02	PHYSPROP	1.1E+00	2.7E-02	PHYSPROP	6.0E-02	PHYSPROP	9.0E+00	PHYSPROP	1.7E+00	CRC89	2.7E-02	7.2E-06	WATER9	1.7E+00	2.1E+01	1.6E+00	PHYSPROP	1.4E+03	EPI	5.0E+00	PHYSPROP	1.8E+00	PHYSPROP	6.5E-01	3.5E+00	1.4E+01	1.0E-01	EPI
Hexachloroethane	67-72-1	2.4E+02	PHYSPROP	1.6E-01	3.9E-03	PHYSPROP	2.1E-01	PHYSPROP	1.9E+02	PHYSPROP	2.1E+00	CRC89	3.2E-02	8.9E-06	WATER9	2.1E+00	2.5E+01	1.6E+00	PHYSPROP	2.0E+02	EPI	4.1E+00	PHYSPROP	5.0E+01	PHYSPROP	2.5E-01	2.2E+00	5.3E+00	4.0E-02	EPI
Hexachlorophene	70-30-4	4.1E+02	PHYSPROP	2.2E-11	5.5E-13	PHYSPROP	1.0E-10	PHYSPROP	1.7E+02	PHYSPROP	1.7E+00	CRC89	3.5E-02	4.0E-06	WATER9	1.7E+00	2.1E+01	1.6E+00	PHYSPROP	6.7E+05	EPI	7.5E+00	PHYSPROP	1.4E+02	PHYSPROP	6.5E+00	2.0E+01	8.9E+01	8.4E-01	EPI
Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	2.2E+02	PHYSPROP	8.2E-10	2.0E-11	EPI	4.1E-09	EPI	2.1E+02	PHYSPROP	1.8E+00	CRC89	3.1E-02	8.5E-06	WATER9	1.8E+00	2.2E+01	1.6E+00	PHYSPROP	8.9E+01	EPI	8.7E-01	PHYSPROP	6.0E+01	PHYSPROP	1.9E-03	1.0E+00	4.4E+00	3.4E-04	EPI
Hexamethylene Diisocyanate, 1,6-Hexamethylolurea	822-06-0	1.7E+02	PHYSPROP	2.0E-03	4.8E-05	PHYSPROP	3.0E-02	PHYSPROP	6.7E+01	PHYSPROP	1.1E+00	CRC89	4.0E-02	7.2E-06	WATER9	1.1E+00	1.4E+01	1.6E+00	PHYSPROP	4.8E+03	EPI	3.2E+00	PHYSPROP	1.2E+02	PHYSPROP	1.2E-01	9.2E-01	2.2E+00	2.4E-02	EPI
Hexamethylolurea	680-31-9	1.8E+02	PHYSPROP	8.6E-07	2.0E-08	PHYSPROP	4.6E-02	PHYSPROP	7.2E+00	PHYSPROP	1.0E+00	CRC89	3.5E-02	6.9E-06	WATER9	1.0E+00	1.3E+01	1.6E+00	PHYSPROP	1.0E+01	EPI	2.8E-01	PHYSPROP	1.0E+06	PHYSPROP	1.2E-03	1.1E+01	2.5E+00	2.4E-04	EPI
Hexane, N-	110-54-3	8.6E+01	PHYSPROP	7.4E+01	1.8E+00	EPI	1.5E+02	PHYSPROP	9.5E+01	PHYSPROP	6.6E-01	CRC89	7.3E-02	8.2E-06	WATER9	6.6E-01	8.0E+00	1.6E+00	PHYSPROP	1.3E+02	EPI	3.9E+00	PHYSPROP	9.5E+00	PHYSPROP	7.2E-01	3.2E-01	1.2E+00	2.0E-01	EPI
Hexachloroacid	124-04-9	1.5E+02	PHYSPROP	1.9E-10	4.7E-12	EPI	3.2E-07	EPI	1.5E+02	PHYSPROP	1.4E+00	CRC89	1.8E-02	4.6E-06	WATER9	1.4E+00	1.7E+01	1.6E+00	PHYSPROP	3.1E+04	PHYSPROP	3.1E+04	PHYSPROP	1.2E-02	PHYSPROP	4.1E-01	3.7E+00	2.7E-04	2.7E-04	EPI
Hexane, 2-	591-78-6	1.0E+02	PHYSPROP	3.8E-03	9.3E-05	EPI	1.2E+01	PHYSPROP	5.6E+01	PHYSPROP	8.1E-01	CRC89	7.0E-02	8.4E-06	WATER9	8.1E-01	9.7E+00	1.6E+00	PHYSPROP	1.5E+01	EPI	1.4E+00	PHYSPROP	1.7E+04	PHYSPROP	1.4E-02	9.8E-01	9.2E-01	3.6E-03	EPI
Hexane, 2,2,4-trimethyl-	51235-04-2	2.5E+02	PHYSPROP	9.2E-11	2.3E-12	EPI	2.3E-07	EPI	1.2E+02	PHYSPROP	1.3E+00	CRC89	2.5E-02	6.3E-06	WATER9	1.3E+00	1.6E+01	1.6E+00	PHYSPROP	1.1E+02	EPI	1.9E+00	PHYSPROP	3.3E+04	PHYSPROP	6.2E-03	2.7E+00	6.5E+00	1.0E-03	EPI
Hexylthiazole	78587-05-0	3.5E+02	PHYSPROP	9.7E-07	2.4E-08	EPI	2.6E-08	PHYSPROP	1.1E+02	PHYSPROP	1.3E+00	CRC89	3.8E-02	4.4E-06	WATER9	1.3E+00	1.6E+01	1.6E+00	PHYSPROP	2.1E+03	EPI	5.6E+00	PHYSPROP	5.0E-01	PHYSPROP	6.0E-01	1.0E+01	2.4E+01	8.3E-02	EPI
Hydramethylnon	67485-29-4	4.9E+02	PHYSPROP	9.0E-05	2.2E-06	EPI	2.0E-08	PHYSPROP	1.9E+02	PHYSPROP	1.3E+00	CRC89	3.0E-02	3.6E-06	WATER9	1.3E+00	1.6E+01	1.6E+00	PHYSPROP	1.8E+08	EPI	2.3E+00	PHYSPROP	6.0E-03	PHYSPROP	7.7E-04	6.2E+01	1.5E+02	9.0E-05	EPI
Hydrazine	302-01-2	3.2E+01	PHYSPROP	2.5E-05	6.1E-07	PubChem	1.4E+01	PHYSPROP	2.0E+00	PHYSPROP	1.0E+00	CRC89	1.7E-01	1.9E-05	WATER9	1.0E+00	1.3E+01	1.6E+00	PHYSPROP	-2.1E+00	PHYSPROP	1.0E+06	PHYSPROP	1.0E+06	PHYSPROP	9.5E-05	1.6E-01	3.8E-01	4.4E-05	RAGSE
Hydrazine Sulfate	10034-93-2	1.3E+02	EPI																											
Hydrogen Chloride	7647-01-0	3.5E+01	EPI	8.3E-07	2.0E+06	Toxnet HSDB	3.5E+04	PubChem	-1.1E+02	CRC89	1.5E+00	CRC89	1.9E-01	2.3E-05	WATER9	1.5E+00	1.8E+01	1.6E+00	PHYSPROP	6.7E+05	Toxnet HSDB	3.1E+04	PHYSPROP	1.4E+01	PHYSPROP	2.4E-03	5.5E-01	1.3E+00	1.0E-03	RAGSE
Hydrogen Fluoride	7664-39-3	2.0E+01	PHYSPROP	4.3E-03	1.0E-04	PHYSPROP	9.2E+02	PHYSPROP	-8.4E+01	PHYSPROP	8.2E-01	CRC89	2.2E-01	2.2E-05	WATER9	8.2E-01	9.7E+00	1.6E+00	PHYSPROP	2.3E-01	OTHER	1.0E+06	PHYSPROP	1.7E-04	1.4E-01	3.3E-01	1.0E-03	RAGSE		
Hydrogen Sulfide	7783-06-4	3.4E+01	PHYSPROP	3.5E-01	8.6E-03	PHYSPROP	1.6E+04	PHYSPROP	-8.5E+01	PHYSPROP	1.4E+00	CRC89	1.9E-01	2.2E-05	WATER9	1.4E+00	1.7E+01	1.6E+00	PHYSPROP	2.3E-01	OTHER	1.0E+06	PHYSPROP	2.2E-03	1.6E-01	3.9E-01	1.0E-03	RAGSE		
Hydroquinone	123-31-9	1.1E+02	PHYSPROP	1.9E-09	4.7E-11	EPI	2.4E-05	EPI	1.7E+02	PHYSPROP	1.3E+00	CRC89	8.0E-02	1.1E-05	WATER9	1.3E+00	1.6E+01	1.6E+00	PHYSPROP	2.4E+02	EPI	5.9E-01	PHYSPROP	7.2E+04	PHYSPROP	3.8E-03	4.3E-01	1.0E+00	9.3E-04	EPI
Imazalil	35554-44-0	3.0E+02	PHYSPROP	1.1E-07	2.6E-09	EPI	1.2E-06	PHYSPROP	5.3E+01	PHYSPROP	1.2E+00	CRC89	2.2E-02	5.7E-06	WATER9	1.2E+00	1.5E+01	1.6E+00	PHYSPROP	8.5E+03	EPI	3.8E+00	PHYSPROP	9.0E+01	PHYSPROP	7.7E-02	4.9E-01	1.2E+01	1.2E-02	EPI
Imazaquin	81335-37-7	3.1E+02	PHYSPROP	2.8E-16	6.9E-18	PHYSPROP	1.0E-13	PHYSPROP	2.2E+02	PHYSPROP	1.2E+00	CRC89	4.1E-02	4.8E-06	WATER9	1.2E+00	1.5E+01	1.6E+00	PHYSPROP	9.0E+01	EPI	1.9E+00	PHYSPROP	9.0E+01	PHYSPROP	3.3E-03	5.8E+00	1.4E+01	4.8E-04	EPI
Imazethapyr	81335-77-5	2.9E+02	PHYSPROP	4.3E-15	1.0E-16	PHYSPROP	2.2E-11	PHYSPROP	1.7E+02	PHYSPROP	1.2E+00	CRC89	4.3E-02	5.1E-06	WATER9															

Regional Removal Management Level (RML) Chemical-specific Parameters Supporting Table May 2016

Contaminant	Molecular Weight		Volatility Parameters						Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients				Water Solubility		Tapwater Dermal Parameters						
	Analyte	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m ³ /mole)	H' and HLC Ref	VP	VP Ref	MP	MP Ref	Density (g/cm ³)	Density Ref	Dia (cm ² /s)	Diw (cm ² /s)	D _a and D _w Ref	K _{oc} (L/kg)	K _{oc} Ref	K _{oc}	K _{oc} Ref	log K _{ow} (unitless)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/event)	t* (hr)	K _p (cm ² /hr)	K _p Ref
Methyl Acrylate	96-33-3	8.6E+01	PHYSPROP	8.1E-03	2.0E-04	EPI	8.7E+01	PHYSPROP	-7.7E+01	PHYSPROP	9.5E-01	CR89	8.6E-02	1.0E-05	WATER9	5.8E+00	EPI	8.0E-01	PHYSPROP	4.9E+04	PHYSPROP	4.9E+04	PHYSPROP	6.2E-03	3.2E-01	7.7E-01	1.8E-03	EPI	
Methyl Ethyl Ketone (2-Butanone)	78-93-3	7.2E+01	PHYSPROP	2.3E-03	5.7E-05	EPI	9.1E+01	PHYSPROP	-8.7E+01	PHYSPROP	8.0E-01	CR89	9.1E-02	1.0E-05	WATER9	4.5E+00	EPI	2.9E-01	PHYSPROP	2.2E+05	PHYSPROP	2.2E+05	PHYSPROP	3.1E-03	2.7E-01	6.4E-01	9.6E-04	EPI	
Methyl Hydrazine	60-34-4	4.6E+01	PHYSPROP	1.2E-04	3.0E-06	PHYSPROP	5.0E+01	PHYSPROP	-5.2E+01	PHYSPROP	8.7E-01	LANGE	1.3E-01	1.4E-05	WATER9	1.3E+01	EPI	1.1E+00	PHYSPROP	1.0E+06	PHYSPROP	1.0E+06	PHYSPROP	4.5E-04	1.9E-01	4.6E-01	1.7E-04	EPI	
Methyl isobutyl Ketone (4-methyl-2-pentanone)	108-10-1	1.0E+02	PHYSPROP	1.5E-03	1.4E-04	EPI	2.0E+01	PHYSPROP	-8.4E+01	PHYSPROP	8.0E-01	CR89	7.0E-02	8.3E-06	WATER9	1.3E+01	EPI	1.3E+00	PHYSPROP	1.9E+04	PHYSPROP	1.9E+04	PHYSPROP	1.2E-02	3.8E-01	9.2E-01	3.2E-03	EPI	
Methyl isocyanate	624-83-9	5.7E+01	PHYSPROP	3.8E-02	9.3E-04	PHYSPROP	3.5E+02	PHYSPROP	-4.5E+01	PHYSPROP	9.6E-01	CR89	1.2E-01	1.3E-05	WATER9	4.0E+01	EPI	7.9E-01	PHYSPROP	2.9E+04	PHYSPROP	2.9E+04	PHYSPROP	7.3E-03	2.2E-01	5.3E-01	2.5E-03	EPI	
Methyl Methacrylate	80-62-4	1.0E+02	PHYSPROP	1.3E-02	3.2E-04	EPI	1.0E+01	PHYSPROP	-8.8E+01	PHYSPROP	9.4E-01	CR89	7.5E-02	9.2E-06	WATER9	9.1E+00	EPI	1.4E+00	PHYSPROP	1.5E+04	PHYSPROP	1.5E+04	PHYSPROP	1.4E-02	3.8E-01	9.2E-01	3.6E-03	EPI	
Methyl Parathion	298-00-0	2.6E+02	PHYSPROP	4.1E-06	1.0E-07	PHYSPROP	3.5E+06	PHYSPROP	1.6E+01	PHYSPROP	1.4E+00	CR89	2.5E-02	5.4E-06	WATER9	7.3E+02	EPI	2.9E+00	PHYSPROP	3.8E+01	PHYSPROP	3.8E+01	PHYSPROP	1.6E-02	3.1E+00	7.5E+00	4.2E-03	EPI	
Methyl Phosphonic Acid	993-13-5	9.6E+01	PHYSPROP	5.0E-10	1.2E-11	PHYSPROP	3.3E-04	EPI	1.1E+02	PHYSPROP			9.1E-02	1.1E-05	WATER9	1.4E+00	EPI	7.0E-01	PHYSPROP	2.0E+04	PHYSPROP	2.0E+04	PHYSPROP	3.7E-04	3.6E-01	8.7E-01	9.8E-05	EPI	
Methyl Styrene (Mixed Isomers)	25013-15-4	1.5E+02	PHYSPROP	1.1E-01	2.6E-03	EPI	8.9E-01	PHYSPROP	-8.6E+01	EPI	8.9E-01	HSDB	1.7E-02	4.2E-06	WATER9	7.2E+00	EPI	3.4E+00	PHYSPROP	8.9E+01	PHYSPROP	8.9E+01	PHYSPROP	4.8E-01	1.0E+01	2.4E+01	6.6E-02	EPI	
Methyl methanesulfonate	66-27-3	1.1E+02	PHYSPROP	1.6E-04	4.0E-06	PHYSPROP	3.1E-01	PHYSPROP	2.0E+01	PHYSPROP	1.3E+00	CR89	7.9E-02	1.1E-05	WATER9	4.3E+00	EPI	6.6E-01	PHYSPROP	2.0E+05	LANGE	2.0E+05	LANGE	5.6E-04	4.4E-01	1.0E+00	1.4E-04	EPI	
Methyl tert-Butyl Ether (MTBE)	1634-04-4	8.8E+01	PHYSPROP	2.4E-02	5.9E-04	PHYSPROP	2.5E+02	PHYSPROP	-1.1E+02	PHYSPROP	7.4E-01	CR89	7.5E-02	8.6E-06	WATER9	1.2E+01	EPI	9.4E-01	PHYSPROP	5.1E+04	PHYSPROP	5.1E+04	PHYSPROP	7.6E-03	3.3E-01	7.9E-01	2.1E-03	EPI	
Methyl-1,4-benzenediamine dihydrochloride, 2-	615-45-2	2.0E+02	PHYSPROP	2.6E-16	6.4E-18	PHYSPROP	4.1E-12	PHYSPROP	2.4E+02	EPI			5.6E-02	6.6E-06	WATER9	2.0E+02	EPI	-2.1E+00	PHYSPROP	1.0E+06	PHYSPROP	1.0E+06	PHYSPROP	2.9E-05	1.3E+00	3.1E+00	5.4E-06	EPI	
Methyl-5-Nitroaniline, 2-	99-55-8	1.5E+02	PHYSPROP	3.4E-07	8.3E-09	PHYSPROP	9.8E-04	PHYSPROP	1.1E+02	PHYSPROP			6.7E-02	7.8E-06	WATER9	1.8E+02	EPI	1.9E+00	PHYSPROP	1.0E+06	PHYSPROP	1.0E+06	PHYSPROP	1.9E-05	7.5E-01	1.8E+00	3.4E-03	EPI	
Methyl-N-nitro-N-nitrosoguanidine, N-	70-25-7	1.5E+02	PHYSPROP	5.0E-11	1.2E-12	PHYSPROP	1.2E-04	PHYSPROP	1.2E+02	EPI			6.8E-02	8.0E-06	WATER9	7.2E+01	EPI	-9.2E-01	PHYSPROP	2.7E+05	PHYSPROP	2.7E+05	PHYSPROP	2.7E-04	7.0E-01	1.7E+00	5.7E-05	EPI	
Methylaniline Hydrochloride, 2-	636-21-5	1.4E+02	PHYSPROP	8.6E-05	2.1E-06	PHYSPROP	2.9E-01	PHYSPROP	2.2E+02	PHYSPROP			6.9E-02	8.1E-06	WATER9	1.2E+02	EPI	1.6E+00	PHYSPROP	8.3E+03	PHYSPROP	8.3E+03	PHYSPROP	4.8E-05	6.7E-01	1.6E+00	1.1E-05	EPI	
Methylarsonic acid	124-58-3	1.4E+02	PHYSPROP	1.6E+02	OTHER		1.6E-03	PHYSPROP	1.6E+02	PHYSPROP			7.0E-02	8.2E-06	WATER9	4.4E+01	EPI	-1.2E+00	PHYSPROP	2.6E+05	PHYSPROP	2.6E+05	PHYSPROP	1.9E-04	6.4E-01	1.5E+00	4.2E-05	EPI	
Methylbenzene,1,4-diamine monohydrochloride, 2-	74612-12-7	1.6E+02	OTHER										6.5E-02	7.6E-06	WATER9					2.7E+05	PHYSPROP	2.7E+05	PHYSPROP						
Methylbenzene-1,4-diamine sulfate, 2-	615-50-9	2.2E+02	OTHER										5.2E-02	6.1E-06	WATER9					2.9E-03	PHYSPROP	2.9E-03	PHYSPROP						
Methylcholanthrene, 3-	56-49-5	2.7E+02	PHYSPROP	2.1E-04	5.2E-06	EPI	4.3E-08	EPI	1.8E+02	PHYSPROP	1.3E+00	CR89	2.4E-02	6.1E-06	WATER9	9.6E+05	EPI	6.4E+00	PHYSPROP	2.9E-03	PHYSPROP	2.9E-03	PHYSPROP	5.7E+00	3.3E+00	1.5E+01	9.0E-01	EPI	
Methylene Chloride	75-09-2	8.5E+01	PHYSPROP	1.3E-01	3.3E-03	EPI	4.4E+02	PHYSPROP	-9.5E+01	PHYSPROP	1.3E+00	CR89	1.0E-01	1.3E-05	WATER9	2.2E+01	EPI	1.3E+00	PHYSPROP	1.3E+04	PHYSPROP	1.3E+04	PHYSPROP	1.3E-02	3.1E-01	7.5E-01	3.5E-03	EPI	
Methylene-bis(2-chloroaniline), 4,4'	103-14-1	2.7E+02	PHYSPROP	1.7E-09	4.4E-11	PHYSPROP	2.9E-07	PHYSPROP	1.1E+02	PHYSPROP			4.6E-02	5.4E-06	WATER9	5.7E+03	EPI	3.9E+00	PHYSPROP	1.4E+01	PHYSPROP	1.4E+01	PHYSPROP	1.2E-01	3.2E+00	7.9E+00	2.0E-02	EPI	
Methylene-bis(N,N-dimethyl Aniline, 4,4'	101-61-1	2.7E+02	PHYSPROP	4.4E-08	1.1E-09	PHYSPROP	1.8E-05	PHYSPROP	6.2E+01	PHYSPROP			4.7E-02	5.5E-06	WATER9	2.7E+03	EPI	4.4E+00	PHYSPROP	6.1E+00	PHYSPROP	6.1E+00	PHYSPROP	5.2E-01	2.8E+00	6.7E+00	8.4E-02	RAGSE	
Methylenedibenzideneamine, 4,4'	101-77-9	2.0E+02	PHYSPROP	2.2E-09	5.3E-11	PHYSPROP	2.0E-07	PHYSPROP	9.3E+01	PHYSPROP			5.6E-02	6.5E-06	WATER9	2.1E+03	EPI	1.6E+00	PHYSPROP	1.0E+03	PHYSPROP	1.0E+03	PHYSPROP	7.5E-03	1.4E+00	3.3E+00	1.4E-03	EPI	
Methylenediphenyl Diisocyanate	101-68-8	2.5E+02	PHYSPROP	3.7E-05	9.0E-07	PHYSPROP	5.0E-06	PHYSPROP	3.8E+01	PHYSPROP	1.2E+00	CR89	2.4E-02	6.2E-06	WATER9	2.8E+05	EPI	5.2E+00	PHYSPROP	8.3E-03	PHYSPROP	8.3E-03	PHYSPROP	1.1E+00	2.7E+00	1.0E+01	1.8E-01	EPI	
Methylstyrene, Alpha-	98-83-9	1.2E+02	PHYSPROP	1.0E-01	2.6E-03	EPI	1.9E+00	EPI	-2.3E+01	PHYSPROP	9.1E-01	CR89	6.3E-02	8.2E-06	WATER9	7.0E+02	EPI	3.5E+00	PHYSPROP	1.2E+02	PHYSPROP	1.2E+02	PHYSPROP	2.9E-01	4.8E-01	1.2E+00	7.0E-02	EPI	
Metolachlor	51218-45-2	2.8E+02	PHYSPROP	3.7E-07	9.0E-09	PHYSPROP	3.1E-05	PHYSPROP	-6.2E+01	PHYSPROP	1.1E+00	CR89	2.2E-02	5.5E-06	WATER9	4.9E+02	EPI	3.1E+00	PHYSPROP	5.3E+02	PHYSPROP	5.3E+02	PHYSPROP	2.2E-02	4.1E+00	9.8E+00	3.4E-03	EPI	
Metribuzin	21087-64-9	2.1E+02	PHYSPROP	4.8E-09	1.2E-10	EPI	4.4E-07	PHYSPROP	1.3E+02	PHYSPROP	1.3E+00	CR89	2.7E-02	7.1E-06	WATER9	5.3E+01	EPI	1.7E+00	PHYSPROP	1.1E+03	PHYSPROP	1.1E+03	PHYSPROP	7.4E-03	1.7E+00	4.0E+00	1.3E-04	EPI	
Methylsulfonyl-methyl	74223-64-6	3.8E+02	PHYSPROP	5.4E-15	1.3E-16	EPI	2.5E-12	PHYSPROP	1.6E+02	PHYSPROP			3.6E-02	4.2E-06	WATER9	9.3E+01	EPI	2.2E+00	PHYSPROP	9.5E+03	PHYSPROP	9.5E+03	PHYSPROP	2.5E-03	1.4E+01	3.4E+01	3.3E-04	EPI	
Mineral oils	8012-95-1	1.7E+02	EPI	3.3E+02	8.2E+00	EPI	1.4E-01	EPI	-9.6E+00	EPI	8.8E-01	ChemNet	3.6E-02	6.4E-06	WATER9	4.8E+03	EPI	6.1E+00	EPI	3.7E-03	EPI	3.7E-03	EPI	9.8E+00	1.9E+01	4.3E+00	2.0E+00	EPI	
Mirex	2385-85-5	5.5E+02	PHYSPROP	3.3E-02	8.1E-04	PHYSPROP	5.0E-07	PHYSPROP	4.9E+02	CR89	2.3E+00	ChemNet	2.2E-02	5.6E-06	WATER9	3.6E+05	EPI	6.9E+00	PHYSPROP	6.9E+00	PHYSPROP	6.9E+00	PHYSPROP	4.6E-01	9.5E-02	2.9E+02	5.2E-02	EPI	
Molinate	2212-67-1	1.9E+02	PHYSPROP	1.7E-04	4.1E-06	PHYSPROP	5.6E-03	PHYSPROP	7.0E+01	EPI	1.1E+00	CR89	3.2E-02	6.8E-06	WATER9	1.8E+02	EPI	3.2E+00	PHYSPROP	9.7E+02	PHYSPROP	9.7E+02	PHYSPROP	9.9E-02	1.2E+00	2.8E+00	1.9E-02	EPI	
Molybdenum	7439-98-7	9.6E+01	PHYSPROP				0.0E+00	NIOSH	2.6E+03	PHYSPROP	1.0E+01	CR89						2.0E+01	BAES					3.8E-03	3.6E-01	8.7E-01	1.0E-03	RAGSE	
Monochloramine	10599-90-3	5.1E+01	EPI						-6.6E+01	CR89														2.8E-03	2.0E-01	4.9E-01	1.0E-03	RAGSE	
Monomethylamine	100-61-8	1.1E+02	PHYSPROP	3.6E-04	8.9E-06	PHYSPROP	4.5E-01	PHYSPROP	-5.7E+01	PHYSPROP	9.9E-01	CR89	7.2E-02	9.1E-06	WATER9	8.2E+01	EPI	1.7E+00	PHYSPROP	5.6E+03	PHYSPROP	5.6E+03	PHYSPROP	2.0E-02	4.2E-01	1.0E+00	5.0E-03	EPI	
Nitrobutyl	88671-89-0	2.7E+02	PHYSPROP	1.7E-07	4.3E-09	EPI	1.6E-06	PHYSPROP	6.6E+01	PHYSPROP			4.5E-02	5.3E-06	WATER9	6.1E+03	EPI	2.9E+00	PHYSPROP	1.4E+02	PHYSPROP	1.4E+02	PHYSPROP	2.1E-02	3.6E+00	8.7E+00	3.4E-03	EPI	
N,N'-Diphenyl-1,4-benzenediamine	74-31-7	2.6E+02	PHYSPROP	8.4E-09	2.1E-10	PHYSPROP	6.4E-09	EPI	1.4E+02	PHYSPROP			4.7E-02	5.4E-06	WATER9	5.2E+04	EPI	4.0E+00	PHYSPROP	7.4E+00	PHYSPROP	7.4E+00	PHYSPROP	1.6E-01	3.0E+00	7.2E+00	2.6E-02	EPI	
Nitrobenzene	98-06-5	3.0E+02	PHYSPROP	2.7E-03	6.5E-05	EPI	2.0E-04	PHYSPROP	2.7E+01	PHYSPROP	2.0E+00	CR89	2.5E-02	6.4E-06	WATER9	1.3E+02	EPI	1.4E+00	PHYSPROP	1.5E+00	PHYSPROP	1.							

Regional Removal Management Level (RML) Chemical-specific Parameters Supporting Table May 2016

Contaminant		Molecular Weight		Volatility Parameters					Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients				Water Solubility		Tapwater Dermal Parameters							
Analyte	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m ³ /mole)	H' and HLC Ref	VP	VP Ref	MP	MP Ref	Density (g/cm ³)	Density Ref	Dia (cm ² /s)	Diw (cm ² /s)	D ₁₀ and D ₁₀₀ Ref	K _{oc} (L/kg)	K _{oc} Ref	K _{oc}	K _{oc} Ref	log K _{ow} (unitless)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/event)	t* (hr)	K _p (cm/hr)	K Ref		
Paclitaxel	76738-62-0	2.9E+02	PHYSPROP	3.4E-09	8.3E-11	EPI	7.5E-09	PHYSPROP	1.7E+02	PHYSPROP	1.2E+00	CRC89	2.2E-02	5.7E-06	WATER9	9.2E+02	EPI	3.2E+00	PHYSPROP	2.6E+01	PHYSPROP	2.6E+01	PHYSPROP	3.1E-02	4.6E+00	1.1E+01	4.7E-03	EPI		
Parquat Dichloride	1910-42-5	2.6E+02	PHYSPROP	1.3E-11	3.2E-13	PHYSPROP	7.5E-08	PHYSPROP	3.0E+02	EPI	1.2E+00	CRC89	4.7E-02	5.5E-06	WATER9	6.8E+03	EPI	4.5E+00	PHYSPROP	6.2E+05	PHYSPROP	6.2E+05	PHYSPROP	3.6E-07	2.9E+00	7.0E+00	5.8E-08	EPI		
Parathion	56-38-2	2.9E+02	PHYSPROP	1.2E-05	3.0E-07	PHYSPROP	6.7E-06	PHYSPROP	6.1E+00	PHYSPROP	1.3E+00	CRC89	2.3E-02	5.8E-06	WATER9	2.4E+03	EPI	3.8E+00	PHYSPROP	1.1E+01	PHYSPROP	1.1E+01	PHYSPROP	8.4E-02	4.5E+00	1.1E+01	1.3E-02	EPI		
Pebutate	1114-71-2	2.0E+02	PHYSPROP	9.7E-03	2.4E-04	EPI	8.9E-02	PHYSPROP	7.1E+01	EPI	9.5E-01	CRC89	2.4E-02	6.1E-06	WATER9	3.0E+02	EPI	3.8E+00	PHYSPROP	1.0E+02	PHYSPROP	1.0E+02	PHYSPROP	2.2E-01	1.4E+00	3.5E+00	4.0E-02	EPI		
Pendimethalin	40487-42-1	2.8E+02	PHYSPROP	3.5E-05	8.6E-07	EPI	1.5E-05	PHYSPROP	5.6E+01	PHYSPROP	1.2E+00	CRC89	2.3E-02	5.7E-06	WATER9	5.6E+03	EPI	5.2E+00	PHYSPROP	3.9E-01	PHYSPROP	3.9E-01	PHYSPROP	7.4E-01	4.0E+00	1.5E+01	1.2E-01	EPI		
Pentabromodiphenyl Ether	32348-91-9	5.6E+02	PHYSPROP	4.4E-03	1.1E-04	PHYSPROP	3.1E-08	EPI	8.0E+00	PHYSPROP	1.2E+00	CRC89	2.8E-02	3.2E-06	WATER9	2.2E+04	EPI	6.8E+00	PHYSPROP	2.4E-03	PHYSPROP	2.4E-03	PHYSPROP	3.4E-01	1.5E+02	3.7E+02	3.7E-02	EPI		
Pentabromodiphenyl ether, 2,2',4,4',5'-(BDE-99)	60348-60-9	5.6E+02	PHYSPROP	4.8E-05	1.2E-06	PHYSPROP	3.1E-08	EPI	8.0E+00	EPI	2.3E+00	IRIS Profile	2.2E-02	5.6E-06	WATER9	2.2E+04	EPI	7.7E+00	PHYSPROP	7.9E-05	PHYSPROP	7.9E-05	PHYSPROP	3.4E-01	1.5E+02	3.7E+02	3.7E-02	EPI		
Pentachlorobenzene	608-93-5	2.5E+02	PHYSPROP	2.9E-02	7.0E-04	PHYSPROP	1.0E-03	EPI	8.6E+01	PHYSPROP	1.8E+00	CRC89	2.9E-02	7.9E-06	WATER9	3.7E+03	EPI	5.2E+00	PHYSPROP	8.3E-01	PHYSPROP	8.3E-01	PHYSPROP	1.0E+00	2.7E+00	1.0E+01	1.7E-01	EPI		
Pentachloroethane	76-01-7	2.0E+02	PHYSPROP	7.9E-02	1.9E-03	EPI	3.5E+00	PHYSPROP	2.9E+01	PHYSPROP	1.7E+00	CRC89	3.2E-02	8.6E-06	WATER9	1.4E+02	EPI	3.2E+00	PHYSPROP	4.9E+02	PHYSPROP	4.9E+02	PHYSPROP	8.6E-02	1.4E+00	3.4E+00	1.6E-02	EPI		
Pentachloronitrobenzene	82-68-8	3.0E+02	PHYSPROP	1.8E-03	4.4E-05	EPI	5.0E-05	PHYSPROP	1.4E+02	PHYSPROP	1.7E+00	CRC89	2.6E-02	6.9E-06	WATER9	6.0E+03	EPI	4.6E+00	PHYSPROP	4.4E-01	PHYSPROP	4.4E-01	PHYSPROP	2.8E-01	4.7E+00	1.1E+01	4.2E-02	EPI		
Pentachloronitrophenol	87-86-5	2.7E+02	PHYSPROP	1.0E-06	2.5E-08	PHYSPROP	1.1E-04	PHYSPROP	1.7E+02	PHYSPROP	2.0E+00	CRC89	3.0E-02	8.0E-06	WATER9	5.9E+02	SSL	5.1E+00	PHYSPROP	1.4E+01	PHYSPROP	1.4E+01	PHYSPROP	8.0E-01	3.3E+00	1.3E+01	1.3E-01	EPI		
Pentaerythritol tetranitrate (PETN)	78-11-5	3.2E+02	PHYSPROP	5.4E-08	1.3E-09	PHYSPROP	5.5E-09	EPI	1.4E+02	PHYSPROP	1.8E+00	CRC89	2.6E-02	6.8E-06	WATER9	6.5E+02	EPI	2.4E+00	PHYSPROP	4.3E+01	PHYSPROP	4.3E+01	PHYSPROP	6.9E-03	6.2E+00	1.5E+01	1.0E-03	EPI		
Pentane, n-	109-66-0	7.2E+01	PHYSPROP	5.1E+01	1.3E+00	PHYSPROP	5.1E+02	PHYSPROP	1.3E+02	PHYSPROP	6.3E-01	CRC89	8.2E-02	8.8E-06	WATER9	7.2E+01	EPI	3.4E+00	PHYSPROP	3.8E+01	PHYSPROP	3.8E+01	PHYSPROP	3.6E-01	2.7E-01	6.4E-01	1.1E-01	EPI		
Perchlorates																														
*Ammonium Perchlorate	7790-98-9	1.2E+02	PHYSPROP								2.0E+00	CRC89										2.5E+05	PHYSPROP	4.2E-03	4.8E-01	1.1E+00	1.0E-03	RAGSE		
*Lithium Perchlorate	7791-03-9	1.1E+02	CRC89								2.4E+00	CRC89										5.9E+05	CRC89	4.0E-03	4.1E-01	1.0E+00	1.0E-03	RAGSE		
*Perchlorate and Perchlorate Salts	14797-73-0	1.2E+02	CRC89								2.4E+00	CRC89										2.5E+05	CRC89	4.2E-03	4.8E-01	1.1E+00	1.0E-03	RAGSE		
*Potassium Perchlorate	7778-74-7	1.4E+02	PHYSPROP								2.5E+00	CRC89										1.5E+04	PHYSPROP	9.1E-03	6.3E-01	1.5E+00	2.0E-03	RAGSE		
*Sodium Perchlorate	7601-89-0	1.2E+02	PHYSPROP								2.5E+00	CRC89										2.1E+06	PHYSPROP	4.3E-03	5.1E-01	1.2E+00	1.0E-03	RAGSE		
Perfluorobutane Sulfonate	375-73-5	3.0E+02	PHYSPROP	5.9E-04	1.4E-05	PHYSPROP	5.2E-02	PHYSPROP	3.7E+01	EPI	1.8E+00	LookChem	2.7E-02	7.2E-06	WATER9	1.8E+02	EPI	2.4E+00	PHYSPROP	1.1E+02	PHYSPROP	8.7E-03	5.0E+00	1.2E+01	1.3E-03	EPI				
Pernectin	52645-53-1	3.9E+02	PHYSPROP	7.7E-05	1.9E-06	EPI	2.2E-08	PHYSPROP	3.4E+01	PHYSPROP	1.2E+00	CRC89	1.9E-02	4.8E-06	WATER9	1.2E+05	EPI	6.5E+00	PHYSPROP	6.0E-03	PHYSPROP	6.0E-03	PHYSPROP	1.6E+00	1.6E+01	6.5E+01	2.1E-01	EPI		
Phenacetin	62-44-2	1.8E+02	PHYSPROP	8.7E-09	2.1E-10	EPI	6.9E-07	PHYSPROP	1.4E+02	PHYSPROP	1.2E+00	CRC89	6.0E-02	7.0E-06	WATER9	4.1E+01	EPI	1.6E+00	PHYSPROP	7.7E-02	PHYSPROP	7.7E-02	PHYSPROP	3.9E-03	1.1E+00	2.5E+00	1.7E-03	EPI		
Phenmedipham	13684-63-4	3.0E+02	PHYSPROP	3.4E-11	8.4E-13	EPI	1.0E-11	PHYSPROP	1.4E+02	PHYSPROP	1.1E+00	CRC89	4.2E-02	5.0E-06	WATER9	2.6E+03	EPI	3.6E+00	PHYSPROP	4.7E+00	PHYSPROP	4.7E+00	PHYSPROP	5.2E-02	5.1E+00	1.2E+01	7.9E-03	EPI		
Phenol	108-95-2	9.4E+01	PHYSPROP	1.4E-05	3.3E-07	PHYSPROP	3.5E-01	PHYSPROP	4.1E+01	PHYSPROP	1.1E+00	CRC89	8.9E-02	1.0E-05	WATER9	1.9E+02	EPI	1.5E+00	PHYSPROP	8.3E+04	PHYSPROP	8.3E+04	PHYSPROP	1.6E-02	3.5E-01	8.5E-01	4.3E-03	EPI		
Phenol, 2-(1-methylethoxy), methylcarbamate	114-26-1	2.1E+02	PHYSPROP	5.8E-08	1.4E-09	EPI	2.1E-05	PHYSPROP	9.0E+01	PHYSPROP	1.1E+00	CRC89	2.6E-02	6.6E-06	WATER9	6.0E+01	EPI	1.5E+00	PHYSPROP	1.9E+03	PHYSPROP	1.9E+03	PHYSPROP	6.0E-03	1.6E+00	3.7E+00	1.1E-03	EPI		
Phenothiazine	92-84-2	2.0E+02	PHYSPROP	1.1E-06	2.8E-08	PHYSPROP	8.9E-07	PHYSPROP	1.9E+02	PHYSPROP	1.3E+00	PubChem	2.9E-02	7.5E-06	WATER9	1.5E+03	EPI	4.2E+00	PHYSPROP	1.6E+00	PHYSPROP	1.6E+00	PHYSPROP	3.7E-01	1.4E+00	3.3E+01	6.8E-02	EPI		
Phenylenediamine, m-	108-45-2	1.1E+02	PHYSPROP	5.1E-08	1.3E-09	EPI	2.1E-03	EPI	6.4E+01	PHYSPROP	1.0E+00	CRC89	7.2E-02	9.2E-06	WATER9	3.4E+01	EPI	-3.3E-01	PHYSPROP	2.4E+05	PHYSPROP	9.4E-04	4.2E-01	1.0E+00	2.3E-04	EPI				
Phenylenediamine, o-	95-54-5	1.1E+02	PHYSPROP	2.9E-07	7.2E-09	EPI	2.1E-03	EPI	1.0E+02	PHYSPROP	1.2E+00	CRC89	8.4E-02	9.8E-06	WATER9	3.5E+01	EPI	1.5E-01	PHYSPROP	4.0E+04	PHYSPROP	1.9E-03	4.2E-01	1.0E+00	4.9E-04	EPI				
Phenylenediamine, p-	106-50-3	1.1E+02	PHYSPROP	2.8E-08	6.7E-10	PHYSPROP	5.0E-03	PHYSPROP	1.5E+02	PHYSPROP	1.2E+00	CRC89	8.4E-02	9.8E-06	WATER9	3.4E+01	EPI	-3.0E-01	PHYSPROP	3.7E+04	PHYSPROP	9.8E-04	4.2E-01	1.0E+00	2.5E-04	EPI				
Phenylphenol, 2-	90-43-7	1.7E+02	PHYSPROP	4.3E-05	1.1E-06	EPI	2.0E-03	EPI	5.9E+01	PHYSPROP	1.2E+00	CRC89	4.2E-02	7.8E-06	WATER9	6.7E+03	EPI	3.1E+00	PHYSPROP	7.0E+02	PHYSPROP	9.8E-02	9.4E-01	2.3E+00	2.0E-02	EPI				
Phorate	298-02-2	2.6E+02	PHYSPROP	1.8E-04	4.4E-06	EPI	6.4E-04	PHYSPROP	1.5E+01	CRC89	1.2E+00	CRC89	2.3E-02	5.9E-06	WATER9	4.6E+02	EPI	3.6E+00	PHYSPROP	5.0E+01	PHYSPROP	5.0E+01	PHYSPROP	7.8E-02	3.0E+00	7.2E+00	1.3E-02	EPI		
Phosgene	75-44-5	9.9E+01	PHYSPROP	6.8E-01	1.7E-02	PHYSPROP	1.4E-03	PHYSPROP	1.2E+02	PHYSPROP	1.4E+00	CRC89	8.9E-02	1.2E-05	WATER9	1.0E+00	EPI	-7.1E-01	PHYSPROP	6.8E+03	YAWS	5.6E-04	3.8E-01	9.0E-01	1.5E-04	EPI				
Phosmet	732-11-6	3.2E+02	PHYSPROP	3.4E-07	8.4E-09	EPI	4.9E-07	PHYSPROP	7.2E+01	PHYSPROP	1.4E+00	CRC89	4.1E-02	4.8E-06	WATER9	1.0E+01	EPI	2.8E+00	PHYSPROP	2.4E+01	PHYSPROP	2.4E+01	PHYSPROP	1.3E-02	6.3E+00	1.5E+01	1.8E-03	EPI		
Phosphates, inorganic																														
*Aluminum metaphosphate	13776-88-0	2.6E+02	CRC89								2.8E+00	CRC89												6.2E-03	3.2E+00	7.6E+00	1.0E-03	RAGSE		
*Ammonium polyphosphate	68333-79-9	7790-76-3	2.5E+02	CRC89							3.1E+00	CRC89												6.1E-03	2.8E+00	6.7E+00	1.0E-03	RAGSE		
*Calcium pyrophosphate	7790-76-3	2.5E+02	CRC89								3.1E+00	CRC89												4.4E-03	5.8E-01	1.4E+00	1.0E-03	RAGSE		
*Diammonium phosphate	7783-28-0	1.3E+02	EPI								3.1E+00	CRC89												4.5E-03	6.1E-01	1.5E+00	1.0E-03	RAGSE		
*Dicalcium phosphate	7757-93-9	1.4E+02	EPI								3.1E+00	CRC89												4.5E-03	6.1E-01	1.5E+00	1.0E-03	RAGSE		
*Dimagnesium phosphate	7782-75-4	1.7E+02	CRC89								2.1E+00	CRC89												5.1E-03	1.0E+00	2.4E+00	1.0E-03			

Regional Removal Management Level (RML) Chemical-specific Parameters Supporting Table May 2016

Contaminant	Molecular Weight		Volatility Parameters				Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients				Water Solubility		Tapwater Dermal Parameters									
	Analyle	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m ³ /mole)	H' and HLC Ref	VP	VP Ref	MP	MP Ref	Density (g/cm ³)	Density Ref	Dia (cm ² /s)	Diw (cm ² /s)	D ₁₀ and D ₁₀₀ Ref	K _{oc} (L/kg)	K _{oc} Ref	K _{oc}	K _{oc} Ref	log K _{ow} (unitless)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/event)	t [*] (hr)	K _p (cm ² /hr)	K Ref	
*Aroclor 1232	11141-16-5	1.9E+02	PHYSPROP	3.0E-02	7.4E-04	EPI	4.1E-03	PHYSPROP	3.4E+01	EPI	1.3E+00	ATSDR Profile	3.3E-02	7.5E-06	WATER9			8.4E+03	EPI	4.4E+00	PHYSPROP	1.5E+00	PHYSPROP	2.8E-01	PHYSPROP	8.9E-01	1.2E+00	4.6E+00	1.7E-01	EPI
*Aroclor 1242	53469-21-9	2.9E+02	PHYSPROP	1.4E-02	3.4E-04	PHYSPROP	8.6E-05	EPI	1.2E+02	EPI	1.4E+00	ATSDR Profile	2.4E-02	6.1E-06	WATER9			7.8E+04	EPI	6.3E+00	PHYSPROP	2.8E-01	PHYSPROP	1.5E+00	PHYSPROP	3.6E+00	4.5E+00	1.9E+01	5.5E-01	EPI
*Aroclor 1248	12672-29-6	6.2E+02	PHYSPROP	1.8E-02	4.4E-04	PHYSPROP	4.9E-04	PHYSPROP	1.2E+02	EPI	1.4E+00	HSDB	1.6E-02	3.9E-06	WATER9			7.7E+04	EPI	6.2E+00	PHYSPROP	1.0E-01	PHYSPROP	4.5E+00	3.1E+02	1.3E+03	4.8E-01	EPI		
*Aroclor 1254	11097-69-1	3.3E+02	PHYSPROP	1.2E-02	2.8E-04	PHYSPROP	7.7E-05	PHYSPROP	1.3E+02	EPI	1.5E+00	ATSDR Profile	2.4E-02	6.1E-06	WATER9			1.3E+05	EPI	6.5E+00	PHYSPROP	4.3E-02	PHYSPROP	5.2E+00	7.1E+00	3.1E+01	7.8E-01	EPI		
*Aroclor 1260	11098-82-5	4.0E+02	PHYSPROP	1.4E-02	3.4E-04	PHYSPROP	4.1E-05	PHYSPROP	1.6E+02	EPI	1.6E+00	ATSDR Profile	2.2E-02	5.6E-06	WATER9			3.5E+05	EPI	7.6E+00	PHYSPROP	1.4E-02	PHYSPROP	7.5E+00	1.7E+01	7.7E+01	9.9E-01	EPI		
*Aroclor 1460	11126-42-4	2.9E+02	PHYSPROP	5.1E-03	1.3E-04	PHYSPROP	8.5E-06	PHYSPROP	1.2E+02	EPI	1.6E+00	LookChem	2.6E-02	6.8E-06	WATER9			8.1E+04	EPI	6.3E+00	PHYSPROP	5.3E-02	PHYSPROP	3.8E+00	4.5E+00	2.0E+01	5.8E-01	EPI		
*Heptachlorobiphenyl, 2,3',4',4',5',5'- (PCB 189)	39635-11-9	4.0E+02	PHYSPROP	2.1E-03	5.1E-05	PHYSPROP	1.3E-07	PHYSPROP	1.6E+02	EPI	1.7E+00	LookChem	4.2E-02	5.7E-06	WATER9			3.5E+05	EPI	8.3E+00	PHYSPROP	7.5E-04	PHYSPROP	2.3E+01	1.7E+01	8.0E+01	3.0E+00	EPI		
*Hexachlorobiphenyl, 2,3',4',4',5',5'- (PCB 167)	52663-72-6	3.6E+02	PHYSPROP	2.8E-03	6.9E-05	PHYSPROP	5.8E-07	PHYSPROP	1.5E+02	EPI	1.6E+00	LookChem	4.4E-02	5.9E-06	WATER9			2.1E+05	EPI	7.5E+00	PHYSPROP	2.2E-03	PHYSPROP	1.0E+01	1.1E+01	5.0E+01	1.4E+00	EPI		
*Hexachlorobiphenyl, 2,3',4',4',5',5'- (PCB 157)	69782-90-7	3.6E+02	PHYSPROP	6.6E-03	1.6E-04	EPI	5.8E-07	EPI	1.5E+02	EPI	1.6E+00	I	4.4E-02	5.9E-06	WATER9			2.1E+05	EPI	7.6E+00	PHYSPROP	1.6E-03	EPI	1.2E+01	1.1E+01	5.0E+01	1.7E+00	EPI		
*Hexachlorobiphenyl, 2,3',4',4',5',5'- (PCB 156)	38380-08-4	3.6E+02	PHYSPROP	5.8E-03	1.4E-04	EPI	1.6E-06	PHYSPROP	1.5E+02	EPI	1.6E+00	LookChem	4.4E-02	5.9E-06	WATER9			2.1E+05	EPI	7.6E+00	PHYSPROP	5.3E-03	PHYSPROP	1.2E+01	1.1E+01	5.0E+01	1.7E+00	EPI		
*Hexachlorobiphenyl, 3,3',4',4',5',5'- (PCB 169)	32774-16-6	3.6E+02	PHYSPROP	2.8E-03	6.9E-05	PHYSPROP	5.8E-07	PHYSPROP	1.5E+02	EPI	1.6E+00	LookChem	4.4E-02	5.9E-06	WATER9			2.1E+05	EPI	7.4E+00	PHYSPROP	5.1E-04	PHYSPROP	9.1E+00	1.1E+01	5.0E+01	1.2E+00	EPI		
*Pentachlorobiphenyl, 2',3,4',4',5'- (PCB 123)	65510-44-3	3.3E+02	EPI	7.8E-03	1.9E-04	EPI	5.5E-06	EPI	9.8E+01	EPI	1.5E+00	LookChem	4.7E-02	6.1E-06	WATER9			1.3E+05	EPI	7.0E+00	EPI	1.6E-02	EPI	6.9E+00	7.1E+00	3.2E+01	1.0E+00	EPI		
*Pentachlorobiphenyl, 2',3,4',4',5'- (PCB 118)	31508-00-6	3.3E+02	PHYSPROP	1.2E-02	2.9E-04	EPI	9.0E-06	PHYSPROP	1.3E+02	EPI	1.5E+00	LookChem	4.7E-02	6.1E-06	WATER9			1.3E+05	EPI	7.1E+00	PHYSPROP	1.3E-02	PHYSPROP	8.6E+00	7.1E+00	3.2E+01	1.2E+00	EPI		
*Pentachlorobiphenyl, 2,3',4',4',5'- (PCB 105)	32598-14-4	3.3E+02	PHYSPROP	1.2E-02	2.8E-04	EPI	6.5E-06	PHYSPROP	1.3E+02	EPI	1.5E+00	LookChem	4.7E-02	6.1E-06	WATER9			1.3E+05	EPI	6.8E+00	PHYSPROP	3.4E-03	PHYSPROP	5.2E+00	7.1E+00	3.1E+01	7.5E-01	EPI		
*Pentachlorobiphenyl, 2,3,4',4',5'- (PCB 114)	74472-37-0	3.3E+02	PHYSPROP	3.8E-03	9.2E-05	PHYSPROP	5.5E-06	PHYSPROP	9.8E+01	PHYSPROP	1.5E+00	LookChem	4.7E-02	6.1E-06	WATER9			1.3E+05	EPI	7.0E+00	PHYSPROP	1.6E-02	PHYSPROP	6.9E+00	7.1E+00	3.2E+01	1.0E+00	EPI		
*Pentachlorobiphenyl, 3,3',4',4',5'- (PCB 126)	57465-28-8	3.3E+02	EPI	7.8E-03	1.9E-04	EPI	2.2E-06	EPI	1.3E+02	EPI	1.5E+00	LookChem	4.7E-02	6.1E-06	WATER9			1.3E+05	EPI	7.0E+00	EPI	7.3E-03	EPI	6.9E+00	7.1E+00	3.2E+01	1.0E+00	EPI		
*Polychlorinated Biphenyls (high risk)	1336-36-3	2.9E+02	PHYSPROP	1.7E-02	4.2E-04	PHYSPROP	4.9E-04	PHYSPROP	1.2E+02	EPI	1.4E+00	HSDB	2.4E-02	6.3E-06	WATER9			7.8E+04	EPI	7.1E+00	PHYSPROP	7.0E-01	PHYSPROP	3.6E+00	4.5E+00	1.9E+01	5.5E-01	EPI		
*Polychlorinated Biphenyls (low risk)	1336-36-3	2.9E+02	PHYSPROP	1.7E-02	4.2E-04	PHYSPROP	4.9E-04	PHYSPROP	1.2E+02	EPI	1.4E+00	HSDB	2.4E-02	6.3E-06	WATER9			7.8E+04	EPI	7.1E+00	PHYSPROP	7.0E-01	PHYSPROP	3.6E+00	4.5E+00	1.9E+01	5.5E-01	EPI		
*Polychlorinated Biphenyls (lowest risk)	1336-36-3	2.9E+02	PHYSPROP	1.7E-02	4.2E-04	PHYSPROP	4.9E-04	PHYSPROP	1.2E+02	EPI	1.4E+00	HSDB	2.4E-02	6.3E-06	WATER9			7.8E+04	EPI	7.1E+00	PHYSPROP	7.0E-01	PHYSPROP	3.6E+00	4.5E+00	1.9E+01	5.5E-01	EPI		
*Tetrachlorobiphenyl, 3,3',4',4'- (PCB 77)	32598-13-3	2.9E+02	PHYSPROP	3.8E-04	9.4E-06	PHYSPROP	1.6E-05	PHYSPROP	1.8E+02	CRCB9			4.9E-02	5.0E-06	WATER9			7.8E+04	EPI	6.6E+00	PHYSPROP	5.7E-04	PHYSPROP	6.0E+00	4.5E+00	2.0E+01	9.2E-01	EPI		
*Tetrachlorobiphenyl, 3,4,4',5'- (PCB 81)	70362-50-4	2.9E+02	EPI	9.1E-03	2.2E-04	EPI	8.5E-06	EPI	1.2E+02	EPI	1.4E+00	LookChem	4.9E-02	6.3E-06	WATER9			7.8E+04	EPI	6.3E+00	EPI	3.2E-02	EPI	3.8E+00	4.5E+00	2.0E+01	5.8E-01	EPI		
Polymeric Methylene Diphenyl Diisocyanate (PMDI)	9016-87-9	5.1E+02	EPI	5.4E-10	1.3E-11	EPI	5.4E-13	EPI	7.5E+02	EPI			3.0E-02	3.5E-06	WATER9			1.0E+10	EPI	1.0E+01	EPI	1.8E-06	EPI	1.6E+02	7.8E+01	3.7E+02	1.9E+01	EPI		
Polycyclic Aromatic Hydrocarbons (PAHs)																														
*Acenaphthene	83-32-9	1.5E+02	PHYSPROP	7.5E-03	1.8E-04	PHYSPROP	2.2E-03	PHYSPROP	9.3E+01	PHYSPROP	1.2E+00	CRCB9	5.1E-02	8.3E-06	WATER9			5.0E+03	EPI	3.9E+00	PHYSPROP	3.9E+00	PHYSPROP	4.1E+01	7.7E-01	1.8E+00	8.6E-02	EPI		
*Anthracene	120-12-7	1.8E+02	PHYSPROP	2.3E-03	5.6E-05	PHYSPROP	6.5E-06	EPI	2.2E+02	PHYSPROP	1.3E+00	CRCB9	3.9E-02	7.9E-06	WATER9			1.6E+04	EPI	4.5E+00	PHYSPROP	4.3E-02	PHYSPROP	7.3E-01	1.0E+00	4.1E+00	1.4E-01	EPI		
*Benzo[a]anthracene	56-55-3	2.3E+02	PHYSPROP	4.9E-04	1.2E-05	PHYSPROP	2.1E-07	PHYSPROP	8.4E+01	PHYSPROP	1.3E+00	PubChem	2.6E-02	6.7E-06	WATER9			1.8E+05	EPI	5.8E+00	PHYSPROP	9.4E-03	PHYSPROP	3.2E+00	2.0E+00	8.5E+00	5.5E-01	EPI		
*Benzo[b]fluoranthene	205-82-3	2.5E+02	PHYSPROP	8.3E-06	2.0E-07	PHYSPROP	2.6E-08	PHYSPROP	1.7E+02	PHYSPROP			4.8E-02	5.6E-06	WATER9			6.0E+05	EPI	6.1E+00	PHYSPROP	2.5E-03	PHYSPROP	4.2E+00	2.7E+00	1.2E+01	6.9E-01	EPI		
*Benzo[a]pyrene	50-32-8	2.5E+02	PHYSPROP	1.9E-05	4.6E-07	PHYSPROP	5.5E-09	EPI	1.8E+02	PHYSPROP			4.8E-02	5.6E-06	WATER9			5.9E+05	EPI	6.1E+00	PHYSPROP	1.6E-03	PHYSPROP	4.4E+00	2.7E+00	1.2E+01	7.1E-01	EPI		
*Benzo[k]fluoranthene	205-99-2	2.5E+02	PHYSPROP	2.7E-05	6.6E-07	PHYSPROP	5.0E-07	PHYSPROP	1.7E+02	PHYSPROP			4.8E-02	5.6E-06	WATER9			6.0E+05	EPI	5.8E+00	PHYSPROP	1.5E-03	PHYSPROP	2.5E+00	2.7E+00	1.1E+01	4.2E-01	EPI		
*Benzo[k]fluoranthene	207-08-9	2.5E+02	PHYSPROP	2.4E-05	5.8E-07	PHYSPROP	9.7E-10	EPI	2.2E+02	PHYSPROP			4.8E-02	5.6E-06	WATER9			5.9E+05	EPI	6.1E+00	PHYSPROP	8.0E-04	PHYSPROP	4.2E+00	2.7E+00	1.2E+01	6.9E-01	EPI		
*Chloronaphthalene, Beta-	91-58-7	1.6E+02	PHYSPROP	1.3E-02	3.2E-04	PHYSPROP	1.2E-02	EPI	6.1E+01	PHYSPROP	1.1E+00	CRCB9	4.5E-02	7.7E-06	WATER9			2.5E+03	EPI	3.9E+00	PHYSPROP	1.2E+01	PHYSPROP	3.7E-01	8.6E-01	2.1E+00	7.5E-02	EPI		
*Chrysene	218-01-9	2.3E+02	PHYSPROP	2.1E-04	5.2E-06	PHYSPROP	6.2E-09	PHYSPROP	2.6E+02	PHYSPROP	1.3E+00	CRCB9	2.6E-02	6.7E-06	WATER9			1.8E+05	EPI	5.8E+00	PHYSPROP	2.0E+03	PHYSPROP	3.5E+00	2.0E+00	8.5E+00	6.0E-01	EPI		
*Dibenz[a,h]anthracene	53-70-3	2.8E+02	PHYSPROP	5.8E-06	1.4E-07	EPI	9.6E-10	EPI	2.7E+02	PHYSPROP			4.5E-02	5.2E-06	WATER9			1.9E+06	EPI	6.8E+00	PHYSPROP	2.5E-03	PHYSPROP	6.1E+00	3.8E+00	1.7E+01	9.5E-01	EPI		
*Dibenz[a,e]pyrene	192-65-4	3.0E+02	PHYSPROP	5.8E-07	1.4E-08	PHYSPROP	7.0E-11	PHYSPROP	7.3E+02	PHYSPROP			4.2E-02	4.9E-06	WATER9			5.6E+06	EPI	7.7E+00	EPI	8.0E-05	PHYSPROP	2.8E+01	5.2E+00	2.4E+01	4.2E+00	EPI		
*Dimethylbenz[a]anthracene, 7,12-	57-97-6	2.6E+02	PHYSPROP	1.5E-04	3.8E-06	EPI	6.8E-07	PHYSPROP	1.2E+02	PHYSPROP			4.7E-02	5.5E-06	WATER9			4.9E+05	EPI	5.8E+00	PHYSPROP	6.1E-02	PHYSPROP	2.5E+00	2.9E+00	1.2E+01	4.1E-01	EPI		
*Fluoranthene	206-44-0	2.0E+02	PHYSPROP	3.6E-04	8.9E-06	PHYSPROP	9.2E-06	PHYSPROP	1.1E+02	PHYSPROP	1.3E+00	CRCB9	2.8E-02	7.2E-06	WATER9			5.5E+04	EPI	5.2E+00	PHYSPROP	2.6E-01	PHYSPROP	1.7E+00	1.4E+00	5.7E+00	3.1E-01	EPI		
*Fluorene	86-73-7	1.7E+02	PHYSPROP	0.5E-03	8.6E-05	PHYSPROP	6.0E-04	PHYSPROP	1.1E+02	PHYSPROP	1.2E+00																			

Regional Removal Management Level (RML) Chemical-specific Parameters Supporting Table May 2016

Contaminant	Molecular Weight		Volatility Parameters				Melting Point		Density		Diffusivity in Air and Water				Partition Coefficients				Water Solubility		Tapwater Dermal Parameters								
	Analyte	CAS No.	MW	MW Ref	H' (unitless)	HLC (atm-m ³ /mole)	H' and HLC Ref	VP	VP Ref	MP	MP Ref	Density (g/cm ³)	Density Ref	Dia (cm ² /s)	Diw (cm ² /s)	D _a and D _w Ref	K _{oc} (L/kg)	K _{oc} Ref	K _{ow} (L/kg)	K _{ow} Ref	log K _{ow} (unitless)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/event)	t [*] (hr)	K _p (cm ² /hr)	K Ref
Sodium Tungstate	13472-45-2	2.9E+02	CRC89						7.0E+02	CRC89	4.2E+00	CRC89											7.4E+05	CRC89	6.6E-03	4.6E+00	1.1E+01	1.0E-03	RAGSE
Sodium Tungstate Dihydrate	10213-10-2	3.3E+02	CRC89						1.0E+02	CRC89	3.3E+00	CRC89											7.4E+05	CRC89	7.0E-03	7.4E+00	1.8E+01	1.0E-03	RAGSE
Stirofos (Tetrachloroethylenes)	961-11-5	3.7E+02	PHYSPROP	7.5E-08	1.8E-09	EPI	4.2E-08	PHYSPROP	9.8E+01	PHYSPROP			3.7E-02	4.3E-06	WATER9			1.4E+03	EPI	3.5E+00	PHYSPROP		1.1E+03	PHYSPROP	2.3E-02	1.2E+01	2.8E+01	3.1E-03	EPI
Strontium Chromate	7789-06-2	2.0E+02	CRC89						7.8E+02	PHYSPROP	3.9E+00	CRC89											1.1E+03	CRC89	5.5E-03	1.5E+00	3.5E+00	1.0E-03	RAGSE
Strontium, Stable	7440-24-6	8.8E+01	PHYSPROP								2.6E+00	CRC89						3.5E+01	BAES				1.1E+03	PHYSPROP	3.6E-03	3.3E+01	7.8E+01	1.0E-03	RAGSE
Strychnine	57-24-9	3.3E+02	PHYSPROP	3.1E-12	7.6E-14	PHYSPROP	2.9E-09	PHYSPROP	2.9E+02	PHYSPROP	1.4E+00	CRC89	2.2E-02	5.6E-06	WATER9			5.4E+03	EPI	1.9E+00	PHYSPROP	1.6E+02	PHYSPROP	1.2E-03	7.8E+00	3.9E+01	4.0E-04	EPI	
Styrene	100-42-5	1.0E+02	PHYSPROP	1.1E-01	2.8E-03	PHYSPROP	6.4E+00	PHYSPROP	3.1E+01	PHYSPROP	0.9E-01	CRC89	7.1E-02	2.8E-06	WATER9			4.5E+02	EPI	3.0E+00	PHYSPROP	3.1E+02	PHYSPROP	1.5E-01	4.0E+01	9.7E+01	3.7E-02	EPI	
Styrene-Acrylonitrile (SAN) Trimer	NA	2.1E+02	OTHER								1.1E+00	PRRTV	2.6E-02	6.5E-06	WATER9							8.5E+01	PRRTV	6.6E-02	1.6E+00	3.8E+00	1.2E-02	RAGSE	
Sulfalone	126-33-0	1.2E+02	PHYSPROP	4.1E-03	4.9E-06	PHYSPROP	4.1E-03	EPI	2.8E+01	PHYSPROP	1.3E+00	CRC89	7.2E-02	9.9E-06	WATER9			1.9E+00	EPI	7.7E-01	PHYSPROP	1.0E+06	PHYSPROP	4.3E-04	5.0E-01	1.2E+00	1.0E-04	EPI	
Sulfonylbis(4-chlorobenzene), 1,1'-	80-07-9	2.9E+02	PHYSPROP	5.6E-06	1.4E-07	PHYSPROP	8.1E-07	PHYSPROP	1.5E+02	PHYSPROP	1.3E+00	CRC89	4.4E-02	5.1E-06	WATER9			2.9E+03	EPI	3.9E+00	PHYSPROP	2.4E+00	PHYSPROP	9.7E-02	4.3E+00	1.0E+01	1.5E-02	EPI	
Sulfur Trioxide	7446-11-9	8.0E+01	PHYSPROP						2.6E+02	PHYSPROP	1.9E+00	CRC89	1.2E-01	1.6E-05	WATER9								3.4E+03	PHYSPROP	3.4E-03	3.0E-01	7.1E+01	1.0E-03	RAGSE
Sulfuric Acid	7664-93-9	9.8E+01	PHYSPROP						5.9E-05	PHYSPROP	1.8E+00	CRC89											1.0E+06	PHYSPROP	3.8E-03	3.7E-01	8.9E+01	1.0E-03	RAGSE
Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methyl ester	140-57-8	3.3E+02	PHYSPROP	7.8E-06	1.9E-07	PHYSPROP	2.2E-07	PHYSPROP	3.2E+01	PHYSPROP	1.1E+00	CRC89	2.0E-02	5.0E-06	WATER9			5.6E+03	EPI	4.8E+00	PHYSPROP	5.0E+01	PHYSPROP	2.3E-01	7.9E+00	1.9E+01	3.3E-02	EPI	
TCMTB	21564-17-0	2.4E+02	PHYSPROP	2.7E-10	6.5E-12	PHYSPROP	3.1E-07	PHYSPROP	1.5E+02	EPI			4.9E-02	5.8E-06	WATER9			3.4E+03	EPI	3.3E+00	PHYSPROP	1.3E+02	PHYSPROP	6.7E-02	2.3E+00	5.5E+00	1.1E-02	EPI	
Tebuflufen	34014-18-1	2.3E+02	PHYSPROP	4.9E-09	1.2E-10	PHYSPROP	3.0E-07	PHYSPROP	1.6E+02	PHYSPROP			5.1E-02	5.9E-06	WATER9			4.2E+01	EPI	1.8E+00	PHYSPROP	2.5E+03	PHYSPROP	7.4E-03	2.0E+00	4.8E+00	1.3E-03	EPI	
Temephos	3383-96-8	4.7E+02	PHYSPROP	8.0E-08	2.0E-09	PHYSPROP	7.9E-08	PHYSPROP	3.0E+01	PHYSPROP	1.3E+00	CRC89	1.8E-02	4.5E-06	WATER9			6.0E+00	PHYSPROP	6.0E+00	PHYSPROP	2.7E-01	PHYSPROP	2.9E-01	4.3E+01	1.0E+02	3.5E-02	EPI	
Terbacil	5902-51-2	2.2E+02	PHYSPROP	4.9E-09	1.2E-10	EPI	4.7E-07	PHYSPROP	1.8E+02	PHYSPROP	1.3E+00	CRC89	2.7E-02	7.2E-06	WATER9			1.9E+00	PHYSPROP	1.9E+00	PHYSPROP	7.1E+02	PHYSPROP	9.7E-03	1.7E+00	4.1E+00	1.7E-03	EPI	
Terbufos	13071-79-9	2.9E+02	PHYSPROP	9.8E-04	2.4E-05	EPI	3.2E-04	PHYSPROP	2.9E+01	PHYSPROP	1.1E+00	CRC89	2.2E-02	5.4E-06	WATER9			1.0E+03	EPI	4.5E+00	PHYSPROP	5.1E+00	PHYSPROP	2.3E-01	4.3E+00	1.0E+01	3.6E-02	EPI	
Terbutynol	886-50-0	2.4E+02	PHYSPROP	2.8E-07	2.2E-08	EPI	1.7E-06	PHYSPROP	1.0E+02	PHYSPROP	1.1E+00	CRC89	2.4E-02	6.0E-06	WATER9			6.1E+02	EPI	3.7E+00	PHYSPROP	2.5E+01	PHYSPROP	1.3E-01	2.4E+00	5.7E+00	2.1E-02	EPI	
Tetrabromodiphenyl ether, 2,2',4,4'	5436-43-1	4.9E+02	PHYSPROP	1.2E-04	3.0E-06	PHYSPROP	7.0E-08	EPI	1.6E+02	EPI			3.1E-02	3.6E-06	WATER9			1.3E+04	EPI	6.8E+00	PHYSPROP	1.5E-03	PHYSPROP	7.9E-01	5.5E+00	2.1E+02	9.3E-02	EPI	
Tetrachlorobenzene, 1,2,4,5-	95-91-3	2.2E+02	PHYSPROP	4.1E-02	1.0E-03	PHYSPROP	5.4E-03	EPI	1.4E+02	PHYSPROP	1.9E+00	CRC89	3.2E-02	8.8E-06	WATER9			2.2E+03	EPI	4.6E+00	PHYSPROP	6.0E-01	PHYSPROP	6.4E-01	1.7E+00	6.7E+00	1.2E-01	EPI	
Tetrachloroethane, 1,1,1,2-	630-20-6	1.7E+02	PHYSPROP	1.0E-01	2.5E-03	PHYSPROP	1.2E+01	PHYSPROP	7.0E+01	PHYSPROP	1.5E+00	CRC89	4.8E-02	9.1E-06	WATER9			8.6E+01	EPI	2.9E+00	PHYSPROP	1.1E+03	PHYSPROP	7.9E-02	9.2E-01	2.2E+00	1.6E-02	EPI	
Tetrachloroethane, 1,1,2,2-	79-34-5	1.7E+02	PHYSPROP	1.5E-02	3.7E-04	PHYSPROP	4.6E+00	PHYSPROP	4.4E+01	PHYSPROP	1.6E+00	CRC89	4.9E-02	9.3E-06	WATER9			9.5E+01	EPI	2.4E+00	PHYSPROP	2.8E+03	PHYSPROP	3.5E-02	9.2E-01	2.2E+00	6.9E-03	EPI	
Tetrachloroethylene	127-18-4	1.7E+02	PHYSPROP	7.2E-01	1.8E-02	PHYSPROP	1.9E+01	PHYSPROP	3.2E+01	PHYSPROP	1.6E+00	CRC89	5.0E-02	9.5E-06	WATER9			9.5E+01	EPI	3.4E+00	PHYSPROP	2.1E+03	PHYSPROP	1.7E-01	8.9E-01	2.1E+00	3.3E-02	EPI	
Tetrachlorophenol, 2,3,4,6-	58-90-2	2.3E+02	PHYSPROP	3.6E-04	8.8E-06	EPI	6.7E-04	EPI	7.0E+01	EPI	1.2E+00	CRC89	5.0E-02	5.9E-06	WATER9			2.8E+02	SSL	4.5E+00	PHYSPROP	2.3E+01	PHYSPROP	4.2E-01	2.1E+00	5.0E+00	7.1E-02	EPI	
Tetrachlorophenol, p, alpha, alpha-	5216-25-1	2.3E+02	PHYSPROP	7.9E-03	1.9E-04	PHYSPROP	3.8E-02	PHYSPROP	4.0E+01	EPI	1.4E+00	CRC89	2.8E-02	7.3E-06	WATER9			1.6E+03	EPI	4.5E+00	PHYSPROP	4.0E+00	PHYSPROP	4.9E-01	2.0E+00	4.9E+00	8.4E-02	EPI	
Tetraethyl DiThiopyrophosphate	3689-24-5	3.2E+02	PHYSPROP	1.8E-04	4.5E-06	EPI	1.1E-04	PHYSPROP	3.2E+01	EPI	1.2E+00	CRC89	2.1E-02	5.3E-06	WATER9			2.7E+02	EPI	4.0E+00	PHYSPROP	3.0E+01	PHYSPROP	7.5E-02	6.7E+00	1.6E+01	1.1E-02	EPI	
Tetrafluoroethane, 1,1,1,2-	811-97-2	1.0E+02	PHYSPROP	2.0E+00	5.0E-02	PHYSPROP	5.0E+03	PHYSPROP	-1.0E+02	PHYSPROP	1.2E+00	CRC89	8.2E-02	1.1E-05	WATER9			8.8E+01	EPI	1.7E+00	PHYSPROP	2.0E+03	PHYSPROP	2.1E-02	3.9E-01	9.4E-01	5.5E-03	EPI	
Tetryl (Trinitrophenylmethyl)nitramine	479-45-8	2.9E+02	PHYSPROP	1.1E-07	2.7E-09	PHYSPROP	5.7E-08	PHYSPROP	1.3E+02	PHYSPROP	1.6E+00	CRC89	2.6E-02	6.7E-06	WATER9			4.6E+03	EPI	1.6E+00	PHYSPROP	7.4E+01	PHYSPROP	3.1E-03	4.3E+00	1.0E+01	4.7E-04	EPI	
Thallic Oxide	1314-32-5	4.6E+02	CRC89						8.3E+02	CRC89	1.0E+01	CRC89											9.6E+04	PHYSPROP	8.2E-03	3.8E+01	9.1E+01	1.0E-03	RAGSE
Thallium (I) Nitrate	10102-45-1	2.7E+02	PHYSPROP						2.1E+02	PHYSPROP	5.6E+00	CRC89											9.6E+04	PHYSPROP	6.3E-03	3.3E+00	7.9E+00	1.0E-03	RAGSE
Thallium (Soluble Salts)	7440-28-0	2.1E+02	PHYSPROP						3.0E+02	PHYSPROP	1.2E+01	CRC89											9.6E+04	PHYSPROP	5.5E-03	1.5E+00	3.6E+00	1.0E-03	RAGSE
Thallium Acetate	563-68-8	2.6E+02	PHYSPROP						1.5E+01	PHYSPROP	1.3E+02	CRC89	3.7E+00	1.2E-05	WATER9			1.5E+00	EPI	-1.7E-01	PHYSPROP	2.8E+04	PHYSPROP	2.5E-04	3.1E+00	7.5E+00	4.0E-05	EPI	
Thallium Carbonate	6533-73-9	4.7E+02	PHYSPROP						5.8E+00	PHYSPROP	2.7E+02	CRC89	7.1E+00	1.2E-05	WATER9			2.9E+00	EPI	-8.6E-01	PHYSPROP	2.8E+04	PHYSPROP	8.2E-06	4.4E+01	1.1E+02	9.8E-07	EPI	
Thallium Chloride	7791-32-0	2.4E+02	PHYSPROP						4.3E+02	PHYSPROP	7.0E+00	CRC89	5.2E-02	1.8E-05	WATER9			2.9E+03	PHYSPROP				2.8E+04	PHYSPROP	6.0E-03	2.3E+00	5.6E+00	1.0E-03	RAGSE
Thallium Selenite	12039-52-0	2.8E+02	EPI						3.3E+02	EPI			6.8E+00	CRC89									5.5E+04	CRC89	6.5E-03	4.1E+00	9.7E+00	1.0E-03	RAGSE
Thallium Sulfate	7446-18-6	5.0E+02	PHYSPROP						6.3E+02	PHYSPROP	6.3E+00	CRC89											5.5E+04	CRC89	8.6E-03	7.1E+01	3.7E+02	1.0E-03	RAGSE
Thiobisulfuron-methyl	79277-27-3	3.9E+02	PHYSPROP	1.7E-12	4.1E-14	PHYSPROP	1.3E-10	PHYSPROP	1.8E+02	PHYSPROP			3.6E-02	4.2E-06	WATER9			5.1E+01	EPI	1.6E+00	PHYSPROP	2.2E+03	PHYSPROP	8.6E-04	1.6E+01	3.7E+01	1.1E-04	EPI	
Thioncarnab	28249-77-6	1.2E+02	PHYSPROP	1.1E-05	2.7E-07	EPI	1.2E-05	PHYSPROP	3.3E+00	PHYSPROP	1.2E+00	CRC89	2.3E-02	5.9E-06	WATER9			1.6E+03	EPI	3.4E+00	PHYSPROP	2.8E+03	PHYSPROP	6.3E-02	2.9E+00	7.0E+00	1.0E-02	EPI	
Thiodiglycol	111-48-8	1.2E+02	PHYS																										

Regional Removal Management Level (RML) Chemical-specific Parameters Supporting Table May 2016

Contaminant		Molecular Weight		Volatility Parameters					Melting Point		Density		Diffusivity in Air and Water			Partition Coefficients				Water Solubility		Tapwater Dermal Parameters							
Analyte	CAS No.	MW	MW Ref	H ⁺ (unitless)	HLC (atm-m ³ /mole)	H ⁺ and HLC Ref	VP	VP Ref	MP	MP Ref	Density (g/cm ³)	Density Ref	D _{ia} (cm ² /s)	D _w (cm ² /s)	D _a and D _w Ref	K _{oc} (L/kg)	K _{oc} Ref	K _{ow} (L/kg)	K _{ow} Ref	log K _{ow} (unitless)	log K _{ow} Ref	S (mg/L)	S Ref	B (unitless)	T _{event} (hr/event)	t* (hr)	K _p (cm/hr)	K Ref	
Tricresyl Phosphate (TCP)	1330-78-5	3.7E+02	PHYSPROP	3.3E-05	8.1E-07	EPI	6.0E-07	EPI	-3.3E+01	PHYSPROP	1.2E+00	Yaws	1.9E-02	4.8E-06	WATER9	4.7E+04	EPI	5.1E+00	PHYSPROP	3.6E-01	PHYSPROP	3.6E-01	PHYSPROP	2.5E-01	1.2E+01	2.9E+01	3.3E-02	EPI	
Tridiphane	58138-08-2	3.2E+02	PHYSPROP	1.7E-05	4.1E-07	PHYSPROP	3.9E-04	PHYSPROP	4.3E+01	PHYSPROP	1.1E+00	PubChem	4.1E-02	4.7E-06	WATER9	3.4E+03	EPI	5.2E+00	PHYSPROP	1.1E+00	PHYSPROP	1.1E+00	PHYSPROP	4.7E-01	6.6E+00	1.6E+01	6.9E-02	EPI	
Triethylamine	121-44-8	1.0E+02	PHYSPROP	6.1E-03	1.5E-04	PHYSPROP	5.7E+01	PHYSPROP	-1.1E+02	PHYSPROP	7.3E-01	CRCB9	6.6E-02	7.9E-06	WATER9	5.1E+01	EPI	1.5E+00	PHYSPROP	6.9E+04	PHYSPROP	6.9E+04	PHYSPROP	1.5E-02	3.9E-01	9.3E-01	3.9E-03	EPI	
Triethylene Glycol	112-27-6	1.5E+02	PHYSPROP	1.3E-09	3.2E-11	PHYSPROP	1.3E-03	PHYSPROP	-7.0E+00	PHYSPROP	1.1E+00	CRCB9	5.1E-02	8.1E-06	WATER9	1.0E+01	EPI	1.8E+00	PHYSPROP	1.0E+06	PHYSPROP	1.0E+06	PHYSPROP	7.3E-05	7.3E-01	1.8E+00	1.6E-05	EPI	
Trifluoroethane, 1,1,1-	420-46-2	8.4E+01	PHYSPROP	3.1E+01	7.7E-01	PHYSPROP	9.5E-03	PHYSPROP	-1.1E+02	PHYSPROP	1.1E+00	PubChem	9.9E-02	1.2E-05	WATER9	4.4E+01	EPI	1.7E+00	PHYSPROP	7.6E+02	PHYSPROP	7.6E+02	PHYSPROP	2.7E-02	3.1E-01	7.5E-01	7.6E-03	EPI	
Trifuralin	1582-09-8	3.4E+02	PHYSPROP	4.2E-03	1.0E-04	PHYSPROP	4.6E-05	PHYSPROP	6.9E+01	PHYSPROP	1.4E+00	PubChem	2.2E-02	5.6E-06	WATER9	1.6E+04	EPI	5.3E+00	PHYSPROP	1.8E-01	PHYSPROP	1.8E-01	PHYSPROP	5.1E-01	7.9E+00	1.9E+01	7.3E-02	EPI	
Trimethyl phosphate	512-56-1	1.4E+02	PHYSPROP	2.9E-07	7.2E-09	PHYSPROP	8.5E-01	EPI	-4.6E+01	PHYSPROP	1.2E+00	CRCB9	5.8E-02	8.8E-06	WATER9	1.1E+01	EPI	6.5E-01	PHYSPROP	5.0E+05	PHYSPROP	5.0E+05	PHYSPROP	4.3E-04	6.4E-01	1.5E+00	9.5E-05	EPI	
Trimethylbenzene, 1,2,3-	526-73-8	1.2E+02	PHYSPROP	1.8E-01	4.4E-03	PHYSPROP	1.7E+00	PHYSPROP	-3.5E+01	PHYSPROP	8.9E-01	CRCB9	6.1E-02	8.0E-06	WATER9	6.3E+02	EPI	3.7E+00	PHYSPROP	7.5E+01	PHYSPROP	7.5E+01	PHYSPROP	3.8E-01	5.0E-01	1.2E+00	9.0E-02	EPI	
Trimethylbenzene, 1,2,4	95-63-6	1.2E+02	PHYSPROP	2.5E-01	6.2E-03	PHYSPROP	2.1E+00	PHYSPROP	-4.4E+01	PHYSPROP	8.8E-01	CRCB9	6.1E-02	7.9E-06	WATER9	6.1E+02	EPI	3.6E+00	PHYSPROP	5.7E+01	PHYSPROP	5.7E+01	PHYSPROP	3.6E-01	5.0E-01	1.2E+00	8.6E-02	EPI	
Trimethylbenzene, 1,3,5-	108-67-8	1.2E+02	PHYSPROP	3.6E-01	8.8E-03	PHYSPROP	2.5E+00	PHYSPROP	-4.5E+01	PHYSPROP	8.6E-01	CRCB9	6.0E-02	7.8E-06	WATER9	6.0E+02	EPI	3.4E+00	PHYSPROP	4.8E+01	PHYSPROP	4.8E+01	PHYSPROP	2.6E-01	5.0E-01	1.2E+00	6.2E-02	EPI	
Trimethylpentene, 2,4,4-	25167-70-8	1.1E+02	PHYSPROP	3.0E+01	7.5E-01	PHYSPROP	7.1E+01	PHYSPROP	-8.4E+01	EPI	7.2E-01	PubChem	6.0E-02	7.3E-06	WATER9	2.4E+02	EPI	4.1E+00	PHYSPROP	4.0E+00	PHYSPROP	4.0E+00	PHYSPROP	7.7E-01	4.5E-01	1.7E+00	1.9E-01	RAGSE	
Trinitrobenzene, 1,3,5-	99-35-4	2.1E+02	PHYSPROP	2.7E-07	6.5E-09	EPI	6.4E-06	EPI	1.2E+02	PHYSPROP	1.5E+00	CRCB9	2.9E-02	7.7E-06	WATER9	1.7E+03	EPI	1.2E+00	PHYSPROP	2.8E+02	PHYSPROP	2.8E+02	PHYSPROP	3.4E-03	1.6E+00	3.9E+00	6.1E-04	EPI	
Trinitrotoluene, 2,4,6-	118-96-7	2.3E+02	PHYSPROP	8.5E-07	2.1E-08	EPI	8.0E-06	PHYSPROP	8.0E+01	PHYSPROP	1.7E+00	CRCB9	3.0E-02	7.9E-06	WATER9	2.8E+03	EPI	1.6E+00	PHYSPROP	1.2E+02	PHYSPROP	1.2E+02	PHYSPROP	5.6E-03	2.0E+00	4.7E+00	9.6E-04	EPI	
Triphenylphosphine Oxide	791-28-6	2.8E+02	PHYSPROP	2.2E-08	5.3E-10	PHYSPROP	2.6E-09	EPI	1.6E+02	PHYSPROP	1.2E+00	CRCB9	2.3E-02	5.8E-06	WATER9	2.0E+03	EPI	2.8E+00	PHYSPROP	6.3E+01	PHYSPROP	6.3E+01	PHYSPROP	2.1E-02	3.8E+00	9.1E+00	3.3E-03	EPI	
Tris(1,3-Dichloro-2-propyl) Phosphate	13674-87-8	4.3E+02	PHYSPROP	1.1E-07	2.6E-09	PHYSPROP	7.4E-08	PHYSPROP	2.7E+01	PHYSPROP	1.4E+00	CRCB9	3.3E-02	3.9E-06	WATER9	1.1E+04	EPI	3.7E+00	PHYSPROP	7.0E+00	PHYSPROP	7.0E+00	PHYSPROP	1.3E-02	2.7E+01	6.5E+01	1.6E-03	EPI	
Tris(1-chloro-2-propyl)phosphate	13674-84-5	3.3E+02	PHYSPROP	2.4E-06	6.0E-08	PHYSPROP	2.0E-05	PHYSPROP	-4.0E+01	PHYSPROP	2.3E+00	PubChem	4.0E-02	4.7E-06	WATER9	1.6E+03	EPI	2.6E+00	PHYSPROP	1.2E+03	PHYSPROP	1.2E+03	PHYSPROP	8.4E-03	7.2E+00	1.7E+01	1.2E-03	EPI	
Tris(2,3-dibromopropyl)phosphate	126-72-7	7.0E+02	PHYSPROP	8.9E-04	2.2E-05	EPI	1.9E-04	PHYSPROP	5.5E+00	PHYSPROP	1.4E+00	PubChem	1.9E-02	4.9E-06	WATER9	9.7E+03	EPI	4.3E+00	PHYSPROP	8.0E+00	PHYSPROP	8.0E+00	PHYSPROP	1.4E-03	8.5E+02	2.0E+03	1.4E-04	EPI	
Tris(2-chloroethyl)phosphate	115-96-8	2.9E+02	PHYSPROP	1.3E-04	3.3E-06	EPI	6.1E-02	PHYSPROP	-5.5E+01	PHYSPROP	1.4E+00	CRCB9	2.4E-02	6.2E-06	WATER9	3.9E+02	EPI	1.4E+00	PHYSPROP	7.0E+03	PHYSPROP	7.0E+03	PHYSPROP	2.3E-03	4.2E+00	1.0E+01	3.6E-04	EPI	
Tris(2-ethylhexyl)phosphate	78-42-2	4.3E+02	PHYSPROP	3.2E-06	7.9E-08	EPI	8.3E-08	PHYSPROP	-7.4E+01	PHYSPROP	9.9E-01	CRCB9	1.6E-02	3.9E-06	WATER9	2.5E+06	EPI	9.5E+00	PHYSPROP	6.0E-01	PHYSPROP	6.0E-01	PHYSPROP	9.3E+01	2.9E+01	1.3E+02	1.2E+01	EPI	
Tungsten	7440-33-7	1.8E+02	PHYSPROP	0.0E+00	NIOSH		0.0E+00	NIOSH	3.4E+03	PHYSPROP	1.9E+01	CRCB9	1.5E+02	BAES		1.5E+02	BAES			5.2E-03	1.1E+00	2.7E+00	1.0E-03	RAGSE	5.9E-03	2.3E+00	5.4E+00	1.0E-03	RAGSE
Uranium (Soluble Salts)	NA	2.4E+02	CRCB9	0.0E+00	NIOSH		0.0E+00	NIOSH	1.1E+03	CRCB9	1.9E+01	CRCB9	4.5E+02	BAES		1.0E+03	SSL			5.9E-03	2.3E+00	5.4E+00	1.0E-03	RAGSE	5.9E-03	2.3E+00	5.4E+00	1.0E-03	RAGSE
Urethane	51-79-6	8.9E+01	PHYSPROP	2.6E-06	6.4E-08	EPI	2.6E-01	EPI	4.9E+01	PHYSPROP	9.9E-01	CRCB9	8.5E-02	1.0E-05	WATER9	1.2E+01	EPI	1.5E-01	PHYSPROP	4.8E+05	PHYSPROP	4.8E+05	PHYSPROP	1.4E-03	3.3E-01	8.0E-01	3.9E-04	EPI	
Vanadium Pentoxide	1314-62-1	1.8E+02	EPI	0.0E+00	NIOSH		0.0E+00	NIOSH	6.8E+02	CRCB9	3.4E+00	CRCB9	8.5E-02	1.0E-05	WATER9	1.0E+03	SSL			5.2E-03	1.1E+00	2.6E+00	1.0E-03	RAGSE	5.2E-03	2.0E-01	4.9E-01	1.0E-03	RAGSE
Vanadium and Compounds	7440-62-2	5.1E+01	EPI	0.0E+00	NIOSH		0.0E+00	NIOSH	1.9E+03	CRCB9	6.0E+00	CRCB9	8.5E-02	1.0E-05	WATER9	1.0E+03	SSL			5.2E-03	1.1E+00	2.6E+00	1.0E-03	RAGSE	5.2E-03	2.0E-01	4.9E-01	1.0E-03	RAGSE
Vernolate	1929-77-7	2.0E+02	PHYSPROP	1.3E-03	3.1E-05	EPI	1.0E-02	PHYSPROP	7.1E+01	EPI	9.5E-01	CRCB9	2.4E-02	6.1E-06	WATER9	3.0E+02	EPI	3.8E+00	PHYSPROP	9.0E+01	PHYSPROP	9.0E+01	PHYSPROP	2.2E-01	1.4E+00	3.5E+00	4.0E-02	EPI	
Vindozolin	50471-44-8	2.9E+02	PHYSPROP	1.7E-07	1.7E-08	EPI	1.2E-07	PHYSPROP	1.1E+02	PHYSPROP	1.5E+00	CRCB9	2.5E-02	6.5E-06	WATER9	2.8E+02	EPI	3.1E+00	PHYSPROP	2.6E+00	PHYSPROP	2.6E+00	PHYSPROP	2.9E-02	4.2E+00	1.0E+01	4.5E-03	EPI	
Vinyl Acetate	108-05-4	8.6E+01	PHYSPROP	2.1E-02	5.1E-04	EPI	9.0E+01	PHYSPROP	-9.3E+01	PHYSPROP	9.3E-01	CRCB9	8.5E-02	1.0E-05	WATER9	5.6E+00	EPI	7.3E-01	PHYSPROP	2.0E+04	PHYSPROP	2.0E+04	PHYSPROP	5.6E-03	3.2E-01	7.7E-01	1.6E-03	EPI	
Vinyl Bromide	593-60-2	1.1E+02	PHYSPROP	5.0E-01	1.2E-02	PHYSPROP	1.0E-03	PHYSPROP	-1.4E+02	PHYSPROP	1.5E+00	CRCB9	8.6E-02	1.2E-05	WATER9	2.2E+01	EPI	1.6E+00	PHYSPROP	7.6E+03	PHYSPROP	7.6E+03	PHYSPROP	1.7E-02	4.2E-01	1.0E+00	4.4E-03	EPI	
Vinyl Chloride	75-01-4	6.2E+01	PHYSPROP	1.1E+00	2.8E-02	PHYSPROP	3.0E+03	EPI	-1.5E+02	PHYSPROP	9.1E-01	CRCB9	1.1E-01	1.2E-05	WATER9	2.2E+01	EPI	1.4E+00	CRCB9	8.8E-03	PHYSPROP	8.8E-03	PHYSPROP	2.5E-02	2.4E-01	5.7E-01	8.4E-03	EPI	
Warfarin	81-81-2	3.1E+02	PHYSPROP	1.1E-07	2.8E-09	EPI	1.2E-07	PHYSPROP	1.6E+02	PHYSPROP	1.1E+00	CRCB9	4.2E-02	4.9E-06	WATER9	4.3E+02	EPI	2.7E+00	PHYSPROP	1.7E+01	PHYSPROP	1.7E+01	PHYSPROP	1.2E-02	5.6E+00	3.3E+01	1.8E-03	EPI	
Xylene, p-	106-42-3	1.1E+02	PHYSPROP	2.8E-01	6.9E-03	PHYSPROP	8.8E+00	PHYSPROP	1.3E+01	PHYSPROP	8.6E-01	CRCB9	6.8E-02	8.4E-06	WATER9	3.8E+02	EPI	3.2E+00	PHYSPROP	1.6E+02	PHYSPROP	1.6E+02	PHYSPROP	2.0E-01	4.1E-01	9.9E-01	4.9E-02	EPI	
Xylene, m-	108-38-3	1.1E+02	PHYSPROP	2.9E-01	7.2E-03	PHYSPROP	8.3E+00	PHYSPROP	-4.8E+01	PHYSPROP	8.6E-01	CRCB9	6.8E-02	8.4E-06	WATER9	3.8E+02	EPI	3.2E+00	PHYSPROP	1.6E+02	PHYSPROP	1.6E+02	PHYSPROP	2.1E-01	4.1E-01	9.9E-01	5.3E-02	EPI	
Xylene, o-	95-47-6	1.1E+02	PHYSPROP	2.1E-01	5.2E-03	PHYSPROP	6.6E+00	PHYSPROP	-2.5E+01	PHYSPROP	8.8E-01	CRCB9	6.9E-02	8.5E-06	WATER9	3.8E+02	EPI	3.1E+00	PHYSPROP	1.8E+02	PHYSPROP	1.8E+02	PHYSPROP	1.9E-01	4.1E-01	9.9E-01	4.7E-02	EPI	
Xylenes	1330-20-7	1.1E+02	PHYSPROP	2.7E-01	6.6E-03	PHYSPROP	8.0E+00	PHYSPROP	-2.5E+01	EPI	8.6E-01	ATSDR Profile	6.9E-02	8.5E-06	WATER9	3.8E+02	EPI	3.2E+00	PHYSPROP	1.1E+02	PHYSPROP	1.1E+02	PHYSPROP	2.0E-01	4.1E-01	9.9E-01	5.0E-02	EPI	
Zinc Phosphide	1314-84-7	2.6E+02	CRCB9	0.0E+00	NIOSH		0.0E+00	NIOSH	1.2E+03	CRCB9	4.6E+00	CRCB9	6.2E+01	SSL		6.2E+01	SSL			3.7E-03	2.9E+00	7.0E+00	6.0E-04	RAGSE	3.7E-03	2.9E+00	7.0E+00	6.0E-04	RAGSE
Zinc and Compounds	7440-66-6	6.5E+01	PHYSPROP	1.1E-07	2.7E-09	PHYSPROP	7.5E-0																						