

Key : SFo<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS p=PPRTV c=California EPA n=NCEA h=HEAST x=Withdrawn r=Route-extrapolation ca=Cancer PRG nc= Noncancer PRG ca\* (where: nc PRG < 100X ca PRG)  
 ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES							CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS	
SFo	RfDo	SFi	RfDi	V	O	CAS No.		Residential Soil (mg/kg)	"Direct Contact Exposure Pathways"			"Migration to Ground Water"		
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	skin abs.	C			Industrial Soil (mg/kg)	Ambient Air (ug/m^3)	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)		
8.7E-03	i 4.0E-03	i 8.7E-03	r 4.0E-03	r 0.1	0.1	30560-19-1	Acephate	5.6E+01	ca** 2.0E+02	ca* 7.7E-01	ca* 7.7E+00	ca*		
		7.7E-03	i 2.6E-03	i y		75-07-0	Acetaldehyde	1.1E+01	ca** 2.3E+01	ca** 8.7E-01	ca* 1.7E+00	ca		
	2.0E-02	i	2.0E-02	r	0.1	34256-82-1	Acetochlor	1.2E+03	nc 1.2E+04	nc 7.3E+01	nc 7.3E+02	nc		
	9.0E-01	i	9.0E-01	r y		67-64-1	<b>Acetone</b>	1.4E+04	nc 5.4E+04	nc 3.3E+03	nc 5.5E+03	nc	1.6E+01 8.0E-01	
	8.0E-04	h	8.0E-04	r	0.1	75-86-5	Acetone cyanohydrin	4.9E+01	nc 4.9E+02	nc 2.9E+00	nc 2.9E+01	nc		
	1.7E-02	r	1.7E-02	i y		75-05-8	Acetonitrile	4.2E+02	nc 1.8E+03	nc 6.2E+01	nc 1.0E+02	nc		
	5.0E-04	i	5.7E-06	i y		107-02-8	<b>Acrolein</b>	1.0E-01	nc 3.4E-01	nc 2.1E-02	nc 4.2E-02	nc		
4.5E+00	i 2.0E-04	i 4.5E+00	i 2.0E-04	r	0.1	79-06-1	Acrylamide	1.1E-01	ca 3.8E-01	ca 1.5E-03	ca 1.5E-02	ca		
	5.0E-01	i	2.9E-04	i	0.1	79-10-7	Acrylic acid	2.9E+04	nc 1.0E+05	max 1.0E+00	nc 1.8E+04	nc		
5.4E-01	i 1.0E-03	h 2.4E-01	i 5.7E-04	i y		107-13-1	Acrylonitrile	2.1E-01	ca* 4.9E-01	ca* 2.8E-02	ca* 3.9E-02	ca*		
1.0E+00	r	1.0E+00	c	y			<b>"CAL-Modified PRG"</b>	5.5E-02	ca 1.2E-01	ca 6.7E-03	ca 1.1E-02	ca		
8.1E-02	h 1.0E-02	i 8.0E-02	r 1.0E-02	r	0.1	15972-60-8	Alachlor	6.0E+00	ca 2.1E+01	ca 8.4E-02	ca 8.4E-01	ca		
	1.5E-01	i	1.5E-01	r	0.1	1596-84-5	Alar	9.2E+03	nc 9.2E+04	nc 5.5E+02	nc 5.5E+03	nc		
	1.0E-03	i	1.0E-03	r	0.1	116-06-3	Aldicarb	6.1E+01	nc 6.2E+02	nc 3.7E+00	nc 3.6E+01	nc		
	1.0E-03	i	1.0E-03	r	0.1	1646-88-4	Aldicarb sulfone	6.1E+01	nc 6.2E+02	nc 3.7E+00	nc 3.6E+01	nc		
1.7E+01	i 3.0E-05	i 1.7E+01	i 3.0E-05	r	0.1	309-00-2	Aldrin	2.9E-02	ca* 1.0E-01	ca 3.9E-04	ca 4.0E-03	ca	5.0E-01 2.0E-02	
	2.5E-01	i	2.5E-01	r	0.1	74223-64-6	Allyl	1.5E+04	nc 1.0E+05	max 9.1E+02	nc 9.1E+03	nc		
	5.0E-03	i	5.0E-03	r	0.1	107-18-6	Allyl alcohol	3.1E+02	nc 3.1E+03	nc 1.8E+01	nc 1.8E+02	nc		
	2.9E-04	r	2.9E-04	i	0.1	107-05-1	<b>Allyl chloride</b>	1.7E+01	nc 1.8E+02	nc 1.0E+00	nc 1.0E+01	nc		
	1.0E+00	p	1.4E-03	p		7429-90-5	Aluminum	7.6E+04	nc 1.0E+05	max 5.1E+00	nc 3.6E+04	nc		
	4.0E-04	i				20859-73-8	Aluminum phosphide	3.1E+01	nc 4.1E+02	nc	1.5E+01	nc		
	3.0E-04	i	3.0E-04	r	0.1	67485-29-4	Amdro	1.8E+01	nc 1.8E+02	nc 1.1E+00	nc 1.1E+01	nc		
	9.0E-03	i	9.0E-03	r	0.1	834-12-8	Ametryn	5.5E+02	nc 5.5E+03	nc 3.3E+01	nc 3.3E+02	nc		
	2.0E-04	n	2.0E-04	r	0.1	1321-12-6	<b>Aminodinitrotoluene</b>	1.2E+01	nc 1.2E+02	nc 7.3E-01	nc 7.3E+00	nc		
	7.0E-02	h	7.0E-02	r	0.1	591-27-5	m-Aminophenol	4.3E+03	nc 4.3E+04	nc 2.6E+02	nc 2.6E+03	nc		
	2.0E-05	h	2.0E-05	r	0.1	504-24-5	4-Aminopyridine	1.2E+00	nc 1.2E+01	nc 7.3E-02	nc 7.3E-01	nc		
	2.5E-03	i	2.5E-03	r	0.1	33089-61-1	Amitraz	1.5E+02	nc 1.5E+03	nc 9.1E+00	nc 9.1E+01	nc		
			2.9E-02	i		7664-41-7	Ammonia			1.0E+02	nc			
	2.0E-01	i			0.1	7773-06-0	Ammonium sulfamate	1.2E+04	nc 1.0E+05	max	7.3E+03	nc		
5.7E-03	i 7.0E-03	p 5.7E-03	r 2.9E-04	i	0.1	62-53-3	Aniline	8.5E+01	ca** 3.0E+02	ca* 1.0E+00	nc 1.2E+01	ca*		
	4.0E-04	i				7440-36-0	Antimony and compounds	3.1E+01	nc 4.1E+02	nc	1.5E+01	nc	5.0E+00 3.0E-01	
	1.3E-02	i	1.3E-02	r	0.1	74115-24-5	Apollo	7.9E+02	nc 8.0E+03	nc 4.7E+01	nc 4.7E+02	nc		
2.5E-02	i 5.0E-02	h 2.5E-02	i 5.0E-02	r	0.1	140-57-8	Aramite	1.9E+01	ca 6.9E+01	ca 2.7E-01	ca 2.7E+00	ca		
1.5E+00	i 3.0E-04	i 1.5E+01	i		0.03	7440-38-2	Arsenic	3.9E-01	ca* 1.6E+00	ca 4.5E-04	ca 4.5E-02	ca	2.9E+01 1.0E+00	
9.5E+00	c	1.2E+01	c		0.03		<b>"CAL-Modified PRG"</b>	6.2E-02	ca 2.5E-01	ca 5.6E-04	ca 7.1E-03	ca		
			1.4E-05	i		7784-42-1	Arsine (see arsenic for cancer endpoint)			5.2E-02	nc			
	9.0E-03	i	9.0E-03	r	0.1	76578-14-8	Assure	5.5E+02	nc 5.5E+03	nc 3.3E+01	nc 3.3E+02	nc		
	5.0E-02	i	5.0E-02	r	0.1	3337-71-1	Asulam	3.1E+03	nc 3.1E+04	nc 1.8E+02	nc 1.8E+03	nc		
2.2E-01	h 3.5E-02	i 2.2E-01	r 3.5E-02	r	0.1	1912-24-9	Atrazine	2.2E+00	ca 7.8E+00	ca 3.1E-02	ca 3.0E-01	ca		
	4.0E-04	i	4.0E-04	r	0.1	71751-41-2	Avermectin B1	2.4E+01	nc 2.5E+02	nc 1.5E+00	nc 1.5E+01	nc		
1.1E-01	i	1.1E-01	i		0.1	103-33-3	Azobenzene	4.4E+00	ca 1.6E+01	ca 6.2E-02	ca 6.1E-01	ca		
	7.0E-02	i	1.4E-04	h		7440-39-3	Barium and compounds	5.4E+03	nc 6.7E+04	nc 5.2E-01	nc 2.6E+03	nc	1.6E+03 8.2E+01	

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TOXICITY VALUES							CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS							
SFo	RfDo	SFi	RfDi	V	O	CAS No.		Residential	"Direct Contact Exposure Pathways"			"Migration to Ground Water"								
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	C	soils	Soil (mg/kg)		Soil (mg/kg)	Ambient Air	Tap Water	Soil (mg/kg)	DAF 1								
								(ug/m^3)	(ug/l)	(mg/kg)	(mg/kg)									
	4.0E-03	i	4.0E-03	r	0.1	114-26-1	Baygon	2.4E+02	nc	2.5E+03	nc	1.5E+01	nc	1.5E+02	nc					
	3.0E-02	i	3.0E-02	r	0.1	43121-43-3	Bayleton	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc					
	2.5E-02	i	2.5E-02	r	0.1	68359-37-5	Baythroid	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc					
	3.0E-01	i	3.0E-01	r	0.1	1861-40-1	Benefin	1.8E+04	nc	1.0E+05	max	1.1E+03	nc	1.1E+04	nc					
	5.0E-02	i	5.0E-02	r	0.1	17804-35-2	Benomyl	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc					
	3.0E-02	i	3.0E-02	r	0.1	25057-89-0	Bentazon	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc					
	1.0E-01	i	1.0E-01	r	0.1	100-52-7	Benzaldehyde	6.1E+03	nc	6.2E+04	nc	3.7E+02	nc	3.6E+03	nc					
5.5E-02	i	4.0E-03	i	2.7E-02	i	8.6E-03	i	y	71-43-2	<b>Benzene</b>	6.4E-01	ca*	1.4E+00	ca*	2.5E-01	ca	3.0E-02	2.0E-03		
2.3E+02	i	3.0E-03	i	2.3E+02	i	3.0E-03	r	0.1	92-87-5	<b>Benzidine</b>	2.1E-03	ca	7.5E-03	ca	2.9E-05	ca	2.9E-04	ca		
	4.0E+00	i	4.0E+00	r	0.1	65-85-0	Benzoic acid	1.0E+05	max	1.0E+05	max	1.5E+04	nc	1.5E+05	nc	4.0E+02	2.0E+01			
1.3E+01	i	1.3E+01	r	0.1	98-07-7	Benzotrichloride	3.7E-02	ca	1.3E-01	ca	5.2E-04	ca	5.2E-03	ca						
	3.0E-01	h	3.0E-01	r	0.1	100-51-6	Benzyl alcohol	1.8E+04	nc	1.0E+05	max	1.1E+03	nc	1.1E+04	nc					
1.7E-01	i	2.9E-03	r	1.7E-01	r	2.9E-03	n	y	100-44-7	Benzyl chloride	8.9E-01	ca*	2.2E+00	ca	4.0E-02	ca	6.6E-02	ca		
	2.0E-03	i	8.4E+00	i	5.7E-06	i	7440-41-7	Beryllium and compounds	1.5E+02	nc	1.9E+03	ca**	8.0E-04	ca*	7.3E+01	nc	6.3E+01	3.0E+00		
	1.0E-04	i	1.0E-04	r	0.1	141-66-2	Bidrin	6.1E+00	nc	6.2E+01	nc	3.7E-01	nc	3.6E+00	nc					
	1.5E-02	i	1.5E-02	r	0.1	82657-04-3	Biphenthrin (Talstar)	9.2E+02	nc	9.2E+03	nc	5.5E+01	nc	5.5E+02	nc					
1.1E+00	i	1.1E+00	i	y	111-44-4	<b>1,1-Biphenyl</b>	3.0E+03	nc	2.3E+04	nc	1.8E+02	nc	3.0E+02	nc						
	7.0E-02	x	4.0E-02	x	4.0E-02	r	y	108-60-1	Bis(2-chloroethyl)ether	2.2E-01	ca	5.8E-01	ca	6.1E-03	ca	1.0E-02	ca	4.0E-04	2.0E-05	
7.0E-02	x	4.0E-02	i	3.5E-02	x	4.0E-02	r	y	108-60-1	Bis(2-chloroisopropyl)ether	2.9E+00	ca	7.4E+00	ca	1.9E-01	ca	2.7E-01	ca		
2.2E+02	i	2.2E+02	i	y	542-89-1	Bis(chloromethyl)ether	1.9E-04	ca	4.3E-04	ca	3.1E-05	ca	5.2E-05	ca						
7.0E-02	x	4.0E-02	i	3.5E-02	x	4.0E-02	r	y	108-60-1	Bis(2-chloro-1-methylethyl)ether	2.9E+00	ca	7.4E+00	ca	1.9E-01	ca	2.7E-01	ca		
1.4E-02	i	2.0E-02	i	1.4E-02	r	2.0E-02	r	0.1	117-81-7	Bis(2-ethylhexyl)phthalate (DEHP)	3.5E+01	ca*	1.2E+02	ca	4.8E-01	ca	4.8E+00	ca		
	5.0E-02	i	5.0E-02	r	0.1	80-05-7	Bisphenol A	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc					
	2.00E-01	i	5.7E-03	h	7440-42-8	Boron	1.6E+04	nc	1.0E+05	max	2.1E+01	nc	7.3E+03	nc						
			2.0E-04	h	7637-07-2	Boron trifluoride					7.3E-01	nc								
7.0E-01	i	4.0E-03	i	7.0E-01	r	4.0E-03	r	0.1	15541-45-4	<b>Bromate</b>	6.9E-01	ca	2.5E+00	ca	9.6E-03	ca	9.6E-02	ca		
	2.0E-02	p	2.9E-03	p	y	108-86-1	Bromobenzene	2.8E+01	nc	9.2E+01	nc	1.0E+01	nc	2.0E+01	nc					
6.2E-02	i	2.0E-02	i	6.2E-02	r	2.0E-02	r	y	75-27-4	Bromodichloromethane	8.2E-01	ca	1.8E+00	ca	1.1E-01	ca	1.8E-01	ca	6.0E-01	3.0E-02
7.9E-03	i	2.0E-02	i	3.9E-03	i	2.0E-02	r	0.1	75-25-2	Bromoform (tribromomethane)	6.2E+01	ca*	2.2E+02	ca*	1.7E+00	ca*	8.5E+00	ca*	8.0E-01	4.0E-02
	1.4E-03	i	1.4E-03	i	y	74-83-9	Bromomethane (Methyl bromide)	3.9E+00	nc	1.3E+01	nc	5.2E+00	nc	8.7E+00	nc	2.0E-01	1.0E-02			
	5.0E-03	h	5.0E-03	r	0.1	2104-96-3	Bromophos	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc					
	2.0E-02	i	2.0E-02	r	0.1	1689-84-5	Bromoxynil	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc					
	2.0E-02	i	2.0E-02	r	0.1	1689-99-2	Bromoxynil octanoate	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc					
1.1E-01	r	5.7E-04	r	1.1E-01	i	5.7E-04	i	y	106-99-0	<b>1,3-Butadiene</b>	5.8E-02	ca*	1.2E-01	ca*	6.1E-02	ca*	1.0E-01	ca*		
6.0E-01	r	5.7E-03	r	6.0E-01	c	5.7E-03	c	y	106-99-0	<b>"CAL-Modified PRG"</b>	1.1E-02	ca	2.3E-02	ca	1.1E-02	ca	1.9E-02	ca		
	1.0E-01	i	2.6E-03	n	0.1	71-36-3	1-Butanol	6.1E+03	nc	6.1E+04	nc	9.5E+00	nc	3.6E+03	nc	1.7E+01	9.0E-01			
	5.0E-02	i	5.0E-02	r	0.1	2008-41-5	Butylate	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc					
	4.0E-02	n	4.0E-02	r	y	104-51-8	n-Butylbenzene	2.4E+02	sat	2.4E+02	sat	1.5E+02	nc	2.4E+02	nc					
	4.0E-02	n	4.0E-02	r	y	135-9-88	sec-Butylbenzene	2.2E+02	sat	2.2E+02	sat	1.5E+02	nc	2.4E+02	nc					
	4.0E-02	n	4.0E-02	r	y	98-06-6	tert-Butylbenzene	3.9E+02	sat	3.9E+02	sat	1.5E+02	nc	2.4E+02	nc					
	2.0E-01	i	2.0E-01	r	0.1	85-68-7	Butyl benzyl phthalate	1.2E+04	nc	1.0E+05	max	7.3E+02	nc	7.3E+03	nc	9.3E+02	8.1E+02			
	1.0E+00	i	1.0E+00	r	0.1	85-70-1	Butylphthalyl butylglycolate	6.1E+04	nc	1.0E+05	max	3.7E+03	nc	3.6E+04	nc					

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SFo	RfDo	SFi	RfDi	V <sub>skin</sub>	CAS No.		Residential Soil (mg/kg)	"Direct Contact Exposure Pathways"			"Migration to Ground Water"								
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	O abs.			Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)								
	5.0E-04	i	6.3E+00	i	0.001	7440-43-9	Cadmium and compounds	3.7E+01	nc	4.5E+02	nc	1.1E-03	ca	1.8E+01	nc	8.0E+00	4.0E-01		
	5.0E-01	i		5.0E-01	r	0.1	105-60-2	Caprolactam	3.1E+04	nc	1.0E+05	max	1.8E+03	nc	1.8E+04	nc			
8.6E-03	h	2.0E-03	i	8.6E-03	r	2.0E-03	r	0.1	2425-06-1	Captafol	5.7E+01	ca**	2.0E+02	ca**	7.8E-01	ca**	7.8E+00	ca**	
3.5E-03	h	1.3E-01	i	3.5E-03	r	1.3E-01	r	0.1	133-06-2	Captan	1.4E+02	ca*	4.9E+02	ca	1.9E+00	ca	1.9E+01	ca	
		1.0E-01	i		1.1E-01	r	0.1	63-25-2	Carbaryl	6.1E+03	nc	6.2E+04	nc	4.0E+02	nc	3.6E+03	nc		
2.0E-02	h		2.0E-02	r	0.1	86-74-8	Carbazole	2.4E+01	ca	8.6E+01	ca	3.4E-01	ca	3.4E+00	ca	6.0E-01	3.0E-02		
	5.0E-03	i		5.0E-03	r	0.1	1563-66-2	Carbofuran	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc			
	1.0E-01	i		2.0E-01	i	y	75-15-0	Carbon disulfide	3.6E+02	nc	7.2E+02	sat	7.3E+02	nc	1.0E+03	nc	3.2E+01	2.0E+00	
1.3E-01	i	7.0E-04	i	5.3E-02	i	7.0E-04	r	y	56-23-5	Carbon tetrachloride	2.5E-01	ca**	5.5E-01	ca*	1.3E-01	ca*	1.7E-01	ca*	
	1.0E-02	i		1.0E-02	r	0.1	55285-14-8	Carbosulfan	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc			
	1.0E-01	i		1.0E-01	r	0.1	5234-68-4	Carboxin	6.1E+03	nc	6.2E+04	nc	3.7E+02	nc	3.6E+03	nc			
	1.5E-02	i		1.5E-02	r	0.1	133-90-4	Chloramben	9.2E+02	nc	9.2E+03	nc	5.5E+01	nc	5.5E+02	nc			
4.0E-01	h		4.0E-01	r	0.1	118-75-2	Chloranil	1.2E+00	ca	4.3E+00	ca	1.7E-02	ca	1.7E-01	ca				
3.5E-01	i	5.0E-04	i	3.5E-01	i	2.0E-04	i	0.04	12789-03-6	Chlordane (technical)	1.6E+00	ca*	6.5E+00	ca*	1.9E-02	ca*	1.9E-01	ca*	
		2.0E-02	i		2.0E-02	r	0.1	90982-32-4	Chlorimuron-ethyl	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
	1.0E-01	i		5.7E-05	n		7782-50-5	Chlorine				2.1E-01	nc						
	3.0E-02	i		5.7E-05	i		10049-04-4	Chlorine dioxide				2.1E-01	nc						
	2.0E-03	h		2.0E-03	r	0.1	79-11-8	Chloroacetic acid	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc			
	8.6E-06	r		8.6E-06	i	y	532-27-4	2-Chloroacetophenone	3.3E-02	nc	1.1E-01	nc	3.1E-02	nc	5.2E-02	nc			
	4.0E-03	i		4.0E-03	r	0.1	106-47-8	4-Chloroaniline	2.4E+02	nc	2.5E+03	nc	1.5E+01	nc	1.5E+02	nc	7.0E-01	3.0E-02	
	2.0E-02	i		1.7E-02	n	y	108-90-7	Chlorobenzene	1.5E+02	nc	5.3E+02	nc	6.2E+01	nc	1.1E+02	nc	1.0E+00	7.0E-02	
2.7E-01	h	2.0E-02	i	2.7E-01	h	2.0E-02	r	0.1	510-15-6	Chlorobenzilate	1.8E+00	ca	6.4E+00	ca	2.5E-02	ca	2.5E-01	ca	
		2.0E-01	h		2.0E-01	r	0.1	74-11-3	p-Chlorobenzoic acid	1.2E+04	nc	1.0E+05	max	7.3E+02	nc	7.3E+03	nc		
		2.0E-02	h		2.0E-02	r	0.1	98-56-6	4-Chlorobenzotrifluoride	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
	2.0E-02	h		2.0E-03	h	y	126-99-8	2-Chloro-1,3-butadiene	3.6E+00	nc	1.2E+01	nc	7.3E+00	nc	1.4E+01	nc			
	4.0E-01	h		4.0E-01	r	y	109-69-3	1-Chlorobutane	4.8E+02	sat	4.8E+02	sat	1.5E+03	nc	2.4E+03	nc			
	1.4E+01	r		1.4E+01	i	y	75-68-3	1-Chloro-1,1-difluoroethane (HCFC-142b)	3.4E+02	sat	3.4E+02	sat	5.2E+04	nc	8.7E+04	nc			
	1.4E+01	r		1.4E+01	i	y	75-45-6	Chlorodifluoromethane	3.4E+02	sat	3.4E+02	sat	5.1E+04	nc	8.5E+04	nc			
2.9E-03	n	4.0E-01	n	2.9E-03	r	2.9E+00	i	y	75-00-3	Chloroethane	3.0E+00	ca	6.5E+00	ca	2.3E+00	ca	4.6E+00	ca	
		1.0E-02	i	8.1E-02	i	1.4E-02	n	y	67-66-3	Chloroform	2.2E-01	ca	4.7E-01	ca	8.3E-02	ca	1.7E-01	ca	
3.1E-02	c		1.9E-02	c		y		"CAL-Modified PRG"	9.4E-01	ca	2.0E+00	ca	3.5E-01	ca	5.3E-01	ca			
		2.6E-02	r		2.6E-02	i	y	74-87-3	Chloromethane (methyl chloride)	4.7E+01	nc	1.6E+02	nc	9.5E+01	nc	1.6E+02	nc		
5.8E-01	h		5.8E-01	r		0.1	95-69-2	4-Chloro-2-methylaniline	8.4E-01	ca	3.0E+00	ca	1.2E-02	ca	1.2E-01	ca			
4.6E-01	h		4.6E-01	r		0.1	3165-93-3	4-Chloro-2-methylaniline hydrochloride	1.1E+00	ca	3.7E+00	ca	1.5E-02	ca	1.5E-01	ca			
		8.0E-02	i		8.0E-02	r	y	91-58-7	beta-Chloronaphthalene	4.9E+03	nc	2.3E+04	nc	2.9E+02	nc	4.9E+02	nc		
9.7E-03	p	1.0E-03	p	9.7E-03	r	2.0E-05	p	y	88-73-3	o-Chloronitrobenzene	1.4E+00	nc**	4.5E+00	nc**	7.3E-02	nc**	1.5E-01	nc**	
6.7E-03	p	1.0E-03	p	6.7E-03	r	1.7E-04	p	y	100-00-5	p-Chloronitrobenzene	1.0E+01	nc**	3.7E+01	nc**	6.2E-01	nc**	1.2E+00	nc**	
		5.0E-03	i		5.0E-03	r	y	95-57-8	2-Chlorophenol	6.3E+01	nc	2.4E+02	nc	1.8E+01	nc	3.0E+01	nc	4.0E+00	2.0E-01
		2.9E-02	r		2.9E-02	h	y	75-29-6	2-Chloropropane	1.7E+02	nc	5.9E+02	nc	1.0E+02	nc	1.7E+02	nc		
1.1E-02	h	1.5E-02	i	1.1E-02	r	1.5E-02	r	0.1	1897-45-6	Chlorothalonil	4.4E+01	ca*	1.6E+02	ca*	6.1E-01	ca*	6.1E+00	ca*	
		2.0E-01	i		2.0E-02	r	y	95-49-8	o-Chlorotoluene	1.6E+02	nc	5.6E+02	nc	7.3E+01	nc	1.2E+02	nc		
		2.0E-01	i		2.0E-01	r	0.1	101-21-3	Chlorpropham	1.2E+04	nc	1.0E+05	max	7.3E+02	nc	7.3E+03	nc		

Key : SFo<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS p=PPRTV c=California EPA n=NCEA h=HEAST x=Withdrawn r=Route-extrapolation ca=Cancer PRG nc= Noncancer PRG ca\* (where: nc PRG < 100X ca PRG) ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES							CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS					
SFo 1/(mg/kg-d)	RfDo (mg/kg-d)	SFi 1/(mg/kg-d)	RfDi (mg/kg-d)	V O C	skin abs. soils	CAS No.		Residential Soil (mg/kg)	"Direct Contact Exposure Pathways" Industrial Soil (mg/kg) Ambient Air (ug/m^3) Tap Water (ug/l)			"Migration to Ground Water" DAF 20 (mg/kg) DAF 1 (mg/kg)						
	3.0E-03	i	3.0E-03	r	0.1	2921-88-2	Chlorpyrifos	1.8E+02	nc	1.8E+03	nc	1.1E+01	nc	1.1E+02	nc			
	1.0E-02	h	1.0E-02	r	0.1	5598-13-0	Chlorpyrifos-methyl	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc			
	5.0E-02	i	5.0E-02	r	0.1	64902-72-3	Chlorsulfuron	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc			
	8.0E-04	h	8.0E-04	r	0.1	60238-56-4	Chlorthiophos	4.9E+01	nc	4.9E+02	nc	2.9E+00	nc	2.9E+01	nc			
			4.2E+01	i			Total Chromium (1:6 ratio Cr VI:Cr III)+++	2.1E+02	ca	4.5E+02	ca	1.6E-04	ca			3.8E+01	2.0E+00	
	1.5E+00	i				16065-83-1	Chromium III	1.0E+05	max	1.0E+05	max		5.5E+04	nc				
	3.0E-03	i	2.9E+02	i	2.2E-06	i	18540-29-9	Chromium VI+++	3.0E+01	ca**	6.4E+01	ca	2.3E-05	ca	1.1E+02	nc	3.8E+01	2.0E+00
	2.0E-02	p	9.8E+00	p	5.7E-06	p	7440-48-4	Cobalt	9.0E+02	ca**	1.9E+03	ca*	6.9E-04	ca*	7.3E+02	nc		
			2.2E+00	i			8007-45-2	Coke Oven Emissions				3.1E-03	ca					
	4.0E-02	h					7440-50-8	Copper and compounds	3.1E+03	nc	4.1E+04	nc		1.5E+03	nc			
1.9E+00	h	1.9E+00	r		y		123-73-9	Crotonaldehyde	5.3E-03	ca	1.1E-02	ca	3.5E-03	ca	5.9E-03	ca		
	1.0E-01	i	1.1E-01	i	y		98-82-8	Cumene (isopropylbenzene)	5.7E+02	nc	2.0E+03	nc	4.0E+02	nc	6.6E+02	nc		
8.4E-01	h	2.0E-03	h	8.4E-01	r	2.0E-03	r	0.1	21725-46-2	Cyanazine	5.8E-01	ca	2.1E+00	ca	8.0E-03	ca	8.0E-02	ca
	2.0E-02	i				0.1	57-12-5	Cyanide (free)	1.2E+03	nc	1.2E+04	nc		7.3E+02	nc			
	2.0E-02	i	8.6E-04	i	y		74-90-8	Cyanide (hydrogen)	1.1E+01	nc	3.5E+01	nc	3.1E+00	nc	6.2E+00	nc		
	4.0E-02	i	4.0E-02	r	y		460-19-5	Cyanogen	1.3E+02	nc	4.3E+02	nc	1.5E+02	nc	2.4E+02	nc		
	9.0E-02	i	9.0E-02	r	y		506-68-3	Cyanogen bromide	2.9E+02	nc	9.7E+02	nc	3.3E+02	nc	5.5E+02	nc		
	5.0E-02	i	5.0E-02	r	y		506-77-4	Cyanogen chloride	1.6E+02	nc	5.4E+02	nc	1.8E+02	nc	3.0E+02	nc		
	1.7E+00	r	1.7E+00	i	y		110-82-7	Cyclohexane	1.4E+02	sat	1.4E+02	sat	6.2E+03	nc	1.0E+04	nc		
	5.0E+00	i	5.0E+00	r	0.1	108-94-1	Cyclohexanone	1.0E+05	max	1.0E+05	max	1.8E+04	nc	1.8E+05	nc			
	2.0E-01	i	2.0E-01	r	0.1	108-91-8	Cyclohexylamine	1.2E+04	nc	1.0E+05	max	7.3E+02	nc	7.3E+03	nc			
	5.0E-03	i	5.0E-03	r	0.1	68085-85-8	Cyhalothrin/Karate	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc			
	1.0E-02	i	1.0E-02	r	0.1	52315-07-8	Cypermethrin	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc			
	7.5E-03	i	7.5E-03	r	0.1	66215-27-8	Cyromazine	4.6E+02	nc	4.6E+03	nc	2.7E+01	nc	2.7E+02	nc			
	1.0E-02	i	1.0E-02	r	0.1	1861-32-1	Dacthal	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc			
	3.0E-02	i	3.0E-02	r	0.1	75-99-0	Dalapon	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc			
	2.5E-02	i	2.5E-02	r	0.1	39515-41-8	Danitol	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc			
2.4E-01	i	2.4E-01	r			0.03	72-54-8	DDD	2.4E+00	ca	1.0E+01	ca	2.8E-02	ca	2.8E-01	ca	1.6E+01	8.0E-01
3.4E-01	i	3.4E-01	r			0.03	72-55-9	DDE	1.7E+00	ca	7.0E+00	ca	2.0E-02	ca	2.0E-01	ca	5.4E+01	3.0E+00
3.4E-01	i	5.0E-04	i	3.4E-01	i	5.0E-04	r	0.03	50-29-3	DDT	1.7E+00	ca*	7.0E+00	ca*	2.0E-02	ca*	2.0E-01	ca*
	1.0E-02	i	1.0E-02	r	0.1	1163-19-5	Decabromodiphenyl ether	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc			
	4.0E-05	i	4.0E-05	r	0.1	8065-48-3	Demeton	2.4E+00	nc	2.5E+01	nc	1.5E-01	nc	1.5E+00	nc			
6.1E-02	h	6.1E-02	r			0.1	2303-16-4	Diallate	8.0E+00	ca	2.8E+01	ca	1.1E-01	ca	1.1E+00	ca		
	9.0E-04	h	9.0E-04	r	0.1	333-41-5	Diazinon	5.5E+01	nc	5.5E+02	nc	3.3E+00	nc	3.3E+01	nc			
	2.0E-03	n	2.0E-03	r	y		132-64-9	Dibenzofuran	1.5E+02	nc	1.6E+03	nc	7.3E+00	nc	1.2E+01	nc		
	1.0E-02	i	1.0E-02	r	0.1	106-37-6	1,4-Dibromobenzene	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc			
8.4E-02	i	2.0E-02	i	8.4E-02	r	2.0E-02	r	y	124-48-1	Dibromochloromethane	1.1E+00	ca	2.6E+00	ca	8.0E-02	ca	1.3E-01	ca
1.4E+00	h	5.7E-05	r	2.4E-03	x	5.7E-05	i	y	96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	4.6E-01	ca**	2.0E+00	ca**	2.1E-01	nc	4.8E-02	ca**
7.0E+00	c	7.0E+00	c		y		96-12-8	"CAL-Modified PRG"	3.0E-02	ca	7.6E-02	ca	9.6E-04	ca	1.6E-03	ca		
2.0E+00	i	9.0E-03	i	2.0E+00	i	2.6E-03	i	y	106-93-4	1,2-Dibromoethane (EDB)	3.2E-02	ca	7.3E-02	ca	3.4E-03	ca	5.6E-03	ca
	1.0E-01	i	1.0E-01	r	0.1	84-74-2	Dibutyl phthalate	6.1E+03	nc	6.2E+04	nc	3.7E+02	nc	3.6E+03	nc	2.3E+03	2.7E+02	
	3.0E-02	i	3.0E-02	r	0.1	1918-00-9	Dicamba	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc			

Key : SFo<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS p=PPRTV c=California EPA n=NCEA h=HEAST x=Withdrawn r=Route-extrapolation ca=Cancer PRG nc= Noncancer PRG ca\* (where: nc PRG < 100X ca PRG)  
 ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES						CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS				
SFo	RfDo	SFi	RfDi	V <sub>skin</sub>	CAS No.		Residential Soil (mg/kg)	"Direct Contact Exposure Pathways"			"Migration to Ground Water"					
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	O abs. C soils			Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)		DAF 20 (mg/kg)	DAF 1 (mg/kg)				
	9.0E-02	i	5.7E-02	h y	95-50-1	<b>1,2-Dichlorobenzene</b>	6.0E+02	sat	6.0E+02	sat	2.1E+02	nc	3.7E+02	nc	1.7E+01	9.0E-01
	3.0E-02	n	3.0E-02	r y	541-73-1	<b>1,3-Dichlorobenzene</b>	5.3E+02	nc	6.0E+02	sat	1.1E+02	nc	1.8E+02	nc		
2.4E-02	h 3.0E-02	n 2.2E-02	n 2.3E-01	i y	106-46-7	<b>1,4-Dichlorobenzene</b>	3.4E+00	ca	7.9E+00	ca	3.1E-01	ca	5.0E-01	ca	2.0E+00	1.0E-01
4.5E-01	i	4.5E-01	r		0.1 91-94-1	<b>3,3-Dichlorobenzidine</b>	1.1E+00	ca	3.8E+00	ca	1.5E-02	ca	1.5E-01	ca	7.0E-03	3.0E-04
	3.0E-02	n	3.0E-02	r	0.1 90-98-2	<b>4,4'-Dichlorobenzophenone</b>	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc		
9.3E+00	r	9.3E+00	h	y	764-41-0	<b>1,4-Dichloro-2-butene</b>	7.9E-03	ca	1.8E-02	ca	7.2E-04	ca	1.2E-03	ca		
	2.0E-01	i	5.7E-02	h y	75-71-8	<b>Dichlorodifluoromethane</b>	9.4E+01	nc	3.1E-02	nc	2.1E+02	nc	3.9E+02	nc		
	1.0E-01	h	1.4E-01	h y	75-34-3	<b>1,1-Dichloroethane</b>	5.1E+02	nc	1.7E+03	nc	5.2E+02	nc	8.1E+02	nc	2.3E+01	1.0E+00
5.7E-03	c	5.7E-03	c	y		<b>"CAL-Modified PRG"</b>	2.8E+00	ca	6.0E+00	ca	1.2E+00	ca	2.0E+00	ca		
9.1E-02	i 2.0E-02	n 9.1E-02	i 1.4E-03	n y	107-06-2	<b>1,2-Dichloroethane (EDC)</b>	2.8E-01	ca*	6.0E-01	ca*	7.4E-02	ca*	1.2E-01	ca*	2.0E-02	1.0E-03
	5.0E-02	i	5.7E-02	i y	75-35-4	<b>1,1-Dichloroethylene</b>	1.2E+02	nc	4.1E+02	nc	2.1E+02	nc	3.4E+02	nc	6.0E-02	3.0E-03
	1.0E-02	p	1.0E-02	r y	156-59-2	<b>1,2-Dichloroethylene (cis)</b>	4.3E+01	nc	1.5E+02	nc	3.7E+01	nc	6.1E+01	nc	4.0E-01	2.0E-02
	2.0E-02	i	2.0E-02	r y	156-60-5	<b>1,2-Dichloroethylene (trans)</b>	6.9E+01	nc	2.3E+02	nc	7.3E+01	nc	1.2E+02	nc	7.0E-01	3.0E-02
	3.0E-03	i	3.0E-03	r	0.1 120-83-2	<b>2,4-Dichlorophenol</b>	1.8E+02	nc	1.8E+03	nc	1.1E+01	nc	1.1E+02	nc	1.0E+00	5.0E-02
	8.0E-03	i	8.0E-03	r	0.1 94-82-6	<b>4-(2,4-Dichlorophenoxy)butyric Acid (2,4-DB)</b>	4.9E+02	nc	4.9E+03	nc	2.9E+01	nc	2.9E+02	nc		
	1.0E-02	i	1.0E-02	r	0.05 94-75-7	<b>2,4-Dichlorophenoxyacetic Acid (2,4-D)</b>	6.9E+02	nc	7.7E+03	nc	3.7E+01	nc	3.6E+02	nc		
6.8E-02	h 1.1E-03	r 6.8E-02	r 1.1E-03	i y	78-87-5	<b>1,2-Dichloropropane</b>	3.4E-01	ca*	7.4E-01	ca*	9.9E-02	ca*	1.6E-01	ca*	3.0E-02	1.0E-03
	2.0E-02	p	2.0E-02	r y	142-28-9	<b>1,3-Dichloropropane</b>	1.0E+02	nc	3.6E+02	nc	7.3E+01	nc	1.2E+02	nc		
1.0E-01	i 3.0E-02	i 1.4E-02	i 5.7E-03	i y	542-75-6	<b>1,3-Dichloropropene</b>	7.8E-01	ca	1.8E+00	ca	4.8E-01	ca	4.0E-01	ca	4.0E-03	2.0E-04
	3.0E-03	i	3.0E-03	r	0.1 616-23-9	<b>2,3-Dichloropropanol</b>	1.8E+02	nc	1.8E+03	nc	1.1E+01	nc	1.1E+02	nc		
2.9E-01	i 5.0E-04	i 2.9E-01	r 1.4E-04	i	0.1 62-73-7	<b>Dichlorvos</b>	1.7E+00	ca*	5.9E+00	ca*	2.3E-02	ca*	2.3E-01	ca*		
4.4E-01	x	4.4E-01	r		0.1 115-32-2	<b>Dicofol</b>	1.1E+00	ca	3.9E+00	ca	1.5E-02	ca	1.5E-01	ca		
	3.0E-02	h	5.7E-05	x y	77-73-6	<b>Dicyclopentadiene</b>	5.4E-01	nc	1.8E+00	nc	2.1E-01	nc	4.2E-01	nc		
1.6E+01	i 5.0E-05	i 1.6E+01	i 5.0E-05	r	0.1 60-57-1	<b>Dieldrin</b>	3.0E-02	ca	1.1E-01	ca	4.2E-04	ca	4.2E-03	ca	4.0E-03	2.0E-04
	1.0E-02	p	5.7E-03	p	0.1 112-34-5	<b>Diethylene glycol, monobutyl ether</b>	6.1E+02	nc	6.2E+03	nc	2.1E+01	nc	3.6E+02	nc		
	6.0E-02	p	8.6E-04	p	0.1 111-90-0	<b>Diethylene glycol, monoethyl ether</b>	3.7E+03	nc	3.7E+04	nc	3.1E+00	nc	2.2E+03	nc		
	4.0E-04	p	4.0E-04	r	0.1 617-84-5	<b>Diethylformamide</b>	2.4E+01	nc	2.5E+02	nc	1.5E+00	nc	1.5E+01	nc		
1.2E-03	i 6.0E-01	i 1.2E-03	r 6.0E-01	r	0.1 103-23-1	<b>Di(2-ethylhexyl)adipate</b>	4.1E+02	ca	1.4E+03	ca	5.6E+00	ca	5.6E+01	ca		
	8.0E-01	i	8.0E-01	r	0.1 84-66-2	<b>Diethyl phthalate</b>	4.9E+04	nc	1.0E+05	max	2.9E+03	nc	2.9E+04	nc		
4.7E+03	h	4.7E+03	r		0.1 56-53-1	<b>Diethylstilbestrol</b>	1.0E-04	ca	3.7E-04	ca	1.4E-06	ca	1.4E-05	ca		
	8.0E-02	i	8.0E-02	r	0.1 43222-48-6	<b>Difenzozquat (Avenge)</b>	4.9E+03	nc	4.9E+04	nc	2.9E+02	nc	2.9E+03	nc		
	2.0E-02	i	2.0E-02	r	0.1 35367-38-5	<b>Diflubenzuron</b>	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
	1.1E+01	r	1.1E+01	i y	75-37-6	<b>1,1-Difluoroethane</b>					4.2E+04	nc	6.9E+04	nc		
	2.0E-02	n	2.0E-02	r	0.1 28553-12-0	<b>Diisononyl phthalate</b>	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
			1.1E-01	p	108-20-3	<b>Diisopropyl ether</b>					4.0E+02	nc				
	8.0E-02	i	8.0E-02	r	0.1 1445-75-6	<b>Diisopropyl methylphosphonate</b>	4.9E+03	nc	4.9E+04	nc	2.9E+02	nc	2.9E+03	nc		
	2.0E-02	i	2.0E-02	r	0.1 55290-64-7	<b>Dimethipin</b>	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
	2.0E-04	i	2.0E-04	r	0.1 60-51-5	<b>Dimethoate</b>	1.2E+01	nc	1.2E+02	nc	7.3E-01	nc	7.3E+00	nc		
1.4E-02	h	1.4E-02	r		0.1 119-90-4	<b>3,3'-Dimethoxybenzidine</b>	3.5E+01	ca	1.2E+02	ca	4.8E-01	ca	4.8E+00	ca		
	5.7E-06	r	5.7E-06	x y	124-40-3	<b>Dimethylamine</b>	6.7E-02	nc	2.5E-01	nc	2.1E-02	nc	3.5E-02	nc		
	2.0E-03	i	2.0E-03	r	0.1 121-69-7	<b>N-N-Dimethylaniline</b>	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc		
7.5E-01	h	7.5E-01	r		0.1 95-68-1	<b>2,4-Dimethylaniline</b>	6.5E-01	ca	2.3E+00	ca	9.0E-03	ca	9.0E-02	ca		

Key : SFo<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS p=PPRTV c=California EPA n=NCEA h=HEAST x=Withdrawn r=Route-extrapolation ca=Cancer PRG nc= Noncancer PRG ca\* (where: nc PRG < 100X ca PRG)  
 ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES						CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS				
SFo	RfDo	SFi	RfDi	V <sub>skin</sub>	CAS No.		Residential Soil (mg/kg)	"Direct Contact Exposure Pathways"			"Migration to Ground Water"					
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	O abs. C soils			Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)		DAF 20 (mg/kg)	DAF 1 (mg/kg)				
5.8E-01	h	5.8E-01	r	0.1	21436-96-4	2,4-Dimethylaniline hydrochloride	8.4E-01	ca	3.0E+00	ca	1.2E-02	ca	1.2E-01	ca		
2.3E+00	p	2.3E+00	r	0.1	119-93-7	<b>3,3'-Dimethylbenzidine</b>	2.1E-01	ca	7.5E-01	ca	2.9E-03	ca	2.9E-02	ca		
1.0E-01	h	8.6E-03	i	0.1	68-12-2	N,N-Dimethylformamide	6.1E+03	nc	6.2E+04	nc	3.1E+01	nc	3.6E+03	nc		
1.0E-03	n	1.0E-03	r	0.1	122-09-8	Dimethylphenethylamine	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc		
2.0E-02	i	2.0E-02	r	0.1	105-67-9	2,4-Dimethylphenol	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
6.0E-04	i	6.0E-04	r	0.1	576-26-1	2,6-Dimethylphenol	3.7E+01	nc	3.7E+02	nc	2.2E+00	nc	2.2E+01	nc		
1.0E-03	i	1.0E-03	r	0.1	95-65-8	3,4-Dimethylphenol	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc		
1.0E+01	h	1.0E+01	r	0.1	131-11-3	Dimethyl phthalate	1.0E+05	max	1.0E+05	max	3.7E+04	nc	3.6E+05	nc		
1.0E-01	i	1.0E-01	r	0.1	120-61-6	Dimethyl terephthalate	6.1E+03	nc	6.2E+04	nc	3.7E+02	nc	3.6E+03	nc		
1.0E-04	p	1.0E-04	r	0.1	534-52-1	<b>4,6-Dinitro-o-cresol</b>	6.1E+00	nc	6.2E+01	nc	3.7E-01	nc	3.6E+00	nc		
2.0E-03	i	2.0E-03	r	0.1	131-89-5	4,6-Dinitro-o-cyclohexyl phenol	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc		
1.0E-04	p	1.0E-04	r	0.1	528-29-0	1,2-Dinitrobenzene	6.1E+00	nc	6.2E+01	nc	3.7E-01	nc	3.6E+00	nc		
1.0E-04	i	1.0E-04	r	0.1	99-65-0	1,3-Dinitrobenzene	6.1E+00	nc	6.2E+01	nc	3.7E-01	nc	3.6E+00	nc		
1.0E-04	p	1.0E-04	r	0.1	100-25-4	1,4-Dinitrobenzene	6.1E+00	nc	6.2E+01	nc	3.7E-01	nc	3.6E+00	nc		
2.0E-03	i	2.0E-03	r	0.1	51-28-5	2,4-Dinitrophenol	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc		
6.8E-01	i	6.8E-01	r	0.1	25321-14-6	Dinitrotoluene mixture	7.2E-01	ca	2.5E+00	ca	9.9E-03	ca	9.9E-02	ca		
2.0E-03	i	2.0E-03	r	0.1	121-14-2	2,4-Dinitrotoluene (also see Dinitrotoluene mixture)	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc		
1.0E-03	h	1.0E-03	r	0.1	606-20-2	2,6-Dinitrotoluene (also see Dinitrotoluene mixture)	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc		
1.0E-03	i	1.0E-03	r	0.1	88-85-7	Dinoseb	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc		
4.0E-02	p	4.0E-02	r	0.1	117-84-0	di-n-Octyl phthalate	2.4E+03	nc	2.5E+04	nc	1.5E+02	nc	1.5E+03	nc		
1.1E-02	i	1.1E-02	r	0.1	123-91-1	1,4-Dioxane	4.4E+01	ca	1.6E+02	ca	6.1E-01	ca	6.1E+00	ca		
1.5E+05	h	1.5E+05	h	0.03	1746-01-6	Dioxin (2,3,7,8-TCDD)+++	3.9E-06	ca	1.6E-05	ca	4.5E-08	ca	4.5E-07	ca		
3.0E-02	i	3.0E-02	r	0.1	957-51-7	Diphenamid	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc		
2.5E-02	i	2.5E-02	r	0.1	122-39-4	Diphenylamine	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc		
3.0E-04	p	3.0E-04	r	0.1	74-31-7	N,N-Diphenyl-1,4 benzenediamine (DPPD)	1.8E+01	nc	1.8E+02	nc	1.1E+00	nc	1.1E+01	nc		
8.0E-01	i	8.0E-01	i	0.1	122-66-7	1,2-Diphenylhydrazine	6.1E-01	ca	2.2E+00	ca	8.4E-03	ca	8.4E-02	ca		
3.0E-03	p	3.0E-03	r	0.1	127-63-9	Diphenyl sulfone	1.8E+02	nc	1.8E+03	nc	1.1E+01	nc	1.1E+02	nc		
2.2E-03	i	2.2E-03	r	0.1	85-00-7	Diquat	1.3E+02	nc	1.4E+03	nc	8.0E+00	nc	8.0E+01	nc		
8.6E+00	h	8.6E+00	r	0.1	1937-37-7	Direct black 38	5.7E-02	ca	2.0E-01	ca	7.8E-04	ca	7.8E-03	ca		
8.1E+00	h	8.1E+00	r	0.1	2602-46-2	Direct blue 6	6.0E-02	ca	2.1E-01	ca	8.3E-04	ca	8.3E-03	ca		
9.3E+00	h	9.3E+00	r	0.1	16071-86-6	Direct brown 95	5.2E-02	ca	1.9E-01	ca	7.2E-04	ca	7.2E-03	ca		
4.0E-05	i	4.0E-05	r	0.1	298-04-4	Disulfoton	2.4E+00	nc	2.5E+01	nc	1.5E-01	nc	1.5E+00	nc		
1.0E-02	i	1.0E-02	r	0.1	505-29-3	1,4-Dithiane	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc		
2.0E-03	i	2.0E-03	r	0.1	330-54-1	Diuron	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc		
4.0E-03	i	4.0E-03	r	0.1	2439-10-3	Dodine	2.4E+02	nc	2.5E+03	nc	1.5E+01	nc	1.5E+02	nc		
1.0E-01	n				7429-91-6	Dysprosium	7.8E+03	nc	1.0E+05	max			3.6E+03	nc		
6.0E-03	i	6.0E-03	r	0.1	115-29-7	Endosulfan	3.7E+02	nc	3.7E+03	nc	2.2E+01	nc	2.2E+02	nc		
2.0E-02	i	2.0E-02	r	0.1	145-73-3	Endothall	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
3.0E-04	i	3.0E-04	r	0.1	72-20-8	Endrin	1.8E+01	nc	1.8E+02	nc	1.1E+00	nc	1.1E+01	nc		
9.9E-03	i	2.0E-03	h	4.2E-03	h	2.9E-04	i	y	106-89-8	Epichlorohydrin	7.6E+00	nc	2.6E+01	nc	1.0E+00	5.0E-02

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TOXICITY VALUES							CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS				
SFo	RfDo	SFi	RfDi	V	skin	CAS No.		Residential	"Direct Contact Exposure Pathways"			"Migration to Ground Water"					
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	O	abs.			Soil (mg/kg)	Soil (mg/kg)	Ambient Air	Tap Water	DAF 20	DAF 1				
				C	soils				(ug/m^3)	(ug/l)	(mg/kg)	(mg/kg)					
8.00E-02	r	8.00E-02	c	y			<b>"CAL-Modified PRG"</b>	1.3E+00	nc	2.9E+00	nc	8.4E-02	nc	1.4E-01	nc		
	5.7E-03	r	5.7E-03	i	0.1	106-88-7	1,2-Epoxybutane	3.5E+02	nc	3.5E+03	nc	2.1E+01	nc	2.1E+02	nc		
	2.5E-02	i	2.5E-02	r	0.1	759-94-4	EPTC (S-Ethyl dipropylthiocarbamate)	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc		
	5.0E-03	i	5.0E-03	r	0.1	16672-87-0	Ethephon (2-chloroethyl phosphonic acid)	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc		
	5.0E-04	i	5.0E-04	r	0.1	563-12-2	Ethion	3.1E+01	nc	3.1E+02	nc	1.8E+00	nc	1.8E+01	nc		
	4.0E-01	h	5.7E-02	i	0.1	110-80-5	2-Ethoxyethanol	2.4E+04	nc	1.0E+05	max	2.1E+02	nc	1.5E+04	nc		
	3.0E-01	h	3.0E-01	r	0.1	111-15-9	2-Ethoxyethanol acetate	1.8E+04	nc	1.0E+05	max	1.1E+03	nc	1.1E+04	nc		
	9.0E-01	i	9.0E-01	r	y	141-78-6	Ethyl acetate	1.9E+04	nc	3.7E+04	sat	3.3E+03	nc	5.5E+03	nc		
4.8E-02	h	4.8E-02	r	y		140-88-5	Ethyl acrylate	2.1E-01	ca	4.5E-01	ca	1.4E-01	ca	2.3E-01	ca		
	1.0E-01	i	2.9E-01	i	y	100-41-4	<b>Ethylbenzene</b>	4.0E+02	sat	4.0E+02	sat	1.1E+03	nc	1.3E+03	nc	1.3E+01 7.0E-01	
2.9E-03	n	4.0E-01	n	2.9E-03	r	2.9E+00	Ethyl chloride	3.0E+00	ca	6.5E+00	ca	2.3E+00	ca	4.6E+00	ca		
	3.0E-01	h	3.0E-01	r	0.1	109-78-4	Ethylene cyanohydrin	1.8E+04	nc	1.0E+05	max	1.1E+03	nc	1.1E+04	nc		
	9.0E-02	p	9.0E-02	r	0.1	107-15-3	<b>Ethylene diamine</b>	5.5E+03	nc	5.5E+04	nc	3.3E+02	nc	3.3E+03	nc		
	2.0E+00	i	2.0E+00	r	0.1	107-21-1	Ethylene glycol	1.0E+05	max	1.0E+05	max	7.3E+03	nc	7.3E+04	nc		
	5.0E-01	i	3.7E+00	i	0.1	111-76-2	Ethylene glycol, monobutyl ether	3.1E+04	nc	1.0E+05	max	1.4E+04	nc	1.8E+04	nc		
1.0E+00	h	3.5E-01	h	y		75-21-8	Ethylene oxide	1.4E-01	ca	3.4E-01	ca	1.9E-02	ca	2.4E-02	ca		
1.1E-01	h	8.0E-05	i	1.1E-01	r	8.0E-05	Ethylene thiourea (ETU)	4.4E+00	ca**	1.6E+01	ca**	6.1E-02	ca**	6.1E-01	ca**		
	2.0E-01	i	2.0E-01	r	y	60-29-7	Ethyl ether	1.8E+03	sat	1.8E+03	sat	7.3E+02	nc	1.2E+03	nc		
	9.0E-02	h	9.0E-02	r	y	97-63-2	Ethyl methacrylate	1.4E+02	sat	1.4E+02	sat	3.3E+02	nc	5.5E+02	nc		
	1.0E-05	i	1.0E-05	r	0.1	2104-64-5	Ethyl p-nitrophenyl phenylphosphorothioate	6.1E-01	nc	6.2E+00	nc	3.7E-02	nc	3.6E-01	nc		
	3.0E+00	i	3.0E+00	r	0.1	84-72-0	Ethylphthalyl ethyl glycolate	1.0E+05	max	1.0E+05	max	1.1E+04	nc	1.1E+05	nc		
	8.0E-03	i	8.0E-03	r	0.1	101200-48-0	Express	4.9E+02	nc	4.9E+03	nc	2.9E+01	nc	2.9E+02	nc		
	2.5E-04	i	2.5E-04	r	0.1	22224-92-6	Fenamiphos	1.5E+01	nc	1.5E+02	nc	9.1E-01	nc	9.1E+00	nc		
	1.3E-02	i	1.3E-02	r	0.1	2164-17-2	Fluometuron	7.9E+02	nc	8.0E+03	nc	4.7E+01	nc	4.7E+02	nc		
	6.0E-02	i			0.1	16984-48-8	Fluorine (soluble fluoride)	3.7E+03	nc	3.7E+04	nc			2.2E+03	nc		
	8.0E-02	i	8.0E-02	r	0.1	59756-60-4	Fluoridone	4.9E+03	nc	4.9E+04	nc	2.9E+02	nc	2.9E+03	nc		
	2.0E-02	i	2.0E-02	r	0.1	56425-91-3	Flurprimidol	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
	6.0E-02	i	6.0E-02	r	0.1	66332-96-5	Flutolanil	3.7E+03	nc	3.7E+04	nc	2.2E+02	nc	2.2E+03	nc		
	1.0E-02	i	1.0E-02	r	0.1	69409-94-5	Fluvalinate	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc		
3.5E-03	i	1.0E-01	i	3.5E-03	r	1.0E-01	Folpet	1.4E+02	ca*	4.9E+02	ca	1.9E+00	ca	1.9E+01	ca		
1.9E-01	i	1.9E-01	r		0.1	72178-02-0	Fomesafen	2.6E+00	ca	9.1E+00	ca	3.5E-02	ca	3.5E-01	ca		
	2.0E-03	i	2.0E-03	r	0.1	944-22-9	Fonofos	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc		
	1.5E-01	i	4.6E-02	i		0.1	50-00-0	Formaldehyde	9.2E+03	nc	1.0E+05	nc	1.5E-01	ca	5.5E+03	nc	
	2.0E+00	h	8.6E-04	p	0.1	64-18-6	<b>Formic Acid</b>	1.0E+05	max	1.0E+05	max	3.1E+00	nc	7.3E+04	nc		
	3.0E+00	i	3.0E+00	r	0.1	39148-24-8	Fosetyl-al	1.0E+05	max	1.0E+05	max	1.1E+04	nc	1.1E+05	nc		
	3.0E+01	i	8.6E+00	h	y	76-13-1	Freon 113	5.6E+03	sat	5.6E+03	sat	3.1E+04	nc	5.9E+04	nc		
	1.0E-03	i	1.0E-03	r	y	110-00-9	Furan	2.5E+00	nc	8.5E+00	nc	3.7E+00	nc	6.1E+00	nc		
3.8E+00	h	3.8E+00	r		0.1	67-45-8	Furazolidone	1.3E-01	ca	4.5E-01	ca	1.8E-03	ca	1.8E-02	ca		
	3.0E-03	i	1.4E-02	h	0.1	98-01-1	Furfural	1.8E+02	nc	1.8E+03	nc	5.2E+01	nc	1.1E+02	nc		
5.0E+01	h	5.0E+01	r		0.1	531-82-8	Furium	9.7E-03	ca	3.4E-02	ca	1.3E-04	ca	1.3E-03	ca		
3.0E-02	i	3.0E-02	r		0.1	60568-05-0	Furmecyclox	1.6E+01	ca	5.7E+01	ca	2.2E-01	ca	2.2E+00	ca		
	4.0E-04	i	4.0E-04	r	0.1	77182-82-2	Glufosinate-ammonium	2.4E+01	nc	2.5E+02	nc	1.5E+00	nc	1.5E+01	nc		

Key : SFo<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS p=PPRTV c=California EPA n=NCEA h=HEAST x=Withdrawn r=Route-extrapolation ca=Cancer PRG nc= Noncancer PRG ca\* (where: nc PRG < 100X ca PRG) ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES							CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS							
SFo	RfDo	SFi	RfDi	V	skin	CAS No.		Residential Soil (mg/kg)	"Direct Contact Exposure Pathways"			"Migration to Ground Water"								
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	O	abs.			Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)								
	4.0E-04	i	2.9E-04	h	0.1	765-34-4	Glycidaldehyde	2.4E+01	nc	2.5E+02	nc	1.0E+00	nc	1.5E+01	nc					
	1.0E-01	i	1.0E-01	r	0.1	1071-83-6	Glyphosate	6.1E+03	nc	6.2E+04	nc	3.7E+02	nc	3.6E+03	nc					
	5.0E-05	i	5.0E-05	r	0.1	69806-40-2	Haloxfop-methyl	3.1E+00	nc	3.1E+01	nc	1.8E-01	nc	1.8E+00	nc					
	1.3E-02	i	1.3E-02	r	0.1	79277-27-3	Harmony	7.9E+02	nc	8.0E+03	nc	4.7E+01	nc	4.7E+02	nc					
4.5E+00	i	5.0E-04	i	4.6E+00	i	5.0E-04	r	0.1	76-44-8	Heptachlor	1.1E-01	ca	3.8E-01	ca	1.5E-03	ca	1.5E-02	ca	2.3E+01	1.0E+00
9.1E+00	i	1.3E-05	i	9.1E+00	i	1.3E-05	r	0.1	1024-57-3	Heptachlor epoxide	5.3E-02	ca*	1.9E-01	ca*	7.4E-04	ca*	7.4E-03	ca*	7.0E-01	3.0E-02
	2.0E-03	i	2.0E-03	r	0.1	87-82-1	Hexabromobenzene	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc					
1.6E+00	i	8.0E-04	i	1.6E+00	i	8.0E-04	r	0.1	118-74-1	Hexachlorobenzene	3.0E-01	ca	1.1E+00	ca	4.2E-03	ca	4.2E-02	ca	2.0E+00	1.0E-01
7.8E-02	i	3.0E-04	n	7.8E-02	i	3.0E-04	r	0.1	87-68-3	Hexachlorobutadiene	6.2E+00	ca**	2.2E+01	ca**	8.6E-02	ca*	8.6E-01	ca*	2.0E+00	1.0E-01
6.3E+00	i	5.0E-04	n	6.3E+00	i	5.0E-04	r	0.04	319-84-6	HCH (alpha)	9.0E-02	ca	3.6E-01	ca	1.1E-03	ca	1.1E-02	ca	5.0E-04	3.0E-05
1.8E+00	i	2.0E-04	n	1.8E+00	i	2.0E-04	r	0.04	319-85-7	HCH (beta)	3.2E-01	ca	1.3E+00	ca	3.7E-03	ca	3.7E-02	ca	3.0E-03	1.0E-04
1.3E+00	h	3.0E-04	i	1.3E+00	r	3.0E-04	r	0.04	58-89-9	HCH (gamma) Lindane	4.4E-01	ca*	1.7E+00	ca	5.2E-03	ca	5.2E-02	ca	9.0E-03	5.0E-04
1.8E+00	i	1.8E+00	i			0.04	608-73-1	HCH-technical	3.2E-01	ca	1.3E+00	ca	3.8E-03	ca	3.7E-02	ca	3.0E-03	1.0E-04		
	6.0E-03	i	5.7E-05	i	0.1	77-47-4	Hexachlorocyclopentadiene	3.7E+02	nc	3.7E+03	nc	2.1E-01	nc	2.2E+02	nc	4.0E+02	2.0E+01			
1.4E-02	i	1.0E-03	i	1.4E-02	i	1.0E-03	r	0.1	67-72-1	Hexachloroethane	3.5E+01	ca**	1.2E+02	ca**	4.8E-01	ca**	4.8E+00	ca**	5.0E-01	2.0E-02
	3.0E-04	i	3.0E-04	r	0.1	70-30-4	Hexachlorophene	1.8E+01	nc	1.8E+02	nc	1.1E+00	nc	1.1E+01	nc					
1.1E-01	i	3.0E-03	i	1.1E-01	r	3.0E-03	r	0.1	121-82-4	Hexahydro-1,3,5-trinitro-1,3,5-triazine	4.4E+00	ca*	1.6E+01	ca	6.1E-02	ca	6.1E-01	ca		
	2.9E-06	r	2.9E-06	i	0.1	822-06-0	1,6-Hexamethylene diisocyanate	1.7E-01	nc	1.8E+00	nc	1.0E-02	nc	1.0E-01	nc					
	1.1E+01	p	5.7E-02	i	y	110-54-3	<b>n-Hexane</b>	1.1E+02	sat	1.1E+02	sat	2.1E+02	nc	4.2E+02	nc					
	3.3E-02	i	3.3E-02	r	0.1	51235-04-2	Hexazinone	2.0E+03	nc	2.0E+04	nc	1.2E+02	nc	1.2E+03	nc					
	5.0E-02	i	5.0E-02	r	0.1	2691-41-0	HMX	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc					
3.0E+00	i	1.7E+01	i			0.1	302-01-2	Hydrazine, hydrazine sulfate	1.6E-01	ca	5.7E-01	ca	3.9E-04	ca	2.2E-02	ca				
3.0E+00	n	1.7E+01	n			0.1	60-34-4	Hydrazine, monomethyl	1.6E-01	ca	5.7E-01	ca	4.0E-04	ca	2.2E-02	ca				
3.0E+00	n	1.7E+01	n			0.1	57-14-7	Hydrazine, dimethyl	1.6E-01	ca	5.7E-01	ca	4.0E-04	ca	2.2E-02	ca				
			5.7E-03	i		7647-01-0	Hydrogen chloride					2.1E+01	nc							
	2.0E-02	i	8.6E-04	i	y	74-90-8	Hydrogen cyanide	1.1E+01	nc	3.5E+01	nc	3.1E+00	nc	6.2E+00	nc					
	3.0E-03	i	2.9E-04	i		7783-06-4	Hydrogen sulfide					1.0E+00	nc	1.1E+02	nc					
5.6E-02	p	4.0E-02	p	5.6E-02	r	4.0E-02	r	0.1	123-31-9	<b>p-Hydroquinone</b>	8.7E+00	ca	3.1E+01	ca	1.2E-01	ca	1.2E+00	ca		
	1.3E-02	i	1.3E-02	r	0.1	35554-44-0	Imazalil	7.9E+02	nc	8.0E+03	nc	4.7E+01	nc	4.7E+02	nc					
	2.5E-01	i	2.5E-01	r	0.1	81335-37-7	Imazaquin	1.5E+04	nc	1.0E+05	max	9.1E+02	nc	9.1E+03	nc					
	4.0E-02	i	4.0E-02	r	0.1	36734-19-7	Iprodione	2.4E+03	nc	2.5E+04	nc	1.5E+02	nc	1.5E+03	nc					
	3.0E-01	n				7439-89-6	Iron	2.3E+04	nc	1.0E+05	max			1.1E+04	nc					
	3.0E-01	i	3.0E-01	r	y	78-83-1	Isobutanol	1.3E+04	nc	4.0E+04	sat	1.1E+03	nc	1.8E+03	nc					
9.5E-04	i	2.0E-01	i	9.5E-04	r	2.0E-01	r	0.1	78-59-1	Isophorone	5.1E+02	ca*	5.1E+02	ca*	7.1E+00	ca	7.1E+01	ca	5.0E-01	3.0E-02
	1.5E-02	i	1.5E-02	r	0.1	33820-53-0	Isopropalin	9.2E+02	nc	9.2E+03	nc	5.5E+01	nc	5.5E+02	nc					
	1.0E-01	i	1.1E-01	r	0.1	1832-54-8	Isopropyl methyl phosphonic acid	6.1E+03	nc	6.2E+04	nc	4.0E+02	nc	3.6E+03	nc					
	5.0E-02	i	5.0E-02	r	0.1	82558-50-7	Isoxaben	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc					
8.0E+00	p	2.0E-04	p	8.0E+00	r	2.0E-04	r	0.1	143-50-0	Kepone	6.1E-02	ca	2.2E-01	ca	8.4E-04	ca	8.4E-03	ca		
	2.0E-03	i	2.0E-03	r	0.1	77501-63-4	Lactofen	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc					
						7439-92-1	<b>Lead+++</b>	4.0E+02	nc	8.0E+02	nc									
							"CAL-Modified PRG"+++	1.5E+02	nc											
	1.0E-07	i			0.1	78-00-2	Lead (tetraethyl)	6.1E-03	nc	6.2E-02	nc			3.6E-03	nc					

[www.epa.gov/superfund/programs/lead/ieubk.htm](http://www.epa.gov/superfund/programs/lead/ieubk.htm)  
[www.dtsc.ca.gov/ScienceTechnology/ledspread.html](http://www.dtsc.ca.gov/ScienceTechnology/ledspread.html)

Key : SFo<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS p=PPRTV c=California EPA n=NCEA h=HEAST x=Withdrawn r=Route-extrapolation ca=Cancer PRG nc= Noncancer PRG ca\* (where: nc PRG < 100X ca PRG)  
 ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES							CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS							
SFo 1/(mg/kg-d)	RfDo (mg/kg-d)	SFi 1/(mg/kg-d)	RfDi (mg/kg-d)	V O C	skin abs. soils	CAS No.		Residential Soil (mg/kg)	"Direct Contact Exposure Pathways"			"Migration to Ground Water"								
							Industrial Soil (mg/kg)	Ambient Air (ug/m^3)	Tap Water (ug/l)		DAF 20 (mg/kg)	DAF 1 (mg/kg)								
	2.0E-03	i	2.0E-03	r	0.1	330-55-2	Linuron	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc					
	2.0E-02	x				7439-93-2	Lithium	1.6E+03	nc	2.0E+04	nc			7.3E+02	nc					
	2.0E-01	i	2.0E-01	r	0.1	83055-99-6	Londax	1.2E+04	nc	1.0E+05	max	7.3E+02	nc	7.3E+03	nc					
	2.0E-02	i	2.0E-02	r	0.1	121-75-5	Malathion	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc					
	1.0E-01	i	1.0E-01	r	0.1	108-31-6	Maleic anhydride	6.1E+03	nc	6.2E+04	nc	3.7E+02	nc	3.6E+03	nc					
	5.0E-01	i	5.0E-01	r	y	123-33-1	Maleic hydrazide	1.7E+03	nc	2.4E+03	sat	1.8E+03	nc	3.0E+03	nc					
	1.0E-04	p	1.0E-04	r	0.1	109-77-3	Malononitrile	6.1E+00	nc	6.2E+01	nc	3.7E-01	nc	3.6E+00	nc					
	3.0E-02	h	3.0E-02	r	0.1	8018-01-7	Mancozeb	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc					
6.0E-02	o	5.0E-03	i	6.0E-02	r	5.0E-03	r	0.1	12427-38-2	Maneb	8.1E+00	ca*	2.9E+01	ca	1.1E-01	ca	1.1E+00	ca		
	2.4E-02	i	1.4E-05	i		7439-96-5	Manganese and compounds+++	1.8E+03	nc	1.9E+04	nc	5.1E-02	nc	8.8E+02	nc					
	9.0E-05	h	9.0E-05	r	0.1	950-10-7	Mephosfolan	5.5E+00	nc	5.5E+01	nc	3.3E-01	nc	3.3E+00	nc					
	3.0E-02	i	3.0E-02	r	0.1	24307-26-4	Mepiquat chloride	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc					
2.9E-02	n	1.0E-01	n	2.9E-02	r	1.0E-01	r	0.1	149-30-4	2-Mercaptobenzothiazole	1.7E+01	ca	5.9E+01	ca	2.3E-01	ca	2.3E+00	ca		
	3.0E-04	i				7487-94-7	Mercury and compounds	2.3E+01	nc	3.1E+02	nc			1.1E+01	nc					
			8.6E-05	i		7439-97-6	Mercury (elemental)					3.1E-01	nc							
	1.0E-04	i			0.1	22967-92-6	Mercury (methyl)	6.1E+00	nc	6.2E+01	nc			3.6E+00	nc					
	3.0E-05	i	3.0E-05	r	0.1	150-50-5	Merphos	1.8E+00	nc	1.8E+01	nc	1.1E-01	nc	1.1E+00	nc					
	3.0E-05	i	3.0E-05	r	0.1	78-48-8	Merphos oxide	1.8E+00	nc	1.8E+01	nc	1.1E-01	nc	1.1E+00	nc					
	6.0E-02	i	6.0E-02	r	0.1	57837-19-1	Metalaxyl	3.7E+03	nc	3.7E+04	nc	2.2E+02	nc	2.2E+03	nc					
	1.0E-04	i	2.0E-04	h	y	126-98-7	Methacrylonitrile	2.1E+00	nc	8.4E+00	nc	7.3E-01	nc	1.0E+00	nc					
	5.0E-05	i	5.0E-05	r	0.1	10265-92-6	Methamidophos	3.1E+00	nc	3.1E+01	nc	1.8E-01	nc	1.8E+00	nc					
	5.0E-01	i	5.0E-01	r	0.1	67-56-1	Methanol	3.1E+04	nc	1.0E+05	max	1.8E+03	nc	1.8E+04	nc					
	1.0E-03	i	1.0E-03	r	0.1	950-37-8	Methidathion	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc					
	2.5E-02	i	2.5E-02	r	y	16752-77-5	Methomyl	4.4E+01	nc	1.5E+02	nc	9.1E+01	nc	1.5E+02	nc					
	5.0E-03	i	5.0E-03	r	0.1	72-43-5	Methoxychlor	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc	1.6E+02	8.0E+00			
	1.0E-03	h	5.7E-03	i	0.1	109-86-4	2-Methoxyethanol	6.1E+01	nc	6.2E+02	nc	2.1E+01	nc	3.6E+01	nc					
	2.0E-03	h	2.0E-03	r	0.1	110-49-6	2-Methoxyethanol acetate	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc					
4.6E-02	h		4.6E-02	r		99-59-2	2-Methoxy-5-nitroaniline	1.1E+01	ca	3.7E+01	ca	1.5E-01	ca	1.5E+00	ca					
	1.0E+00	h	1.0E+00	r	y	79-20-9	Methyl acetate	2.2E+04	nc	9.2E+04	nc	3.7E+03	nc	6.1E+03	nc					
	3.0E-02	h	3.0E-02	r	y	96-33-3	Methyl acrylate	7.0E+01	nc	2.3E+02	nc	1.1E+02	nc	1.8E+02	nc					
2.4E-01	h		2.4E-01	r		95-53-4	2-Methylaniline (o-toluidine)	2.0E+00	ca	7.2E+00	ca	2.8E-02	ca	2.8E-01	ca					
1.8E-01	h		1.8E-01	r		636-21-5	2-Methylaniline hydrochloride	2.7E+00	ca	9.6E+00	ca	3.7E-02	ca	3.7E-01	ca					
	5.0E-04	i	5.0E-04	r	0.1	94-74-6	2-Methyl-4-chlorophenoxyacetic acid	3.1E+01	nc	3.1E+02	nc	1.8E+00	nc	1.8E+01	nc					
	1.0E-02	i	1.0E-02	r	0.1	94-81-5	4-(2-Methyl-4-chlorophenoxy) butyric acid	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc					
	1.0E-03	i	1.0E-03	r	0.1	93-65-2	2-(2-Methyl-4-chlorophenoxy) propionic acid	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc					
	1.0E-03	i	1.0E-03	r	0.1	16484-77-8	2-(2-Methyl-1,4-chlorophenoxy) propionic acid	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc					
	8.6E-01	r	8.6E-01	h	y	108-87-2	Methylcyclohexane	2.6E+03	nc	8.7E+03	nc	3.1E+03	nc	5.2E+03	nc					
2.5E-01	h		2.5E-01	r		101-77-9	4,4'-Methylenebisbenzeneamine	1.9E+00	ca	6.9E+00	ca	2.7E-02	ca	2.7E-01	ca					
1.3E-01	h	7.0E-04	h	1.3E-01	h	7.0E-04	r	0.1	101-14-4	4,4'-Methylene bis(2-chloroaniline)	3.7E+00	ca*	1.3E+01	ca*	5.2E-02	ca*	5.2E-01	ca*		
4.6E-02	i		4.6E-02	r		101-61-1	4,4'-Methylene bis(N,N'-dimethyl)aniline	1.1E+01	ca	3.7E+01	ca	1.5E-01	ca	1.5E+00	ca					
	1.0E-02	h	1.0E-02	r	y	74-95-3	Methylene bromide	6.7E+01	nc	2.3E+02	nc	3.7E+01	nc	6.1E+01	nc					
7.5E-03	i	6.0E-02	i	1.6E-03	i	8.6E-01	h	y	75-09-2	Methylene chloride	9.1E+00	ca	2.1E+01	ca	4.1E+00	ca	4.3E+00	ca	2.0E-02	1.0E-03

Key : SFo<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS p=PPRTV c=California EPA n=NCEA h=HEAST x=Withdrawn r=Route-extrapolation ca=Cancer PRG nc= Noncancer PRG ca\* (where: nc PRG < 100X ca PRG)  
 ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES						CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS			
SFo	RfDo	SFi	RfDi	V <sub>skin</sub>	CAS No.		Residential Soil (mg/kg)	Industrial Soil (mg/kg)	"Direct Contact Exposure Pathways"		"Migration to Ground Water"				
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	O abs.				Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)				
	1.7E-04	r	1.7E-04	i	0.1	101-68-8	4,4'-Methylene diphenyl diisocyanate	1.0E+01	nc	1.0E+02	nc	6.2E-01	nc	6.2E+00	nc
	6.0E-01	i	1.4E+00	i	y	78-93-3	<b>Methyl ethyl ketone (2-Butanone)</b>	2.2E+04	nc	1.1E+05	nc	5.1E+03	nc	7.0E+03	nc
	8.0E-02	h	8.6E-01	i	y	108-10-1	<b>Methyl isobutyl ketone</b>	5.3E+03	nc	4.7E+04	nc	3.1E+03	nc	2.0E+03	nc
	5.7E-04	r	5.7E-04	n	0.1	74-93-1	Methyl Mercaptan	3.5E+01	nc	3.5E+02	nc	2.1E+00	nc	2.1E+01	nc
	1.4E+00	i	2.0E-01	i	y	80-62-6	Methyl methacrylate	2.2E+03	nc	2.7E+03	sat	7.3E+02	nc	1.4E+03	nc
3.3E-02	h	3.3E-02	r		0.1	99-55-8	2-Methyl-5-nitroaniline	1.5E+01	ca	5.2E+01	ca	2.0E-01	ca	2.0E+00	ca
	2.5E-04	i	2.5E-04	r	0.1	298-00-0	Methyl parathion	1.5E+01	nc	1.5E+02	nc	9.1E-01	nc	9.1E+00	nc
	5.0E-02	i	5.0E-02	r	0.1	95-48-7	2-Methylphenol	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc
	5.0E-02	i	5.0E-02	r	0.1	108-39-4	3-Methylphenol	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc
	5.0E-03	h	5.0E-03	r	0.1	106-44-5	4-Methylphenol	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc
	2.0E-02	p	2.0E-02	r	0.1	993-13-5	Methyl phosphonic acid	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc
	6.0E-03	h	1.1E-02	h	y	25013-15-4	Methyl styrene (mixture)	1.3E+02	nc	5.4E+02	nc	4.2E+01	nc	6.0E+01	nc
	7.0E-02	h	7.0E-02	r	y	98-83-9	Methyl styrene (alpha)	6.8E+02	sat	6.8E+02	sat	2.6E+02	nc	4.3E+02	nc
1.8E-03	c	8.6E-01	r	9.1E-04	c	8.6E-01	<b>Methyl tertbutyl ether (MTBE)</b>	3.2E+01	ca	7.0E+01	ca	7.4E+00	ca	1.1E+01	ca
	1.5E-01	i	1.5E-01	r	0.1	51218-45-2	Metolacloer (Dual)	9.2E+03	nc	9.2E+04	nc	5.5E+02	nc	5.5E+03	nc
	2.5E-02	i	2.5E-02	r	0.1	21087-64-9	Metribuzin	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc
1.8E+00	x	2.0E-04	i	1.8E+00	r	2.0E-04	Mirex	2.7E-01	ca*	9.6E-01	ca	3.7E-03	ca	3.7E-02	ca
	2.0E-03	i	2.0E-03	r	0.1	2212-67-1	Molinate	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc
	5.0E-03	i				7439-98-7	Molybdenum	3.9E+02	nc	5.1E+03	nc			1.8E+02	nc
	1.0E-01	i	1.0E-01	r	0.1	10599-90-3	Monochloramine	6.1E+03	nc	6.2E+04	nc	3.7E+02	nc	3.6E+03	nc
	2.0E-03	i	2.0E-03	r	0.1	300-76-5	Naled	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc
	1.0E-01	i	1.0E-01	r	0.1	15299-99-7	Napropamide	6.1E+03	nc	6.2E+04	nc	3.7E+02	nc	3.6E+03	nc
	2.0E-02	i				7440-02-0	Nickel (soluble salts)	1.6E+03	nc	2.0E+04	nc			7.3E+02	nc
	8.4E-01	i					Nickel refinery dust					8.0E-03	ca		
	1.7E+00	i				12035-72-2	Nickel subsulfide			1.1E+04	ca	4.0E-03	ca		
	Tap Water PRG Based on Infant NOAEL (see IRIS)					14797-55-8	Nitrate+++							1.0E+04	nc
	Tap Water PRG Based on Infant NOAEL (see IRIS)					14797-65-0	Nitrite+++							1.0E+03	nc
	3.0E-03	p	3.0E-05	p	0.1	88-74-4	<b>2-Nitroaniline</b>	1.8E+02	nc	1.8E+03	nc	1.1E-01	nc	1.1E+02	nc
2.1E-02	p	3.0E-04	p	2.1E-02	r	3.0E-04	<b>3-Nitroaniline</b>	1.8E+01	nc	8.2E+01	ca**	3.2E-01	ca**	3.2E+00	ca**
2.1E-02	p	3.0E-03	p	2.1E-02	r	1.0E-03	<b>4-Nitroaniline</b>	2.3E+01	ca**	8.2E+01	ca*	3.2E-01	ca*	3.2E+00	ca*
	5.0E-04	i	5.7E-04	h	y	98-95-3	Nitrobenzene	2.0E+01	nc	1.0E+02	nc	2.1E+00	nc	3.4E+00	nc
	7.0E-02	h	7.0E-02	r	0.1	67-20-9	Nitrofurantoin	4.3E+03	nc	4.3E+04	nc	2.6E+02	nc	2.6E+03	nc
1.5E+00	h	1.5E+00	r		0.1	59-87-0	Nitrofurazone	3.2E-01	ca	1.1E+00	ca	4.5E-03	ca	4.5E-02	ca
1.4E-02	n	1.4E-02	r		0.1	55-63-0	Nitroglycerin	3.5E+01	ca	1.2E+02	ca	4.8E-01	ca	4.8E+00	ca
	1.0E-01	i	1.0E-01	r	0.1	556-88-7	Nitroguanidine	6.1E+03	nc	6.2E+04	nc	3.7E+02	nc	3.6E+03	nc
9.4E+00	r	5.7E-03	r	9.4E+00	h	5.7E-03	2-Nitropropane					7.2E-04	ca	1.2E-03	ca
5.4E+00	i	5.6E+00	i		y	924-16-3	N-Nitrosodi-n-butylamine	2.4E-02	ca	5.8E-02	ca	1.2E-03	ca	2.0E-03	ca
2.8E+00	i	2.8E+00	r		0.1	1116-54-7	N-Nitrosodiethanolamine	1.7E-01	ca	6.2E-01	ca	2.4E-03	ca	2.4E-02	ca
1.5E+02	i	1.5E+02	i		0.1	55-18-5	N-Nitrosodiethylamine	3.2E-03	ca	1.1E-02	ca	4.5E-05	ca	4.5E-04	ca
5.1E+01	i	8.0E-06	p	4.9E+01	i	8.0E-06	<b>N-Nitrosodimethylamine</b>	9.5E-03	ca*	3.4E-02	ca	1.4E-04	ca	1.3E-03	ca
4.9E-03	i	2.0E-02	p	4.9E-03	r	2.0E-02	<b>N-Nitrosodiphenylamine</b>	9.9E+01	ca*	3.5E+02	ca*	1.4E+00	ca*	1.4E+01	ca*
7.0E+00	i	7.0E+00	r		0.1	621-64-7	N-Nitroso di-n-propylamine	6.9E-02	ca	2.5E-01	ca	9.6E-04	ca	9.6E-03	ca

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 ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES						CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)				SOIL SCREENING LEVELS						
SFo	RfDo	SFi	RfDi	V	CAS No.		Residential Soil (mg/kg)	"Direct Contact Exposure Pathways"			"Migration to Ground Water"						
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	O abs. C soils			Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)						
2.2E+01	i	2.2E+01	r	0.1	10595-95-6	N-Nitroso-N-methylethylamine	2.2E-02	ca	7.8E-02	ca	3.1E-04	ca	3.1E-03	ca			
2.1E+00	i	2.1E+00	i	0.1	930-55-2	N-Nitrosopyrrolidine	2.3E-01	ca	8.2E-01	ca	3.1E-03	ca	3.2E-02	ca			
	2.0E-02	p	2.0E-02	r y	99-08-1	<b>m-Nitrotoluene</b>	7.3E+02	nc	1.0E+03	sat	7.3E+01	nc	1.2E+02	nc			
2.3E-01	p	1.0E-02	h	2.3E-01	r	1.0E-02	r y	88-72-2	<b>o-Nitrotoluene</b>	8.8E-01	ca	2.2E+00	ca	2.9E-02	ca	4.9E-02	ca
1.7E-02	p	1.0E-02	p	1.7E-02	r	1.0E-02	r y	99-99-0	<b>p-Nitrotoluene</b>	1.2E+01	ca*	3.0E+01	ca*	4.0E-01	ca*	6.6E-01	ca*
	4.0E-02	i	4.0E-02	r	0.1	27314-13-2	Norflurazon	2.4E+03	nc	2.5E+04	nc	1.5E+02	nc	1.5E+03	nc		
7.0E-04	i	7.0E-04	r	0.1	85509-19-9	NuStar	4.3E+01	nc	4.3E+02	nc	2.6E+00	nc	2.6E+01	nc			
3.0E-03	i	3.0E-03	r	0.1	32536-52-0	Octabromodiphenyl ether	1.8E+02	nc	1.8E+03	nc	1.1E+01	nc	1.1E+02	nc			
2.0E-03	h	2.0E-03	r	0.1	152-16-9	Octamethylpyrophosphoramidate	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc			
5.0E-02	i	5.0E-02	r	0.1	19044-88-3	Oryzalin	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc			
5.0E-03	i	5.0E-03	r	0.1	19666-30-9	Oxadiazon	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc			
2.5E-02	i	2.5E-02	r	0.1	23135-22-0	Oxamyl	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc			
3.0E-03	i	3.0E-03	r	0.1	42874-03-3	Oxyfluorfen	1.8E+02	nc	1.8E+03	nc	1.1E+01	nc	1.1E+02	nc			
1.3E-02	i	1.3E-02	r	0.1	76738-62-0	Paclobutrazol	7.9E+02	nc	8.0E+03	nc	4.7E+01	nc	4.7E+02	nc			
4.5E-03	i	4.5E-03	r	0.1	4685-14-7	Paraquat	2.7E+02	nc	2.8E+03	nc	1.6E+01	nc	1.6E+02	nc			
6.0E-03	h	6.0E-03	r	0.1	56-38-2	Parathion	3.7E+02	nc	3.7E+03	nc	2.2E+01	nc	2.2E+02	nc			
5.0E-02	h	5.0E-02	r	0.1	1114-71-2	Pebulate	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc			
4.0E-02	i	4.0E-02	r	0.1	40487-42-1	Pendimethalin	2.4E+03	nc	2.5E+04	nc	1.5E+02	nc	1.5E+03	nc			
2.3E-02	h	2.3E-02	r	0.1	87-84-3	Pentabromo-6-chloro cyclohexane	2.1E+01	ca	7.5E+01	ca	2.9E-01	ca	2.9E+00	ca			
	2.0E-03	i	2.0E-03	r	0.1	32534-81-9	Pentabromodiphenyl ether	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc		
	8.0E-04	i	8.0E-04	r	0.1	608-93-5	Pentachlorobenzene	4.9E+01	nc	4.9E+02	nc	2.9E+00	nc	2.9E+01	nc		
2.6E-01	h	3.0E-03	i	2.6E-01	r	3.0E-03	r	82-68-8	Pentachloronitrobenzene	1.9E+00	ca*	6.6E+00	ca	2.6E-02	ca	2.6E-01	ca
1.2E-01	i	3.0E-02	i	1.2E-01	r	3.0E-02	r	87-86-5	Pentachlorophenol	3.0E+00	ca	9.0E+00	ca	5.6E-02	ca	5.6E-01	ca
	1.0E-04	n			7601-90-3	Perchlorate	7.8E+00	ca/nc	1.0E+02	ca/nc			3.6E+00	ca/nc			
5.0E-02	i	5.0E-02	r	0.1	52645-53-1	Permethrin	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc			
2.5E-01	i	2.5E-01	r	0.1	13684-63-4	Phenmedipham	1.5E+04	nc	1.0E+05	max	9.1E+02	nc	9.1E+03	nc			
3.0E-01	i	3.0E-01	r	0.1	108-95-2	<b>Phenol</b>	1.8E+04	nc	1.0E+05	max	1.1E+03	nc	1.1E+04	nc			
	2.0E-03	n	2.0E-03	r	0.1	92-84-2	Phenothiazine	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc		
	6.0E-03	i	6.0E-03	r	0.1	108-45-2	m-Phenylenediamine	3.7E+02	nc	3.7E+03	nc	2.2E+01	nc	2.2E+02	nc		
4.7E-02	h	4.7E-02	r	0.1	95-54-5	<b>o-Phenylenediamine</b>	1.0E+01	ca	3.7E+01	ca	1.4E-01	ca	1.4E+00	ca			
	1.9E-01	h	1.9E-01	r	0.1	106-50-3	p-Phenylenediamine	1.2E+04	nc	1.0E+05	max	6.9E+02	nc	6.9E+03	nc		
	8.0E-05	i	8.0E-05	r	0.1	62-38-4	Phenylmercuric acetate	4.9E+00	nc	4.9E+01	nc	2.9E-01	nc	2.9E+00	nc		
1.9E-03	h	1.9E-03	r	0.1	90-43-7	2-Phenylphenol	2.5E+02	ca	8.9E+02	ca	3.5E+00	ca	3.5E+01	ca			
	2.0E-04	h	2.0E-04	r	0.1	298-02-2	Phorate	1.2E+01	nc	1.2E+02	nc	7.3E-01	nc	7.3E+00	nc		
	2.0E-02	i	2.0E-02	r	0.1	732-11-6	Phosmet	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
	3.0E-04	i	8.6E-05	i	0.1	7803-51-2	Phosphine	1.8E+01	nc	1.8E+02	nc	3.1E-01	nc	1.1E+01	nc		
			2.9E-03	i	7664-38-2	Phosphoric acid					1.0E+01	nc					
	2.0E-05	i			7723-14-0	Phosphorus (white)	1.6E+00	nc	2.0E+01	nc			7.3E-01	nc			
	1.0E+00	h	1.0E+00	r	0.1	100-21-0	p-Phthalic acid	6.1E+04	nc	1.0E+05	max	3.7E+03	nc	3.6E+04	nc		
	2.0E+00	i	3.4E-02	h	0.1	85-44-9	Phthalic anhydride	1.0E+05	max	1.0E+05	max	1.2E+02	nc	7.3E+04	nc		
	7.0E-02	i	7.0E-02	r	0.1	1918-02-1	Picloram	4.3E+03	nc	4.3E+04	nc	2.6E+02	nc	2.6E+03	nc		
	1.0E-02	i	1.0E-02	r	0.1	29232-93-7	Pirimiphos-methyl	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc		

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 ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES							CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS		
SFo	RfDo	SFi	RfDi	V	O	CAS No.		Residential Soil (mg/kg)	"Direct Contact Exposure Pathways"			"Migration to Ground Water"			
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	skin	abs.			Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)			
8.9E+00	h 7.0E-06	h 8.9E+00	r 7.0E-06	r	0.1		Polybrominated biphenyls	5.5E-02	ca**	1.9E-01	ca*	7.6E-04	ca*	7.6E-03	ca*
							Polychlorinated biphenyls (PCBs, see IRIS)								
7.0E-02	i 7.0E-05	i 7.0E-02	i 7.0E-05	r	0.14	12674-11-2	PCBs (unspeciated mixture, low risk, e.g. Aroclor 1016)	3.9E+00	nc	2.1E+01	ca**	9.6E-02	ca**	9.6E-01	ca**
2.0E+00	i 2.0E-05	i 2.0E+00	i 2.0E-05	r	0.14	11097-69-1	PCBs (unspeciated mixture, high risk, e.g. Aroclor 1254)	2.2E-01	ca**	7.4E-01	ca*	3.4E-03	ca*	3.4E-02	ca*
4.5E+00	n	4.5E+00	r		0.1	61788-33-8	Polychlorinated terphenyls	1.1E-01	ca	3.8E-01	ca	1.5E-03	ca	1.5E-02	ca
							Polynuclear aromatic hydrocarbons (PAHs)								
	6.0E-02	i	6.0E-02	r	y	83-32-9	Acenaphthene	3.7E+03	nc	2.9E+04	nc	2.2E+02	nc	3.7E+02	nc
	3.0E-01	i	3.0E-01	r	y	120-12-7	Anthracene	2.2E+04	nc	1.0E+05	max	1.1E+03	nc	1.8E+03	nc
7.3E-01	n	7.3E-01	r		0.13	56-55-3	Benz[a]anthracene	6.2E-01	ca	2.1E+00	ca	9.2E-03	ca	9.2E-02	ca
7.3E-01	n	7.3E-01	r		0.13	205-99-2	Benzo[b]fluoranthene	6.2E-01	ca	2.1E+00	ca	9.2E-03	ca	9.2E-02	ca
7.3E-02	n	7.3E-02	r		0.13	207-08-9	Benzo[k]fluoranthene	6.2E+00	ca	2.1E+01	ca	9.2E-02	ca	9.2E-01	ca
1.2E+00	c	3.9E-01	c		0.13	207-08-9	"CAL-Modified PRG"	3.8E-01	ca	1.3E+00	ca	1.7E-02	ca	5.6E-02	ca
7.3E+00	i	7.3E+00	r		0.13	50-32-8	Benzo[a]pyrene	6.2E-02	ca	2.1E-01	ca	9.2E-04	ca	9.2E-03	ca
7.3E-03	n	7.3E-03	r		0.13	218-01-9	Chrysene	6.2E+01	ca	2.1E+02	ca	9.2E-01	ca	9.2E+00	ca
1.2E-01	c	3.9E-02	c		0.13		"CAL-Modified PRG"	3.8E+00	ca	1.3E+01	ca	1.7E-01	ca	5.6E-01	ca
7.3E+00	n	7.3E+00	r		0.13	53-70-3	Dibenz[ah]anthracene	6.2E-02	ca	2.1E-01	ca	9.2E-04	ca	9.2E-03	ca
	4.0E-02	i	4.0E-02	r	0.13	206-44-0	Fluoranthene	2.3E+03	nc	2.2E+04	nc	1.5E+02	nc	1.5E+03	nc
	4.0E-02	i	4.0E-02	r	y	86-73-7	Fluorene	2.7E+03	nc	2.6E+04	nc	1.5E+02	nc	2.4E+02	nc
7.3E-01	n	7.3E-01	r		0.13	193-39-5	Indeno[1,2,3-cd]pyrene	6.2E-01	ca	2.1E+00	ca	9.2E-03	ca	9.2E-02	ca
	2.0E-02	i	8.6E-04	i	y	91-20-3	Naphthalene	5.6E+01	nc	1.9E+02	nc	3.1E+00	nc	6.2E+00	nc
1.2E-01	r	1.2E-01	c				"CAL-Modified PRG"	1.7E+00	ca	4.2E+00	ca	5.6E-02	ca	9.3E-02	ca
	3.0E-02	i	3.0E-02	r	y	129-00-0	Pyrene	2.3E+03	nc	2.9E+04	nc	1.1E+02	nc	1.8E+02	nc
1.5E-01	i	9.0E-03	i	1.5E-01	r	9.0E-03	Prochloraz	3.2E+00	ca	1.1E+01	ca	4.5E-02	ca	4.5E-01	ca
	6.0E-03	h	6.0E-03	r	0.1	26399-36-0	Profuralin	3.7E+02	nc	3.7E+03	nc	2.2E+01	nc	2.2E+02	nc
	1.5E-02	i	1.5E-02	r	0.1	1610-18-0	Prometon	9.2E+02	nc	9.2E+03	nc	5.5E+01	nc	5.5E+02	nc
	4.0E-03	i	4.0E-03	r	0.1	7287-19-6	Prometryn	2.4E+02	nc	2.5E+03	nc	1.5E+01	nc	1.5E+02	nc
	7.5E-02	i	7.5E-02	r	0.1	23950-58-5	Pronamide	4.6E+03	nc	4.6E+04	nc	2.7E+02	nc	2.7E+03	nc
	1.3E-02	i	1.3E-02	r	0.1	1918-16-7	Propachlor	7.9E+02	nc	8.0E+03	nc	4.7E+01	nc	4.7E+02	nc
	5.0E-03	i	5.0E-03	r	0.1	709-98-8	Propanil	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc
	2.0E-02	i	2.0E-02	r	0.1	2312-35-8	Propargite	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc
	2.0E-03	i	2.0E-03	r	0.1	107-19-7	Propargyl alcohol	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc
	2.0E-02	i	2.0E-02	r	0.1	139-40-2	Propazine	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc
	2.0E-02	i	2.0E-02	r	0.1	122-42-9	Propham	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc
	1.3E-02	i	1.3E-02	r	0.1	60207-90-1	Propiconazole	7.9E+02	nc	8.0E+03	nc	4.7E+01	nc	4.7E+02	nc
						98-82-8	Isopropylbenzene (see Cumene)								
	4.0E-02	n	4.0E-02	r	y	103-65-1	n-Propylbenzene	2.4E+02	sat	2.4E+02	sat	1.5E+02	nc	2.4E+02	nc
	5.0E-01	p	8.6E-04	p	0.1	57-55-6	Propylene glycol	3.0E+04	nc	1.0E+05	max	3.1E+00	nc	1.8E+04	nc
	7.0E-01	h	7.0E-01	r	0.1	52125-53-8	Propylene glycol, monoethyl ether	4.3E+04	nc	1.0E+05	max	2.6E+03	nc	2.6E+04	nc
	7.0E-01	h	5.7E-01	i	0.1	107-98-2	Propylene glycol, monomethyl ether	4.3E+04	nc	1.0E+05	max	2.1E+03	nc	2.6E+04	nc
2.4E-01	i	8.6E-03	r	1.3E-02	i	8.6E-03	Propylene oxide	1.9E+00	ca*	6.6E+00	ca*	5.2E-01	ca*	2.2E-01	ca
	2.5E-01	i	2.5E-01	r	0.1	81335-77-5	Pursuit	1.5E+04	nc	1.0E+05	max	9.1E+02	nc	9.1E+03	nc
	2.5E-02	i	2.5E-02	r	0.1	51630-58-1	Pydrin	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc

Key : SFo<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS p=PPRTV c=California EPA n=NCEA h=HEAST x=Withdrawn r=Route-extrapolation ca=Cancer PRG nc= Noncancer PRG ca\* (where: nc PRG < 100X ca PRG)  
 ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES							CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS					
SFo	RfDo	SFi	RfDi	V	O	CAS No.		Residential Soil (mg/kg)	"Direct Contact Exposure Pathways"			"Migration to Ground Water"						
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	skin	abs.			Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)						
	1.0E-03	i	1.0E-03	r	0.1	110-86-1	Pyridine	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc			
	5.0E-04	i	5.0E-04	r	0.1	13593-03-8	Quinalphos	3.1E+01	nc	3.1E+02	nc	1.8E+00	nc	1.8E+01	nc			
3.0E+00		i	3.0E+00	r		91-22-5	Quinoline	1.6E-01	ca	5.7E-01	ca	2.2E-03	ca	2.2E-02	ca			
1.1E-01	i	3.0E-03	i	1.1E-01	r	3.0E-03	r	0.1	121-82-4	RDx (Cyclonite)	4.4E+00	ca*	1.6E+01	ca	6.1E-02	ca	6.1E-01	ca
	3.0E-02	i	3.0E-02	r	0.1	10453-86-8	Resmethrin	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc			
	5.0E-02	h	5.0E-02	r	0.1	299-84-3	Ronnel	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc			
	4.0E-03	i	4.0E-03	r	0.1	83-79-4	Rotenone	2.4E+02	nc	2.5E+03	nc	1.5E+01	nc	1.5E+02	nc			
	2.5E-02	i	2.5E-02	r	0.1	78587-05-0	Savey	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc			
	5.0E-03	i			0.1	7783-00-8	Selenious Acid	3.1E+02	nc	3.1E+03	nc			1.8E+02	nc			
	5.0E-03	i				7782-49-2	Selenium	3.9E+02	nc	5.1E+03	nc			1.8E+02	nc			
	5.0E-03	h			0.1	630-10-4	Selenourea	3.1E+02	nc	3.1E+03	nc			1.8E+02	nc			
	9.0E-02	i	9.0E-02	r	0.1	74051-80-2	Sethoxydim	5.5E+03	nc	5.5E+04	nc	3.3E+02	nc	3.3E+03	nc			
	5.0E-03	i				7440-22-4	Silver and compounds	3.9E+02	nc	5.1E+03	nc			1.8E+02	nc			
1.2E-01	h	5.0E-03	i	1.2E-01	r	5.00E-03	r	0.1	122-34-9	Simazine	4.1E+00	ca*	1.4E+01	ca	5.6E-02	ca	5.6E-01	ca
	4.0E-03	i				26628-22-8	Sodium azide											
2.7E-01	h	3.0E-02	i	2.7E-01	r	3.0E-02	r	0.1	148-18-5	Sodium diethyldithiocarbamate	1.8E+00	ca	6.4E+00	ca	2.5E-02	ca	2.5E-01	ca
	2.0E-05	i	2.0E-05	r	0.1	62-74-8	Sodium fluoroacetate	1.2E+00	nc	1.2E+01	nc	7.3E-02	nc	7.3E-01	nc			
	1.0E-03	h	1.0E-03	r	0.1	13718-26-8	Sodium metavanadate	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc			
	6.0E-01	i				7440-24-6	Strontium, stable	4.7E+04	nc	1.0E+05	max			2.2E+04	nc			
	3.0E-04	i	3.0E-04	r	0.1	57-24-9	Styrene	1.8E+01	nc	1.8E+02	nc	1.1E+00	nc	1.1E+01	nc			
	2.0E-01	i	2.9E-01	i	y	100-42-5	Styrene	1.7E+03	sat	1.7E+03	sat	1.1E+03	nc	1.6E+03	nc			
	5.0E-03	p	5.0E-03	r		80-07-9	<b>1,1'-Sulfonylbis (4-chlorobenzene)</b>	3.9E+02	nc	5.1E+03	nc	1.8E+01	nc	1.8E+02	nc			
	2.5E-02	i	2.5E-02	r	0.1	88671-89-0	Systhane	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc			
1.5E+05	h	1.5E+05	h		0.03	1746-01-6	2,3,7,8-TCDD (dioxin)	3.9E-06	ca	1.6E-05	ca	4.5E-08	ca	4.5E-07	ca			
	7.0E-02	i	7.0E-02	r	0.1	34014-18-1	Tebuthiuron	4.3E+03	nc	4.3E+04	nc	2.6E+02	nc	2.6E+03	nc			
	2.0E-02	h	2.0E-02	r	0.1	3383-96-8	Temephos	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc			
	1.3E-02	i	1.3E-02	r	0.1	5902-51-2	Terbacil	7.9E+02	nc	8.0E+03	nc	4.7E+01	nc	4.7E+02	nc			
	2.5E-05	h	2.5E-05	r	0.1	13071-79-9	Terbufos	1.5E+00	nc	1.5E+01	nc	9.1E-02	nc	9.1E-01	nc			
	1.0E-03	i	1.0E-03	r	0.1	886-50-0	Terbutryn	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc			
	3.0E-04	i	3.0E-04	r	0.1	95-94-3	1,2,4,5-Tetrachlorobenzene	1.8E+01	nc	1.8E+02	nc	1.1E+00	nc	1.1E+01	nc			
2.6E-02	i	3.0E-02	i	2.6E-02	i	3.0E-02	r	y	630-20-6	1,1,1,2-Tetrachloroethane	3.2E+00	ca	7.3E+00	ca	2.6E-01	ca	4.3E-01	ca
2.0E-01	i	6.0E-02	p	2.0E-01	i	6.0E-02	r	y	79-34-5	1,1,2,2-Tetrachloroethane	4.1E-01	ca	9.3E-01	ca	3.3E-02	ca	5.5E-02	ca
5.4E-01	c	1.0E-02	i	2.1E-02	c	1.0E-02	c	y	127-18-4	<b>Tetrachloroethylene (PCE)</b>	4.8E-01	ca*	1.3E+00	ca	3.2E-01	ca	1.0E-01	ca
	3.0E-02	i	3.0E-02	r	0.1	58-90-2	2,3,4,6-Tetrachlorophenol	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc			
2.0E+01	h	2.0E+01	r		0.1	5216-25-1	p,a,a,a-Tetrachlorotoluene	2.4E-02	ca	8.6E-02	ca	3.4E-04	ca	3.4E-03	ca			
2.4E-02	h	3.0E-02	i	2.4E-02	r	3.0E-02	r	0.1	961-11-5	Tetrachlorovinphos	2.0E+01	ca*	7.2E+01	ca	2.8E-01	ca	2.8E+00	ca
	5.0E-04	i	5.0E-04	r	0.1	3689-24-5	Tetraethylthiopyrophosphate	3.1E+01	nc	3.1E+02	nc	1.8E+00	nc	1.8E+01	nc			
7.6E-03	n	2.1E-01	n	6.8E-03	n	8.6E-02	n	y	109-99-9	Tetrahydrofuran	9.4E+00	ca	2.1E+01	ca	9.9E-01	ca	1.6E+00	ca
	6.6E-05	i				7440-28-0	Thallium and compounds+++	5.2E+00	nc	6.7E+01	nc			2.4E+00	nc			
	1.0E-02	i	1.0E-02	r	0.1	28249-77-6	Thiobencarb	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc			
	5.0E-04	n	5.0E-04	r	0.1	N/A	Thiocyanate	3.1E+03	nc	1.0E+05	max	1.8E+02	nc	1.8E+03	nc			
	3.0E-04	h	3.0E-04	r	0.1	39196-18-4	Thiofanox	1.8E+01	nc	1.8E+02	nc	1.1E+00	nc	1.1E+01	nc			

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 ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES							CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS						
SFo	RfDo	SFi	RfDi	V	O	CAS No.		Residential Soil (mg/kg)	"Direct Contact Exposure Pathways"			"Migration to Ground Water"							
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	skin	abs.			Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)							
	8.0E-02	i	8.0E-02	r	0.1	23564-05-8	Thiophanate-methyl	4.9E+03	nc	4.9E+04	nc	2.9E+02	nc	2.9E+03	nc				
	5.0E-03	i	5.0E-03	r	0.1	137-26-8	Thiram	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc				
	6.0E-01	h				7440-31-5	Tin (inorganic, also see tributyltin oxide)	4.7E+04	nc	1.0E+05	max			2.2E+04	nc				
	4.0E+00	n	8.6E-03	n		7440-32-6	Titanium	1.0E+05	max	1.0E+05	max	3.1E+01	nc	1.5E+05	nc				
	2.0E-01	i	1.1E-01	i	y	108-88-3	Toluene	5.2E+02	sat	5.2E+02	sat	4.0E+02	nc	7.2E+02	nc	1.2E+01 6.0E-01			
3.2E+00	h	3.2E+00	r		0.1	95-80-7	Toluene-2,4-diamine	1.5E-01	ca	5.4E-01	ca	2.1E-03	ca	2.1E-02	ca				
	6.0E-01	h	6.0E-01	r	0.1	95-70-5	Toluene-2,5-diamine	3.7E+04	nc	1.0E+05	max	2.2E+03	nc	2.2E+04	nc				
	2.0E-01	h	2.0E-01	r	0.1	823-40-5	Toluene-2,6-diamine	1.2E+04	nc	1.0E+05	max	7.3E+02	nc	7.3E+03	nc				
1.9E-01	i	1.9E-01	r		0.1	106-49-0	p-Toluidine	2.6E+00	ca	9.1E+00	ca	3.5E-02	ca	3.5E-01	ca				
1.1E+00	i	1.1E+00	i		0.1	8001-35-2	Toxaphene	4.4E-01	ca	1.6E+00	ca	6.0E-03	ca	6.1E-02	ca	3.1E+01 2.0E+00			
	7.5E-03	i	7.5E-03	r	0.1	66841-25-6	Tralometrin	4.6E+02	nc	4.6E+03	nc	2.7E+01	nc	2.7E+02	nc				
	1.3E-02	i	1.3E-02	r	0.1	2303-17-5	Triallate	7.9E+02	nc	8.0E+03	nc	4.7E+01	nc	4.7E+02	nc				
	1.0E-02	i	1.0E-02	r	0.1	82097-50-5	Triasulfuron	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc				
	5.0E-03	i	5.0E-03	r	0.1	615-54-3	1,2,4-Tribromobenzene	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc				
9.2E-03	p	2.0E-01	p	9.2E-03	r	2.0E-01	r	0.1	126-73-8	<b>Tributyl phosphate</b>	5.3E+01	ca	1.9E+02	ca	7.3E-01	ca	7.3E+00	ca	
	3.0E-04	i			0.1	56-35-9	Tributyltin oxide (TBTO)	1.8E+01	nc	1.8E+02	nc			1.1E+01	nc				
3.4E-02	h	3.4E-02	r		0.1	634-93-5	2,4,6-Trichloroaniline	1.4E+01	ca	5.1E+01	ca	2.0E-01	ca	2.0E+00	ca				
2.9E-02	h	2.9E-02	r		0.1	33663-50-2	2,4,6-Trichloroaniline hydrochloride	1.7E+01	ca	5.9E+01	ca	2.3E-01	ca	2.3E+00	ca				
	1.0E-02	i	1.0E-03	p	y	120-82-1	<b>1,2,4-Trichlorobenzene</b>	6.2E+01	nc	2.2E+02	nc	3.7E+00	nc	7.2E+00	nc	5.0E+00 3.0E-01			
	2.8E-01	n	6.3E-01	p	y	71-55-6	1,1,1-Trichloroethane	1.2E+03	sat	1.2E+03	sat	2.3E+03	nc	3.2E+03	nc	2.0E+00 1.0E-01			
5.7E-02	i	4.0E-03	i	5.6E-02	i	4.0E-03	r	y	79-00-5	1,1,2-Trichloroethane	7.3E-01	ca*	1.6E+00	ca*	1.2E-01	ca	2.0E-01	ca	2.0E-02 9.0E-04
4.0E-01	n	3.0E-04	n	4.0E-01	n	1.0E-02	n	y	79-01-6	Trichloroethylene (TCE)	5.3E-02	ca	1.1E-01	ca	1.7E-02	ca	2.8E-02	ca	6.0E-02 3.0E-03
1.3E-02	c	7.0E-03	c	1.7E-01	c	y	79-01-6	<b>"CAL-Modified PRG"</b>	2.9E+00	ca	6.5E+00	ca	9.6E-01	ca	1.4E+00	ca			
	3.0E-01	i	2.0E-01	h	y	75-69-4	Trichlorofluoromethane	3.9E+02	nc	2.0E+03	sat	7.3E+02	nc	1.3E+03	nc				
1.1E-02	i	1.0E-04	n	1.1E-02	i	1.0E-04	r	0.1	95-95-4	2,4,5-Trichlorophenol	6.1E+03	nc	6.2E+04	nc	3.7E+02	nc	3.6E+03	nc	2.7E+02 1.4E+01
	1.0E-04	n	1.1E-02	i	1.0E-04	r	0.1	88-06-2	2,4,6-Trichlorophenol	6.1E+00	nc**	6.2E+01	nc**	3.7E-01	nc**	3.6E+00	nc**	2.0E-01 8.0E-03	
7.0E-02	c	7.0E-02	c		0.1	88-06-2	<b>"CAL-Modified PRG"</b>	6.9E+00	ca	2.5E+01	ca	9.6E-02	ca	9.6E-01	ca				
	1.0E-02	i	1.0E-02	r	0.1	93-76-5	2,4,5-Trichlorophenoxyacetic Acid	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc				
	8.0E-03	i	8.0E-03	r	0.1	93-72-1	2-(2,4,5-Trichlorophenoxy) propionic acid	4.9E+02	nc	4.9E+03	nc	2.9E+01	nc	2.9E+02	nc				
	5.0E-03	i	5.0E-03	r	y	598-77-6	<b>1,1,2-Trichloropropane</b>	7.1E+01	nc	2.7E+02	nc	1.8E+01	nc	3.0E+01	nc				
2.0E+00	n	6.0E-03	i	2.0E+00	r	1.4E-03	n	y	96-18-4	<b>1,2,3-Trichloropropane</b>	3.4E-02	ca	7.6E-02	ca	3.4E-03	ca	5.6E-03	ca	
	1.0E-02	p	3.0E-04	p	y	96-19-5	<b>1,2,3-Trichloropropene</b>	5.2E+00	nc	1.7E+01	nc	1.1E+00	nc	2.2E+00	nc				
	3.0E-03	i	3.0E-03	r	0.1	58138-08-2	Triphane	1.8E+02	nc	1.8E+03	nc	1.1E+01	nc	1.1E+02	nc				
7.7E-03	i	7.5E-03	i	7.7E-03	r	7.5E-03	r	0.1	121-44-8	Triethylamine	2.3E+01	nc	8.6E+01	nc	7.3E+00	nc	1.2E+01	nc	
	1.4E-04	r	1.4E-04	n	0.1	552-30-7	Triethylamine	6.3E+01	ca**	2.2E+02	ca*	8.7E-01	ca*	8.7E+00	ca*				
	5.0E-02	p	1.7E-03	p	y	95-63-6	Trimellitic Anhydride (TMAN)	8.6E+00	nc	8.6E+01	nc	5.1E-01	nc	5.1E+00	nc				
3.7E-02	h	3.7E-02	r		0.1	512-56-1	1,2,4-Trimethylbenzene	5.2E+01	nc	1.7E+02	nc	6.2E+00	nc	1.2E+01	nc				
	5.0E-02	p	1.7E-03	p	y	108-67-8	1,3,5-Trimethylbenzene	2.1E+01	nc	7.0E+01	nc	6.2E+00	nc	1.2E+01	nc				
	3.0E-02	i	3.0E-02	r	0.1	99-35-4	Trimethyl phosphate	1.3E+01	ca	4.7E+01	ca	1.8E-01	ca	1.8E+00	ca				
	1.0E-02	h	1.0E-02	r	0.1	479-45-8	1,3,5-Trinitrobenzene	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc				
3.0E-02	i	5.0E-04	i	3.0E-02	r	5.0E-04	r	0.1	118-96-7	Trinitrophenylmethylnitramine	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc	
	5.0E-04	i	3.0E-02	r	5.0E-04	r	0.1	118-96-7	2,4,6-Trinitrotoluene	1.6E+01	ca**	5.7E+01	ca**	2.2E-01	ca**	2.2E+00	ca**		

Key : SFO,i=Cancer Slope Factor oral, inhalation RfDo,i=Reference Dose oral, inhalation i=IRIS p=PPRTV c=California EPA n=NCEA h=HEAST x=Withdrawn r=Route-extrapolation ca=Cancer PRG nc= Noncancer PRG ca\* (where: nc PRG < 100X ca PRG)  
 ca\*\* (where nc PRG < 10X ca PRG) +++=Non-Standard Method Applied (See User's Guide) sat=Soil Saturation (See User's Guide) max=Ceiling limit (See User's Guide) DAF=Dilution Attenuation Factor (See User's Guide) CAS=Chemical Abstract Services

TOXICITY VALUES							CONTAMINANT	PRELIMINARY REMEDIATION GOALS (PRGs)					SOIL SCREENING LEVELS			
SFO	RfDo	SFi	RfDi	V	CAS No.	Residential Soil (mg/kg)		"Direct Contact Exposure Pathways"				"Migration to Ground Water"				
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	O C		Soil (mg/kg)	Industrial Soil (mg/kg)	Ambient Air (ug/m^3)	Tap Water (ug/l)		DAF 20 (mg/kg)	DAF 1 (mg/kg)				
	2.0E-02	p	2.0E-02	r	0.1	791-28-6	<b>Triphenylphosphine oxide</b>	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc	
1.4E-02	p	3.1E-01	p	1.4E-02	r	3.1E-01	<b>Tris(2-chloroethyl) phosphate</b>	3.5E+01	ca	1.2E+02	ca	4.8E-01	ca	4.8E+00	ca	
3.2E-03	p	1.0E-01	p	3.2E-03	r	1.0E-01	<b>Tris(2-ethylhexyl) phosphate</b>	1.5E+02	ca*	5.4E+02	ca	2.1E+00	ca	2.1E+01	ca	
	2.0E-04	n				7440-61-1	Uranium (chemical toxicity only)	1.6E+01	nc	2.0E+02	nc			7.3E+00	nc	
	1.0E-03	n				7440-62-2	Vanadium and compounds	7.8E+01	nc	1.0E+03	nc			3.6E+01	nc	
	1.0E-03	i	1.0E-03	r	0.1	1929-77-7	Vernam	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc	
	2.5E-02	i	2.5E-02	r	0.1	50471-44-8	Vinclozolin	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc	
	1.0E+00	h	5.7E-02	i	y	108-05-4	Vinyl acetate	4.3E+02	nc	1.4E+03	nc	2.1E+02	nc	4.1E+02	nc	
1.1E-01	r	8.6E-04	r	1.1E-01	h	8.6E-04	Vinyl bromide (bromoethene)	1.9E-01	ca*	4.2E-01	ca*	6.1E-02	ca*	1.0E-01	ca*	
1.5E+00	i	3.0E-03	i	3.1E-02	i	2.9E-02	Vinyl chloride (child/adult)+++	7.9E-02	ca			1.1E-01	ca	2.0E-02	ca	
7.5E-01	i	3.0E-03	i	1.6E-02	i	2.9E-02	Vinyl chloride (adult)			7.5E-01	ca					
	3.0E-04	i	3.0E-04	r	0.1	81-81-2	Warfarin	1.8E+01	nc	1.8E+02	nc	1.1E+00	nc	1.1E+01	nc	
	2.0E-01	i	2.9E-02	i	y	0.1	1330-20-7	<b>Xylenes</b>	2.7E+02	nc	4.2E+02	sat	1.1E+02	nc	2.1E+02	nc
	3.0E-01	i				7440-66-6	Zinc	2.3E+04	nc	1.0E+05	max			1.1E+04	nc	
	3.0E-04	i				1314-84-7	Zinc phosphide	2.3E+01	nc	3.1E+02	nc			1.1E+01	nc	
	5.0E-02	i	5.0E-02	r	0.1	12122-67-7	Zineb	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc	