

Key : SFO<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca) +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION										CONTAMINANT	PRELIMINARY REMEDIAL GOALS (PRGs)				SOIL SCREENING LEVELS				
SFO	RfDo	SFi	RfDi	V	skin		CAS No.	"Direct Contact Exposure Pathways"				"Migration to Ground Water"							
1/(mg/kg-d)	(mg/kg-d)	1/(mg/kg-d)	(mg/kg-d)	O	abs.			Residential	Industrial	Ambient Air	Tap Water	DAF 20	DAF 1						
				C	soils			Soil (mg/kg)	Soil (mg/kg)	(ug/m <sup>3</sup> )	(ug/l)	(mg/kg)	(mg/kg)						
8.7E-03	i	4.0E-03	i	8.7E-03	r	4.0E-03	r	0	0.10	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	1		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	0.10	34256-82-1	Acetochlor	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc
		1.0E-01	i			1.0E-01	r	1		67-64-1	Acetone	1.6E+03	nc	6.0E+03	nc	3.7E+02	nc	6.1E+02	nc
		8.0E-04	h			8.0E-04	r	0	0.10	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	1		75-05-8	Acetonitrile	4.2E+02	nc	1.8E+03	nc	6.2E+01	nc	1.0E+02	nc
		2.0E-02	h			5.7E-06	i	1		107-02-8	Acrolein	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	0.10	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	0.10	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	1		107-13-1	Acrylonitrile	2.1E-01	ca*	4.9E-01	ca*	2.8E-02	ca*	3.9E-02	ca*
8.1E-02	h	1.0E-02	i	8.0E-02	r	1.0E-02	r	0	0.10	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	0.10	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	0.10	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	0.10	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	0.10	309-00-2	Aldrin	2.9E-02	ca*	1.0E-01	ca	3.9E-04	ca	4.0E-03	ca
		2.5E-01	i			2.5E-01	r	0	0.10	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	0.10	107-18-6	Allyl alcohol	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc
		5.0E-02	h			2.9E-04	i	0	0.10	107-05-1	Allyl chloride	3.0E+03	nc	3.0E+04	nc	1.0E+00	nc	1.8E+03	nc
		1.0E+00	n			1.4E-03	n	0		7429-90-5	Aluminum	7.6E+04	nc	1.0E+05	max	5.1E+00	nc	3.6E+04	nc
		4.0E-04	i					0		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	0.10	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	0.10	834-12-8	Ametryn	5.5E+02	nc	5.5E+03	nc	3.3E+01	nc	3.3E+02	nc
		7.0E-02	h			7.0E-02	r	0	0.10	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	0.10	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	0.10	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	0.10	7773-06-0	Ammonium sulfamate	1.2E+04	nc	1.0E+05	max			7.3E+03	nc
5.7E-03	i	7.0E-03	n	5.7E-03	r	2.9E-04	i	0	0.10	62-53-3	Aniline	8.5E+01	ca**	3.0E+02	ca*	1.0E+00	nc	1.2E+01	ca*
		4.0E-04	i					0		7440-36-0	Antimony and compounds	3.1E+01	nc	4.1E+02	nc			1.5E+01	nc
		5.0E-04	h					0		1314-60-9	Antimony pentoxide	3.9E+01	nc	5.1E+02	nc			1.8E+01	nc
		9.0E-04	h					0		28300-74-5	Antimony potassium tartrate	7.0E+01	nc	9.2E+02	nc			3.3E+01	nc
		4.0E-04	h					0		1332-81-6	Antimony tetroxide	3.1E+01	nc	4.1E+02	nc			1.5E+01	nc
		4.0E-04	h			5.7E-05	i	0		1309-64-4	Antimony trioxide	3.1E+01	nc	4.1E+02	nc	2.1E-01	nc	1.5E+01	nc
		1.3E-02	i			1.3E-02	r	0	0.10	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	0.10	140-57-8	Aramite	1.9E+01	ca	6.9E+01	ca	2.7E-01	ca	2.7E+00	ca
		3.0E-04	i					0	0.03	7440-38-2	Arsenic (noncancer endpoint)	2.2E+01	nc	2.6E+02	nc				

Key : SFo<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca)  
 +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION						CONTAMINANT		PRELIMINARY REMEDIAL 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.		"Direct Contact Exposure Pathways"				"Migration to Ground Water"			
								Residential Soil (mg/kg)	Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)		
1.5E+00	i 3.0E-04	i 1.5E+01	i		0	0.03	7440-38-2	Arsenic (cancer endpoint)	3.9E-01	ca* 1.6E+00	ca 4.5E-04	ca 4.5E-02	ca	2.9E+01	1.0E+00
							7784-42-1	Arsine (see arsenic for cancer endpoint)			5.2E-02	nc			
	9.0E-03	i				0.10	76578-12-6	Assure	5.5E+02	nc 5.5E+03	nc 3.3E+01	nc 3.3E+02	nc		
	5.0E-02	i				0.10	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	h 2.2E-01	r			0.10	1912-24-9	Atrazine	2.2E+00	ca 7.8E+00	ca 3.1E-02	ca 3.0E-01	ca		
	4.0E-04	i				0.10	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			0.10	103-33-3	Azobenzene	4.4E+00	ca 1.6E+01	ca 6.2E-02	ca 6.1E-01	ca		
	7.0E-02	i				0	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
	4.0E-03	i				0.10	114-26-1	Baygon	2.4E+02	nc 2.5E+03	nc 1.5E+01	nc 1.5E+02	nc		
	3.0E-02	i				0.10	43121-43-3	Bayleton	1.8E+03	nc 1.8E+04	nc 1.1E+02	nc 1.1E+03	nc		
	2.5E-02	i				0.10	68359-37-5	Baythroid	1.5E+03	nc 1.5E+04	nc 9.1E+01	nc 9.1E+02	nc		
	3.0E-01	i				0.10	1861-40-1	Benefin	1.8E+04	nc 1.0E+05	max 1.1E+03	nc 1.1E+04	nc		
	5.0E-02	i				0.10	17804-35-2	Benomyl	3.1E+03	nc 3.1E+04	nc 1.8E+02	nc 1.8E+03	nc		
	3.0E-02	i				0.10	25057-89-0	Bentazon	1.8E+03	nc 1.8E+04	nc 1.1E+02	nc 1.1E+03	nc		
	1.0E-01	i				0.10	100-52-7	Benzaldehyde	6.1E+03	nc 6.2E+04	nc 3.7E+02	nc 3.6E+03	nc		
5.5E-02	i 3.0E-03	n 2.9E-02	i			1	71-43-2	Benzene	6.0E-01	ca* 1.3E+00	ca* 2.3E-01	ca* 3.4E-01	ca*	3.0E-02	2.0E-03
2.3E+02	i 3.0E-03	i 2.3E+02	i			0.10	92-87-5	Benzidine	2.1E-03	ca 7.5E-03	ca 2.9E-05	ca 2.9E-04	ca		
	4.0E+00	i				0.10	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			0.10	98-07-7	Benzotrithloride	3.7E-02	ca 1.3E-01	ca 5.2E-04	ca 5.2E-03	ca		
	3.0E-01	h				0.10	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			1	100-44-7	<b>Benzyl chloride</b>	8.9E-01	ca* 2.2E+00	ca 4.0E-02	ca 6.6E-02	ca		
	2.0E-03	i 8.4E+00	i			0	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				0.10	141-66-2	Bidrin	6.1E+00	nc 6.2E+01	nc 3.7E-01	nc 3.6E+00	nc		
	1.5E-02	i				0.10	82657-04-3	Biphenthrin (Talstar)	9.2E+02	nc 9.2E+03	nc 5.5E+01	nc 5.5E+02	nc		
	5.0E-02	i				1	92-52-4	1,1-Biphenyl	3.5E+02	sat 3.5E+02	sat 1.8E+02	nc 3.0E+02	nc		
1.1E+00	i	1.2E+00	i			1	111-44-4	Bis(2-chloroethyl)ether	2.1E-01	ca 5.5E-01	ca 5.8E-03	ca 9.8E-03	ca	4.0E-04	2.0E-05
7.0E-02	x 4.0E-02	i 3.5E-02	x			1	39638-32-9	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			1	542-88-1	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			1	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			0.10	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				0.10	80-05-7	Bisphenol A	3.1E+03	nc 3.1E+04	nc 1.8E+02	nc 1.8E+03	nc		
	2.0E-01	i				0	7440-42-8	<b>Boron</b>	1.6E+04	nc 1.0E+05	max 2.1E+01	nc 7.3E+03	nc		
						0	7637-07-2	Boron trifluoride			7.3E-01	nc			
	4.00E-03	i					15541-45-4	<b>Bromate</b>	3.1E+02	nc 4.1E+03	nc 0.0E+00	1.5E+02	nc		
	2.0E-02	n				1	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			1	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

Key : SFo,i=Cancer Slope Factor oral, inhalation RfDo,i=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca) +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION								CONTAMINANT	PRELIMINARY REMEDIAL 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.	"Direct Contact Exposure Pathways"				"Migration to Ground Water"					
							Residential Soil (mg/kg)	Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)				
7.9E-03	i 2.0E-02	i 3.9E-03	i 2.0E-02	r 0	0.10	75-25-2	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 1		74-83-9	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	0.10	2104-96-3	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	0.10	1689-84-5	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	0.10	1689-99-2	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
9.8E-01	r	9.8E-01	i		1	106-99-0	6.5E-03	ca	1.4E-02	ca	6.9E-03	ca	1.1E-02	ca		
	1.0E-01	i	2.6E-03	n 0	0.10	71-36-3	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	0.10	2008-41-5	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc		
	4.00E-02	n	4.00E-02	r 1		104-51-8	2.4E+02	sat	2.4E+02	sat	1.5E+02	nc	2.4E+02	nc		
	4.00E-02	n	4.00E-02	r 1		135-9-88	2.2E+02	sat	2.2E+02	sat	1.5E+02	nc	2.4E+02	nc		
	4.00E-02	n	4.00E-02	r 1		98-06-6	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	0.10	85-68-7	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	0.10	85-70-1	6.1E+04	nc	1.0E+05	max	3.7E+03	nc	3.6E+04	nc		
2.5E-01	h 3.0E-04	h 2.5E-01	r 3.0E-04	r 0	0.10	75-60-5	1.9E+00	ca**	6.9E+00	ca*	2.7E-02	ca*	2.7E-01	ca*		
	5.0E-04	i 6.3E+00	i		0.001	7440-43-9	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	0.10	105-60-2	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	0.10	2425-06-1	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	0.10	133-06-2	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	0.10	63-25-2	6.1E+03	nc	6.2E+04	nc	4.0E+02	nc	3.6E+03	nc		
2.0E-02	h	2.0E-02	r		0.10	86-74-8	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	0.10	1563-66-2	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc		
	1.0E-01	i	2.0E-01	i 1		75-15-0	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 1		56-23-5	2.5E-01	ca**	5.5E-01	ca*	1.3E-01	ca*	1.7E-01	ca*	7.0E-02	3.0E-03
	1.0E-02	i	1.0E-02	r 0	0.10	55285-14-8	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	0.10	5234-68-4	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	0.10	133-90-4	9.2E+02	nc	9.2E+03	nc	5.5E+01	nc	5.5E+02	nc		
4.0E-01	h	4.0E-01	r		0.10	118-75-2	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	0.04	12789-03-6	1.6E+00	ca*	6.5E+00	ca*	1.9E-02	ca*	1.9E-01	ca*	1.0E+01	5.0E-01
	2.0E-02	i	2.0E-02	r 0	0.10	90982-32-4	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
	1.0E-01	i	5.71E-05	n		7782-50-5				2.1E-01	nc					
			5.7E-05	i		10049-04-4				2.1E-01	nc					
	2.0E-03	h	2.0E-03	r 0	0.10	79-11-8	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc		
	8.6E-06	r	8.6E-06	i 1		532-27-4	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	0.10	106-47-8	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 1		108-90-7	1.5E+02	nc	5.3E+02	nc	6.2E+01	nc	1.1E+02	nc	1.0E+00	7.0E-02

Key : SFO<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca) +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION										CONTAMINANT	PRELIMINARY REMEDIAL GOALS (PRGs)				SOIL SCREENING LEVELS						
SFO <sub>1</sub> (mg/kg-d)	RfDo (mg/kg-d)	SFi 1/(mg/kg-d)	RfDi (mg/kg-d)	V	skin	O	abs.	CAS No.	C		soils	Residential Soil (mg/kg)	Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)				
2.7E-01	h	2.0E-02	i	2.7E-01	h	2.0E-02	r	0	0.10	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	0.10	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	0.10	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	1		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	1		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	1		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	1		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	1		75-00-3	Chloroethane	3.0E+00	ca	6.5E+00	ca	2.3E+00	ca	4.6E+00	ca		
		1.0E-02	i			8.6E-04	n	1		67-66-3	Chloroform	3.6E+00	ca/nc	1.2E+01	ca/nc	3.1E+00	ca/nc	6.2E+00	ca/nc	6.0E-01	3.0E-02
3.1E-02				1.9E-02				1			Chloroform "CAL-Modified PRG"	9.4E-01	ca	2.0E+00	ca	3.5E-01	ca	5.3E-01	ca		
1.3E-02	h			6.3E-03	h	8.6E-02	n	1		74-87-3	Chloromethane	1.2E+00	ca	2.6E+00	ca	1.1E+00	ca	1.5E+00	ca		
5.8E-01	h			5.8E-01	r			0	0.10	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	0.10	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	1		91-58-7	beta-Chloronaphthalene	4.9E+03	nc	2.3E+04	nc	2.9E+02	nc	4.9E+02	nc		
9.7E-03	h	1.0E-03	h	9.7E-03	r	2.0E-05	h	1		88-73-3	o-Chloronitrobenzene	1.4E+00	nc**	4.5E+00	nc**	7.3E-02	nc**	1.5E-01	nc**		
6.7E-03	h	1.0E-03	h	6.7E-03	r	1.7E-04	h	1		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	1		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	1		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	0.10	1897-45-6	Chloroethanol	4.4E+01	ca*	1.6E+02	ca*	6.1E-01	ca*	6.1E+00	ca*		
		2.0E-02	i			2.0E-02	r	1		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	0.10	101-21-3	Chlorpropham	1.2E+04	nc	1.0E+05	max	7.3E+02	nc	7.3E+03	nc		
		3.0E-03	i			3.0E-03	r	0	0.10	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	0.10	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	0.10	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	0.10	60238-56-4	Chlorthiophos	4.9E+01	nc	4.9E+02	nc	2.9E+00	nc	2.9E+01	nc		
				4.2E+01	i			0			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	0.0E+00		5.5E+04	nc		
		3.0E-03	i	2.9E+02	i	2.2E-06	i	0		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
		2.00E-02	n	9.8E+00	n	5.7E-06	n			7440-48-4	Cobalt	9.0E+02	ca**	1.9E+03	ca*	6.9E-04	ca*	7.3E+02	nc		
				2.2E+00	i			0		8007-45-2	Coke Oven Emissions					3.1E-03	ca				
		4.00E-02	h					0		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			1		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	1		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	0.10	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	0.10	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	1		74-90-8	Cyanide (hydrogen)	1.1E+01	nc	3.5E+01	nc	3.1E+00	nc	6.2E+00	nc		

Key : SFo<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca)  
 +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION						CONTAMINANT	PRELIMINARY REMEDIAL 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.	"Direct Contact Exposure Pathways"				"Migration to Ground Water"									
							Residential Soil (mg/kg)	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-02	i	4.0E-02	r	1	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	1	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	1	506-77-4	Cyanogen chloride	1.6E+02	nc	5.4E+02	nc	1.8E+02	nc	3.0E+02	nc						
	5.7E+00	r	5.7E+00	n	1	110-82-7	Cyclohexane	1.4E+02	sat	1.4E+02	sat	2.1E+04	nc	3.5E+04	nc						
	5.0E+00	i	5.0E+00	r	0	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	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	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	0.03	50-29-3	DDT	1.7E+00	ca*	7.0E+00	ca*	2.0E-02	ca*	2.0E-01	ca*	3.2E+01	2.0E+00
	1.0E-02	i	1.0E-02	r	0	0.10	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	0.10	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	0.10	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	0.10	333-41-5	Diazinon	5.5E+01	nc	5.5E+02	nc	3.3E+00	nc	3.3E+01	nc					
	4.0E-03	n	4.0E-03	r	1	132-64-9	Dibenzofuran	2.9E+02	nc	3.1E+03	nc	1.5E+01	nc	2.4E+01	nc						
	1.0E-02	i	1.0E-02	r	0	0.10	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	1	124-48-1	Dibromochloromethane	1.1E+00	ca	2.6E+00	ca	8.0E-02	ca	1.3E-01	ca	4.0E-01	2.0E-02	
1.4E+00	h	5.7E-05	r	2.4E-03	x	5.7E-05	i	1	96-12-8	1,2-Dibromo-3-chloropropane	4.5E-01	ca**	2.0E+00	ca**	2.1E-01	nc	4.8E-02	ca**			
7.0E+00		7.0E+00			1	96-12-8	"CAL-Modified PRG"	1.9E-02	ca	4.6E-02	ca	9.6E-04	ca	1.6E-03	ca						
8.5E+01	i	5.7E-05	r	7.7E-01	i	5.7E-05	h	1	106-93-4	1,2-Dibromoethane	6.9E-03	ca	2.8E-02	ca*	8.7E-03	ca*	7.6E-04	ca			
	1.0E-01	i	1.0E-01	r	0	0.10	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	0.10	1918-00-9	Dicamba	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc					
	9.0E-02	i	5.7E-02	h	1	95-50-1	1,2-Dichlorobenzene	3.7E+02	sat	3.7E+02	sat	2.1E+02	nc	3.7E+02	nc	1.7E+01	9.0E-01				
	9.00E-04	n	9.00E-04	r	1	541-73-1	1,3-Dichlorobenzene	1.6E+01	nc	6.3E+01	nc	3.3E+00	nc	5.5E+00	nc						
2.4E-02	h	3.00E-02	n	2.2E-02	n	3.00E-02	i	1	106-46-7	1,4-Dichlorobenzene	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	0.10	91-94-1	3,3-Dichlorobenzidine	1.1E+00	ca	3.8E+00	ca	1.5E-02	ca	1.5E-01	ca	7.0E-03	3.0E-04			
	3.00E-02	n	3.00E-02	r	0.10	90-98-2	4,4'-Dichlorobenzophenone	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc						
9.3E+00	r	9.3E+00	h		1	764-41-0	1,4-Dichloro-2-butene	7.9E-03	ca	1.8E-02	ca	7.2E-04	ca	1.2E-03	ca						
	2.0E-01	i	5.7E-02	h	1	75-71-8	Dichlorodifluoromethane	9.4E+01	nc	3.1E+02	nc	2.1E+02	nc	3.9E+02	nc						
	1.0E-01	h	1.4E-01	h	1	75-34-3	1,1-Dichloroethane	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		5.7E-03			1		"CAL-Modified PRG"	2.8E+00	ca	6.0E+00	ca	1.2E+00	ca	2.0E+00	ca						

Key : SFO<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca) +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION							CONTAMINANT		PRELIMINARY REMEDIAL GOALS (PRGs)				SOIL SCREENING LEVELS								
SFO <sub>i</sub> 1/(mg/kg-d)	RfDo (mg/kg-d)	SFi 1/(mg/kg-d)	RfDi (mg/kg-d)	V	skin	CAS No.	"Direct Contact Exposure Pathways"							"Migration to Ground Water"							
				O	abs.		Residential Soil (mg/kg)	Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)									
9.1E-02	i	3.0E-02	n	9.1E-02	i	1.4E-03	n	1	107-06-2	1,2-Dichloroethane (EDC)	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	1	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	h			1.0E-02	r	1	156-59-2	1,2-Dichloroethylene (cis)	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	1	156-60-5	1,2-Dichloroethylene (trans)	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	0.10	120-83-2	2,4-Dichlorophenol	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	0.10	94-82-6	4-(2,4-Dichlorophenoxy)butyric Acid (2,4-DB)	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	0.05	94-75-7	2,4-Dichlorophenoxyacetic Acid (2,4-D)	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	1	78-87-5	1,2-Dichloropropane	3.4E-01	ca*	7.4E-01	ca*	9.9E-02	ca*	1.6E-01	ca*	3.0E-02	1.0E-03	
1.0E-01	i	3.00E-02	i	1.4E-02	i	5.7E-03	i	1	542-75-6	1,3-Dichloropropene	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	0.10	616-23-9	2,3-Dichloropropanol	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	0.10	62-73-7	Dichlorvos	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	0.10	115-32-2	Dicofol	1.1E+00	ca	3.9E+00	ca	1.5E-02	ca	1.5E-01	ca		
		3.0E-02	h			5.7E-05	x	1	77-73-6	Dicyclopentadiene	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	0.10	60-57-1	Dieldrin	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	h			5.7E-03	h	0	0.10	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	h			8.6E-04	h	0	0.10	111-90-0	<b>Diethylene glycol, monomethyl ether</b>	3.7E+03	nc	3.7E+04	nc	3.1E+00	nc	2.2E+03	nc		
		4.0E-03	h			4.0E-03	r	0	0.10	617-84-5	<b>Diethylformamide</b>	2.4E+02	nc	2.5E+03	nc	1.5E+01	nc	1.5E+02	nc		
1.2E-03	i	6.0E-01	i	1.2E-03	r	6.0E-01	r	0	0.10	103-23-1	Di(2-ethylhexyl)adipate	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	0.10	84-66-2	Diethyl phthalate	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	0.10	56-53-1	Diethylstilbestrol	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	0.10	43222-48-6	Difenzoquat (Avenge)	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	0.10	35367-38-5	Diflubenzuron	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
		1.1E+01	r			1.1E+01	i	1	75-37-6	1,1-Difluoroethane					4.2E+04	nc	6.9E+04	nc			
		2.00E-02	n			2.00E-02	r		0.10	28553-12-0	Diisononyl phthalate	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc		
		8.0E-02	i			8.0E-02	r	0	0.10	1445-75-6	Diisopropyl methylphosphonate	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	0.10	55290-64-7	Dimethipin	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	0.10	60-51-5	Dimethoate	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	0.10	119-90-4	3,3'-Dimethoxybenzidine	3.5E+01	ca	1.2E+02	ca	4.8E-01	ca	4.8E+00	ca		
		5.7E-06	r			5.7E-06	x	1	124-40-3	Dimethylamine	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	0.10	121-69-7	N-N-Dimethylaniline	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	0.10	95-68-1	2,4-Dimethylaniline	6.5E-01	ca	2.3E+00	ca	9.0E-03	ca	9.0E-02	ca		
5.8E-01	h			5.8E-01	r			0	0.10	21436-96-4	2,4-Dimethylaniline hydrochloride	8.4E-01	ca	3.0E+00	ca	1.2E-02	ca	1.2E-01	ca		
9.2E+00	h			9.2E+00	r			0	0.10	119-93-7	3,3'-Dimethylbenzidine	5.3E-02	ca	1.9E-01	ca	7.3E-04	ca	7.3E-03	ca		
		1.0E-01	h			8.6E-03	i	0	0.10	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	0.10	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	0.10	105-67-9	2,4-Dimethylphenol	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc	9.0E+00	4.0E-01

Key : SFO<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca) +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION							CONTAMINANT	PRELIMINARY REMEDIAL GOALS (PRGs)				SOIL SCREENING LEVELS							
SFO <sub>i</sub> 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.		"Direct Contact Exposure Pathways"				"Migration to Ground Water"							
							Residential Soil (mg/kg)	Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)							
	6.0E-04	i	6.0E-04	r	0	0.10	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	0.10	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	0.10	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	0.10	120-61-6	Dimethyl terephthalate	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	0.10	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	h	1.0E-04	r	0	0.10	528-29-0	<b>1,2-Dinitrobenzene</b>	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	0.10	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	h	1.0E-04	r	0	0.10	100-25-4	<b>1,4-Dinitrobenzene</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	0.10	51-28-5	2,4-Dinitrophenol	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc	3.0E-01	1.0E-02	
6.8E-01	i	6.8E-01	r	0	0.10	25321-14-6	Dinitrotoluene mixture	7.2E-01	ca	2.5E+00	ca	9.9E-03	ca	9.9E-02	ca	8.0E-04	4.0E-05		
	2.0E-03	i	2.0E-03	r	0	0.10	121-14-2	2,4-Dinitrotoluene (see DNT mixture for "ca")	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc	8.0E-04	4.0E-05	
	1.0E-03	h	1.0E-03	r	0	0.10	606-20-2	2,6-Dinitrotoluene (see DNT mixture for "ca")	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc	7.0E-04	3.0E-05	
	1.0E-03	i	1.0E-03	r	0	0.10	88-85-7	Dinoseb	6.1E+01	nc	6.2E+02	nc	3.7E+00	nc	3.6E+01	nc			
	4.0E-02	h	4.0E-02	r	0	0.10	117-84-0	<b>di-n-Octyl phthalate</b>	2.4E+03	nc	2.5E+04	nc	1.5E+02	nc	1.5E+03	nc	1.0E+04	1.0E+04	
1.1E-02	i	1.1E-02	r	0	0.10	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	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	0.10	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	0.10	122-39-4	Diphenylamine	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc			
	3.00E-04	n	3.00E-04	r	0	0.10	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	7.7E-01	i	0	0.10	122-66-7	1,2-Diphenylhydrazine	6.1E-01	ca	2.2E+00	ca	8.7E-03	ca	8.4E-02	ca				
	3.0E-03	n	3.0E-03	r	0	0.10	127-63-9	<b>Diphenyl sulfone</b>	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	2439-10-3	Dodine	2.4E+02	nc	2.5E+03	nc	1.5E+01	nc	1.5E+02	nc			
	2.0E-01	n	7429-91-6	r	0	0.10	Dysprosium	1.6E+04	nc	1.0E+05	max	7.3E+03	nc						
	6.0E-03	i	6.0E-03	r	0	0.10	115-29-7	Endosulfan	3.7E+02	nc	3.7E+03	nc	2.2E+01	nc	2.2E+02	nc	1.8E+01	9.0E-01	
	2.0E-02	i	2.0E-02	r	0	0.10	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	0.10	72-20-8	Endrin	1.8E+01	nc	1.8E+02	nc	1.1E+00	nc	1.1E+01	nc	1.0E+00	5.0E-02	
9.9E-03	i	2.0E-03	h	4.2E-03	h	2.9E-04	i	1	106-89-8	Epichlorohydrin	7.6E+00	nc	2.6E+01	nc	1.0E+00	nc	2.0E+00	nc	
	5.7E-03	r	5.7E-03	i	0	0.10	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	0.10	759-94-4	EPTC (S-Ethyl dipropylthiocarbamate)	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 h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca) +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION							CONTAMINANT		PRELIMINARY REMEDIAL 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)	Industrial Soil (mg/kg)	"Direct Contact Exposure Pathways"		"Migration to Ground Water"								
										Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)							
	5.0E-03	i	5.0E-03	r	0	0.10	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	0.10	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	0.10	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	0.10	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	1		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		1	140-88-5	Ethyl acrylate	2.1E-01	ca	4.5E-01	ca	1.4E-01	ca	2.3E-01	ca				
3.85E-03	r	1.0E-01	i	3.85E-03	n	2.9E-01	i	1	100-41-4	<b>Ethylbenzene</b>	8.9E+00	ca	2.0E+01	ca	1.7E+00	ca	2.9E+00	ca	1.3E+01	7.0E-01
2.9E-03	n	4.0E-01	n	2.9E-03	r	2.9E+00	i	1	75-00-3	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	0.10	109-78-4	Ethylene cyanohydrin	1.8E+04	nc	1.0E+05	max	1.1E+03	nc	1.1E+04	nc				
	2.0E-02	h	2.0E-02	r	0	0.10	107-15-3	Ethylene diamine	1.2E+03	nc	1.2E+04	nc	7.3E+01	nc	7.3E+02	nc				
	2.0E+00	i	2.0E+00	r	0	0.10	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	0.10	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			1	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	r	0	0.10	96-45-7	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	1		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	1		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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	16984-48-8	Flouride	3.7E+03	nc	3.7E+04	nc		2.2E+03	nc				
	8.0E-02	i	8.0E-02	r	0	0.10	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	0.10	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	0.10	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	0.10	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	r	0	0.10	133-07-3	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	0.10	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	0.10	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	0.10	50-00-0	Formaldehyde	9.2E+03	nc	1.0E+05	nc	1.5E-01	ca	5.5E+03	nc			
	2.0E+00	h	2.0E+00	r	0	0.10	64-18-6	Formic Acid	1.0E+05	max	1.0E+05	max	7.3E+03	nc	7.3E+04	nc				
	3.0E+00	i	3.0E+00	r	0	0.10	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	1		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	1		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	0.10	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	0.10	98-01-1	Furfural	1.8E+02	nc	1.8E+03	nc	5.2E+01	nc	1.1E+02	nc				



Key : SFO<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca) +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION							CONTAMINANT	PRELIMINARY REMEDIAL GOALS (PRGs)				SOIL SCREENING LEVELS										
SFO <sub>i</sub> 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.		"Direct Contact Exposure Pathways"				"Migration to Ground Water"										
							Residential Soil (mg/kg)	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+01	h	5.0E+01	r	0	0.10	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	0.10	60568-05-0	Furmecycloz	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	77182-82-2	Glufosinate-ammonium	2.4E+01	nc	2.5E+02	nc	1.5E+00	nc	1.5E+01	nc						
		4.0E-04	i	2.9E-04	h	0	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	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	69806-40-2	Haloxypop-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	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	0.10	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	0.10	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	0.10	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	0.10	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	3.00E-04	n	7.8E-02	i	3.00E-04	r	0	0.10	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	0.04	319-84-6	<b>HCH (alpha)</b>	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	0.04	319-85-7	<b>HCH (beta)</b>	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	0.04	58-89-9	<b>HCH (gamma) Lindane</b>	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	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	0.10	77-47-4	<b>Hexachlorocyclopentadiene</b>	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	0.10	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	0.10	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	0.10	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	0.10	822-06-0	1,6-Hexamethylene diisocyanate	1.7E-01	nc	1.8E+00	nc	1.0E-02	nc	1.0E-01	nc					
		6.0E-02	h	5.7E-02	i	1	110-54-3	n-Hexane	1.1E+02	sat	1.1E+02	sat	2.1E+02	nc	3.5E+02	nc						
		3.3E-02	i	3.3E-02	r	0	0.10	51235-04-2	Hexazinone	2.0E+03	nc	2.0E+04	nc	1.2E+02	nc	1.2E+03	nc					
3.0E+00	i	1.7E+01	i		0	0.10	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		0.10	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		0.10	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	1	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							
		4.0E-02	h	4.0E-02	r	0	0.10	123-31-9	p-Hydroquinone	2.4E+03	nc	2.5E+04	nc	1.5E+02	nc	1.5E+03	nc					
		1.3E-02	i	1.3E-02	r	0	0.10	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	0.10	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	0.10	36734-19-7	Iprodione	2.4E+03	nc	2.5E+04	nc	1.5E+02	nc	1.5E+03	nc					
		3.0E-01	n		0	7439-89-6	Iron	2.3E+04	nc	1.0E+05	max			1.1E+04	nc							
		3.0E-01	i	3.0E-01	r	1	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	0.10	78-59-1	Isophorone	5.1E+02	ca*	1.8E+03	ca*	7.1E+00	ca	7.1E+01	ca	5.0E-01	3.0E-02	

Key : SFo,i=Cancer Slope Factor oral, inhalation RfDo,i=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca) +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION								CONTAMINANT		PRELIMINARY REMEDIAL 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)	Industrial Soil (mg/kg)	"Direct Contact Exposure Pathways"		"Migration to Ground Water"							
										Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)						
	1.5E-02	i	1.5E-02	r	0	0.10	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	0.10	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	0.10	82558-50-7	Isosabten	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc			
8.0E+00	n	3.0E-04	n	8.0E+00	r	3.0E-04	r	0	0.10	143-50-0	<b>Kepone</b>	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	0.10	77501-63-4	Lactofen	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc			
For info see: <a href="http://www.epa.gov/oerrpage/superfund/programs/lead/prods.htm#guidance">www.epa.gov/oerrpage/superfund/programs/lead/prods.htm#guidance</a>							7439-92-1	<b>Lead+++</b>	4.0E+02	nc	7.5E+02	nc							
For info see: <a href="http://www.dtsc.ca.gov/ScienceTechnology/leadspred.html">www.dtsc.ca.gov/ScienceTechnology/leadspred.html</a>								<b>Lead "CAL-Modified PRG"+++</b>	1.5E+02										
	1.0E-07	i			0	0.10	78-00-2	Lead (tetraethyl)	6.1E-03	nc	6.2E-02	nc			3.6E-03	nc			
	2.0E-03	i	2.0E-03	r	0	0.10	330-55-2	Linuron	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc			
	2.0E-02	x			0		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	0.10	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	0.10	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	0.10	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	1		123-33-1	Maleic hydrazide	1.7E+03	nc	2.4E+03	sat	1.8E+03	nc	3.0E+03	nc			
	2.0E-05	h	2.0E-05	r	0	0.10	109-77-3	Malononitrile	1.2E+00	nc	1.2E+01	nc	7.3E-02	nc	7.3E-01	nc			
	3.0E-02	h	3.0E-02	r	0	0.10	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	0.10	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	0		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	0.10	950-10-7	Mephosolan	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	0.10	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	0.10	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			0		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	0.10	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	0.10	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	0.10	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	0.10	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	1		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	0.10	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	0.10	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	0.10	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	1		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	0.10	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	0.10	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	0.10	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		0.10	99-59-2	2-Methoxy-5-nitroaniline	1.1E+01	ca	3.7E+01	ca	1.5E-01	ca	1.5E+00	ca			

Key : SFO<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca)  
 +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION						CAS No.	CONTAMINANT	PRELIMINARY REMEDIAL GOALS (PRGs)				SOIL SCREENING LEVELS								
SFO <sub>i</sub> 1/(mg/kg-d)	RfDo (mg/kg-d)	SFi 1/(mg/kg-d)	RfDi (mg/kg-d)	V O C	skin abs. soils			"Direct Contact Exposure Pathways"				"Migration to Ground Water"								
						Residential Soil (mg/kg)	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+00	h	1.0E+00	r	1	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	1	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	0	0.10	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	0	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	1	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	0	0.10	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	0.10	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	0	0.10	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	1	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	1	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
	1.7E-04	r	1.7E-04	i	0	0.10	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	2.9E-01	i	1	78-93-3	Methyl ethyl ketone	7.3E+03	nc	2.7E+04	nc	1.0E+03	nc	1.9E+03	nc					
	8.0E-02	h	2.3E-02	h	1	108-10-1	Methyl isobutyl ketone	7.9E+02	nc	2.8E+03	nc	8.3E+01	nc	1.6E+02	nc					
	5.7E-04	r	5.7E-04	n	0	0.10	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	1	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	0.10	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	0.10	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	0.10	95-48-7	2-Methylphenol	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc	1.5E+01	8.0E-01		
	5.0E-02	i	5.0E-02	r	0	0.10	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	0.10	106-44-5	4-Methylphenol	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc				
	2.0E-02	n	2.0E-02	r	0	0.10	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	1	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	1	98-83-9	Methyl styrene (alpha)	6.8E+02	sat	6.8E+02	sat	2.6E+02	nc	4.3E+02	nc					
3.3E-03	n	8.6E-01	r	3.5E-04	n	8.6E-01	i	1	1634-04-4	<b>Methyl tertbutyl ether (MTBE)</b>	6.2E+01	ca*	1.6E+02	ca	1.9E+01	ca	1.3E+01	ca		
1.8E-03		1.8E-03			1					<b>"CAL-Modified PRG"</b>	1.7E+01	ca	3.6E+01	ca	3.7E+00	ca	6.2E+00	ca		
	1.5E-01	i	1.5E-01	r	0	0.10	51218-45-2	Metolacolor (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	0.10	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	r	0	0.10	2385-85-5	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	0.10	2212-67-1	Molinatate	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc				
	5.0E-03	i			0	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	0.10	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	0.10	300-76-5	Naled	1.2E+02	nc	1.2E+03	nc	7.3E+00	nc	7.3E+01	nc				

Key : SFO<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca)  
 +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION				CONTAMINANT			PRELIMINARY REMEDIAL 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.	"Direct Contact Exposure Pathways"				"Migration to Ground Water"							
							Residential Soil (mg/kg)	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-01	i	1.0E-01	r	0	0.10	15299-99-7	Napropamide	6.1E+03	nc	6.2E+04	nc	3.7E+02	nc	3.6E+03	nc		
	2.0E-02	i			0		7440-02-0	Nickel (soluble salts)	1.6E+03	nc	2.0E+04	nc			7.3E+02	nc	1.3E+02	7.0E+00
		8.4E-01	i		0			Nickel refinery dust					8.0E-03	ca				
		1.7E+00	i		0		12035-72-2	Nickel subsulfide			1.1E+04	ca	4.0E-03	ca				
							14797-55-8	Nitrate+++					1.0E+04	nc				
							14797-65-0	Nitrite+++					1.0E+03	nc				
	2.86E-05	r	2.86E-05	h	0	0.10	88-74-4	2-Nitroaniline	1.7E+00	nc	1.8E+01	nc	1.0E-01	nc	1.0E+00	nc		
	5.0E-04	i	5.7E-04	h	1		98-95-3	Nitrobenzene	2.0E+01	nc	1.0E+02	nc	2.1E+00	nc	3.4E+00	nc	1.0E-01	7.0E-03
	7.0E-02	h	7.0E-02	r	0	0.10	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	0.10	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	0.10	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	0.10	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	i	1	79-46-9	2-Nitropropane			7.2E-04	ca	1.2E-03	ca	
	5.4E+00	i	5.6E+00	i			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	0.10	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	0.10	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	4.9E+01	i	0	0.10	62-75-9	N-Nitrosodimethylamine	9.5E-03	ca	3.4E-02	ca	1.4E-04	ca	1.3E-03	ca		
	4.9E-03	i	4.9E-03	r	0	0.10	86-30-6	N-Nitrosodiphenylamine	9.9E+01	ca	3.5E+02	ca	1.4E+00	ca	1.4E+01	ca	1.0E+00	6.0E-02
	7.0E+00	i	7.0E+00	r	0	0.10	621-64-7	N-Nitroso di-n-propylamine	6.9E-02	ca	2.5E-01	ca	9.6E-04	ca	9.6E-03	ca	5.0E-05	2.0E-06
	2.2E+01	i	2.2E+01	r	0	0.10	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	0.10	930-55-2	N-Nitrosopyrrolidine	2.3E-01	ca	8.2E-01	ca	3.1E-03	ca	3.2E-02	ca		
	1.0E-02	h	1.0E-02	r	1		99-08-1	m-Nitrotoluene	3.7E+02	nc	1.0E+03	sat	3.7E+01	nc	6.1E+01	nc		
	1.0E-02	h	1.0E-02	r	1		99-08-1	o-Nitrotoluene	3.7E+02	nc	1.0E+03	sat	3.7E+01	nc	6.1E+01	nc		
	1.0E-02	h	1.0E-02	r	1		99-99-0	p-Nitrotoluene	3.7E+02	nc	1.0E+03	sat	3.7E+01	nc	6.1E+01	nc		
	4.0E-02	i	4.0E-02	r	0	0.10	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	0.10	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	0.10	32536-52-0	Octabromodiphenyl ether	1.8E+02	nc	1.8E+03	nc	1.1E+01	nc	1.1E+02	nc		
	5.0E-02	i	5.0E-02	r	0	0.10	2691-41-0	Octahydro-1357-tetranitro-1357- tetrazocine (HMX)	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc		
	2.0E-03	h	2.0E-03	r	0	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	76738-62-0	Paclitaxel	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	0.10	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	0.10	56-38-2	Parathion	3.7E+02	nc	3.7E+03	nc	2.2E+01	nc	2.2E+02	nc		

Key : SFo<sub>i</sub>=Cancer Slope Factor oral, inhalation RfDo<sub>i</sub>=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca)  
 +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION							CONTAMINANT	PRELIMINARY REMEDIAL 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.		"Direct Contact Exposure Pathways"				"Migration to Ground Water"							
							Residential Soil (mg/kg)	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-02	h		5.0E-02	r	0	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0	0.10	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	0	0.25	87-86-5	Pentachlorophenol	3.0E+00	ca	9.0E+00	ca	5.6E-02	ca	5.6E-01	ca	3.0E-02	1.0E-03
	1.00E-04	x				0		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	0.10	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	0.10	13684-63-4	Phenmedipham	1.5E+04	nc	1.0E+05	max	9.1E+02	nc	9.1E+03	nc		
	6.0E-01	i		6.0E-01	r	0	0.10	108-95-2	Phenol	3.7E+04	nc	1.0E+05	max	2.2E+03	nc	2.2E+04	nc	1.0E+02	5.0E+00
	2.0E-03	n		2.0E-03	r	0	0.10	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	0.10	108-45-2	m-Phenylenediamine	3.7E+02	nc	3.7E+03	nc	2.2E+01	nc	2.2E+02	nc		
	1.9E-01	h		1.9E-01	r	0	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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				0		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	0.10	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	0.10	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	0.10	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	0.10	29232-93-7	Pirimiphos-methyl	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc		
8.9E+00	h	7.0E-06	h	8.9E+00	r	0	0.10		Polybrominated biphenyls	5.5E-02	ca**	1.9E-01	ca*	7.6E-04	ca*	7.6E-03	ca*		
2.0E+00	i		2.0E+00	i		0	0.14	1336-36-3	Polychlorinated biphenyls (PCBs)	2.2E-01	ca	7.4E-01	ca	3.4E-03	ca	3.4E-02	ca		
7.0E-02	i	7.0E-05	i	7.0E-02	i	0	0.14	12674-11-2	Aroclor 1016	3.9E+00	nc	2.1E+01	ca**	9.6E-02	ca**	9.6E-01	ca**		
2.0E+00	i		2.0E+00	i		0	0.14	11104-28-2	Aroclor 1221	2.2E-01	ca	7.4E-01	ca	3.4E-03	ca	3.4E-02	ca		
2.0E+00	i		2.0E+00	i		0	0.14	11141-16-5	Aroclor 1232	2.2E-01	ca	7.4E-01	ca	3.4E-03	ca	3.4E-02	ca		
2.0E+00	i		2.0E+00	i		0	0.14	53469-21-9	Aroclor 1242	2.2E-01	ca	7.4E-01	ca	3.4E-03	ca	3.4E-02	ca		
2.0E+00	i		2.0E+00	i		0	0.14	12672-29-6	Aroclor 1248	2.2E-01	ca	7.4E-01	ca	3.4E-03	ca	3.4E-02	ca		
2.0E+00	i	2.0E-05	i	2.0E+00	i	0	0.14	11097-69-1	Aroclor 1254	2.2E-01	ca**	7.4E-01	ca*	3.4E-03	ca*	3.4E-02	ca*		
2.0E+00	i		2.0E+00	i		0	0.14	11096-82-5	Aroclor 1260	2.2E-01	ca	7.4E-01	ca	3.4E-03	ca	3.4E-02	ca		

Key : SFo,i=Cancer Slope Factor oral, inhalation RfDo,i=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca)  
 +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION					CONTAMINANT		PRELIMINARY REMEDIAL 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.	"Direct Contact Exposure Pathways"				"Migration to Ground Water"								
							Residential Soil (mg/kg)	Industrial Soil (mg/kg)	Ambient Air (ug/m^3)	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)							
4.5E+00	n	4.5E+00	r			0.10 0.13	61788-33-8	<b>Polychlorinated terphenyls</b>		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	1		83-32-9	3.7E+03	nc	2.9E+04	nc	2.2E+02	nc	3.7E+02	nc	5.7E+02	2.9E+01		
	3.0E-01	i	3.0E-01	r	1		120-12-7	2.2E+04	nc	1.0E+05	max	1.1E+03	nc	1.8E+03	nc	1.2E+04	5.9E+02		
7.3E-01	n	7.3E-01	r		0	0.13	56-55-3	6.2E-01	ca	2.1E+00	ca	9.2E-03	ca	9.2E-02	ca	2.0E+00	8.0E-02		
7.3E-01	n	7.3E-01	r		0	0.13	205-99-2	6.2E-01	ca	2.1E+00	ca	9.2E-03	ca	9.2E-02	ca	5.0E+00	2.0E-01		
7.3E-02	n	7.3E-02	r		0	0.13	207-08-9	6.2E+00	ca	2.1E+01	ca	9.2E-02	ca	9.2E-01	ca	4.9E+01	2.0E+00		
1.2E+00		3.9E-01				0.13	207-08-9	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	0.13	50-32-8	6.2E-02	ca	2.1E-01	ca	9.2E-04	ca	9.2E-03	ca	8.0E+00	4.0E-01		
7.3E-03	n	7.3E-03	r		0	0.13	218-01-9	6.2E+01	ca	2.1E+02	ca	9.2E-01	ca	9.2E+00	ca	1.6E+02	8.0E+00		
1.2E-01		3.9E-02				0.13		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	0.13	53-70-3	6.2E-02	ca	2.1E-01	ca	9.2E-04	ca	9.2E-03	ca	2.0E+00	8.0E-02		
	4.0E-02	i	4.0E-02	r	0	0.13	206-44-0	2.3E+03	nc	2.2E+04	nc	1.5E+02	nc	1.5E+03	nc	4.3E+03	2.1E+02		
	4.0E-02	i	4.0E-02	r	1		86-73-7	2.7E+03	nc	2.6E+04	nc	1.5E+02	nc	2.4E+02	nc	5.6E+02	2.8E+01		
7.3E-01	n	7.3E-01	r		0	0.13	193-39-5	6.2E-01	ca	2.1E+00	ca	9.2E-03	ca	9.2E-02	ca	1.4E+01	7.0E-01		
	2.0E-02	i	8.6E-04	i	1		91-20-3	5.6E+01	nc	1.9E+02	nc	3.1E+00	nc	6.2E+00	nc	8.4E+01	4.0E+00		
	3.0E-02	i	3.0E-02	r	1		129-00-0	2.3E+03	nc	2.9E+04	nc	1.1E+02	nc	1.8E+02	nc	4.2E+03	2.1E+02		
1.5E-01	i	9.0E-03	i	1.5E-01	r	0	0.10	67747-09-5	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	0.10	26399-36-0	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	0.10	1610-18-0	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	0.10	7287-19-6	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	0.10	23950-58-5	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	0.10	1918-16-7	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	0.10	709-98-8	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	0.10	2312-35-8	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	0.10	107-19-7	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	0.10	139-40-2	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	0.10	122-42-9	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	0.10	60207-90-1	7.9E+02	nc	8.0E+03	nc	4.7E+01	nc	4.7E+02	nc				
	4.00E-02	n	4.00E-02	r	1		103-65-1	2.4E+02	sat	2.4E+02	sat	1.5E+02	nc	2.4E+02	nc				
	5.0E-01	h	8.6E-04	h	0	0.10	57-55-6	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	0.10	52125-53-8	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	0.10	107-98-2	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	1	75-56-9	1.9E+00	ca*	6.6E+00	ca*	5.2E-01	ca*	2.2E-01	ca				

Key : SFo,i=Cancer Slope Factor oral, inhalation RfDo,i=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca)  
 +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION							CONTAMINANT	PRELIMINARY REMEDIAL 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.		"Direct Contact Exposure Pathways"				"Migration to Ground Water"							
							Residential Soil (mg/kg)	Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)							
	2.5E-01	i		2.5E-01	r	0	0.10	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	0.10	51630-58-1	Pydrin	1.5E+03	nc	1.5E+04	nc	9.1E+01	nc	9.1E+02	nc		
	1.0E-03	i		1.0E-03	r	0	0.10	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	0.10	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		0	0.10	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	0.10	121-82-4		1.6E+01	ca	6.1E-02	ca	6.1E-01	ca		
	3.0E-02	i		3.0E-02	r	0	0.10	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	0.10	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	0.10	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	0.10	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	0.10	7783-00-8	Selenious Acid	3.1E+02	nc	3.1E+03	nc			1.8E+02	nc		
	5.0E-03	i				0		7782-49-2	Selenium	3.9E+02	nc	5.1E+03	nc			1.8E+02	nc		
	5.0E-03	h				0	0.10	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	0.10	74051-80-2	Sethoxydim	5.5E+03	nc	5.5E+04	nc	3.3E+02	nc	3.3E+03	nc		
	5.0E-03	i				0		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	2.0E-03	r	0	0.10	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	0.10	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	0.10	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	0.10	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				0		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	0.10	57-24-9	Strychnine	1.8E+01	nc	1.8E+02	nc	1.1E+00	nc	1.1E+01	nc		
	2.0E-01	i		2.9E-01	i	1		100-42-5	Styrene	1.7E+03	sat	1.7E+03	sat	1.1E+03	nc	1.6E+03	nc		
	1.00E-03	n		1.00E-03	r			80-07-9	1,1'-Sulfonylbis (4-chlorobenzene)	7.8E+01	nc	1.0E+03	nc	3.7E+00	nc	3.6E+01	nc		
	2.5E-02	i		2.5E-02	r	0	0.10	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	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	1	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	6.00E-02	n	2.0E-01	i	6.00E-02	r	1	79-34-5	1,1,2,2-Tetrachloroethane	4.1E-01	ca	9.3E-01	ca	3.3E-02	ca	5.5E-02	ca	
5.2E-02	n	1.0E-02	i	1.00E-02	n	1.7E-01	n	1	127-18-4	Tetrachloroethylene (PCE)	1.5E+00	ca*	3.4E+00	ca*	6.7E-01	ca	6.6E-01	ca	
	3.0E-02	i		3.0E-02	r	0	0.10	58-90-2	2,3,4,6-Tetrachlorophenol	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc		

Key : SFo,i=Cancer Slope Factor oral, inhalation RfDo,i=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca) +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION										CONTAMINANT		PRELIMINARY REMEDIAL 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.						"Direct Contact Exposure Pathways"				"Migration to Ground Water"					
							Residential Soil (mg/kg)	Industrial Soil (mg/kg)	Ambient Air (ug/m <sup>3</sup> )	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)									
2.0E+01	h	2.0E+01	r	0	0.10	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	0	0.10	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	0.10	3689-24-5	Tetraethyldithiopyrophosphate	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	1	109-99-9	<b>Tetrahydrofuran</b>	9.4E+00	ca	2.1E+01	ca	9.9E-01	ca	1.6E+00	ca					
		6.6E-05	i			0	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	0.10	28249-77-6	Thiobencarb	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc				
		5.0E-02	n	5.0E-02	r	0	0.10	N/A	<b>Thiocyanate</b>	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	0.10	39196-18-4	Thiofanox	1.8E+01	nc	1.8E+02	nc	1.1E+00	nc	1.1E+01	nc				
		8.0E-02	i	8.0E-02	r	0	0.10	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	0.10	137-26-8	Thiram	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc				
		6.0E-01	h			0			Tin (inorganic, see tributyltin oxide for organic tin)	4.7E+04	nc	1.0E+05	max			2.2E+04	nc				
		2.0E-01	i	1.1E-01	i	1	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	0.10	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	0.10	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	0.10	823-40-5	Toluene-2,6-diamine	1.2E+04	nc	1.0E+05	max	7.3E+02	nc	7.3E+03	nc				
2E-01	i	2E-01	r		0	0.10	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	0.10	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	0.10	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	0.10	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	0.10	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	0.10	615-54-3	1,2,4-Tribromobenzene	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc				
		3.0E-04	i		0	0.10	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	0.10	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	0.10	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	5.7E-02	h	1	120-82-1	1,2,4-Trichlorobenzene	6.5E+02	nc	3.0E+03	sat	2.1E+02	nc	1.9E+02	nc	5.0E+00	3.0E-01			
		2.8E-01	n	6.3E-01	n	1	71-55-6	<b>1,1,1-Trichloroethane</b>	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	1	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	
4.00E-01	n	3.00E-04	n	4.00E-01	n	1.00E-02	n	1	79-01-6	<b>Trichloroethylene (TCE)</b>	5.3E-02	ca	1.1E-01	ca	1.7E-02	ca	2.8E-02	ca	6.0E-02	3.0E-03	
		3.0E-01	i	2.0E-01	h	1	75-69-4	Trichlorofluoromethane	3.9E+02	nc	2.0E+03	sat	7.3E+02	nc	1.3E+03	nc					
		1.0E-01	i	1.0E-01	r	0	0.10	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		
1.1E-02	i	1.0E-04	n	1.1E-02	i	1.0E-04	r	0	0.10	88-06-2	<b>2,4,6-Trichlorophenol</b>	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		7.0E-02				0.10	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	0.10	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	0.10	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	1	598-77-6	1,1,2-Trichloropropane	1.5E+01	nc	5.1E+01	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	1	96-18-4	<b>1,2,3-Trichloropropane</b>	5.0E-03	ca	1.1E-02	ca	3.4E-03	ca	5.6E-03	ca			



Key : SFo,i=Cancer Slope Factor oral, inhalation RfDo,i=Reference Dose oral, inhalation i=IRIS h=HEAST n=NCEA x=Withdrawn o=Other EPA Source r=Route-extrapolation ca=Cancer PRG nc=Noncancer PRG ca\* (where: nc < 100X ca) ca\*\* (where: nc < 10X ca) +++=Non-Standard Method Applied (See Section 2.3 of the "Region 9 PRGs Table User's Guide") sat=Soil Saturation (See Section 4.5) max=Ceiling limit (See Section 2.1) DAF=Dilution Attenuation Factor (See Section 2.5) CAS=Chemical Abstract Services

TOXICITY INFORMATION						CONTAMINANT		PRELIMINARY REMEDIAL 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.	"Direct Contact Exposure Pathways"								"Migration to Ground Water"					
							Residential Soil (mg/kg)	Industrial Soil (mg/kg)	Ambient Air (ug/m^3)	Tap Water (ug/l)	DAF 20 (mg/kg)	DAF 1 (mg/kg)								
	5.0E-03	h	5.0E-03	r	1	96-19-5	1,2,3-Trichloropropene	1.2E+01	nc	3.8E+01	nc	1.8E+01	nc	3.0E+01	nc					
	3.0E-03	i	3.0E-03	r	0	0.10	58138-08-2	Tridiphane	1.8E+02	nc	1.8E+03	nc	1.1E+01	nc	1.1E+02	nc				
	2.0E-03	r	2.0E-03	i	1		121-44-8	Triethylamine	2.3E+01	nc	8.6E+01	nc	7.3E+00	nc	1.2E+01	nc				
7.7E-03	i	7.5E-03	i	7.7E-03	r	0	0.10	1582-09-8	Trifluralin	6.3E+01	ca**	2.2E+02	ca*	8.7E-01	ca*	8.7E+00	ca*			
	1.400E-04	r	1.400E-04	n		0.10	552-30-7	Trimellitic Anhydride (TMAN)	8.6E+00	nc	8.6E+01	nc	5.1E-01	nc	5.1E+00					
	5.0E-02	n	1.7E-03	n	1		95-63-6	1,2,4-Trimethylbenzene	5.2E+01	nc	1.7E+02	nc	6.2E+00	nc	1.2E+01	nc				
3.7E-02	h	3.7E-02	r		0	0.10	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	0.10	512-56-1	Trimethyl phosphate	1.3E+01	ca	4.7E+01	ca	1.8E-01	ca	1.8E+00	ca				
	3.0E-02	i	3.0E-02	r	0	0.10	99-35-4	1,3,5-Trinitrobenzene	1.8E+03	nc	1.8E+04	nc	1.1E+02	nc	1.1E+03	nc				
3E-02	i	1.0E-02	h	1.0E-02	r	0	0.10	479-45-8	Trinitrophenylmethyl nitramine	6.1E+02	nc	6.2E+03	nc	3.7E+01	nc	3.6E+02	nc			
	5.0E-04	i	3E-02	r	0	0.10	118-96-7	2,4,6-Trinitrotoluene	1.6E+01	ca**	5.7E+01	ca**	2.2E-01	ca**	2.2E+00	ca**				
	5.00E-03	n	5.00E-03	r	0	0.10	791-28-6	Triphenylphosphine oxide	3.1E+02	nc	3.1E+03	nc	1.8E+01	nc	1.8E+02	nc				
3.2E-03	n	1.1E-01	n	3.2E-03	r	0	0.10	115-96-8	Tris(2-chloroethyl) phosphate	1.5E+02	ca*	5.4E+02	ca	2.1E+00	ca	2.1E+01	ca			
	2.00E-04	n					7440-61-0	Uranium (chemical toxicity only)	1.6E+01	nc	2.0E+02	nc			7.3E+00	nc				
	7.0E-03	h			0		7440-62-2	Vanadium and compounds	5.5E+02	nc	7.2E+03	nc			2.6E+02	nc	6.0E+03	3.0E+02		
	1.0E-03	i	1.0E-03	r	0	0.10	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	0.10	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	1		108-05-4	Vinyl acetate	4.3E+02	nc	1.4E+03	nc	2.1E+02	nc	4.1E+02	nc	1.7E+02	8.0E+00		
1.1E-01	r	8.6E-04	r	1.1E-01	h	8.6E-04	i	1	593-60-2	Vinyl bromide (bromoethene)	1.9E-01	ca*	4.2E-01	ca*	6.1E-02	ca*	1.0E-01	ca*		
1.5E+00	i	3.00E-03	i	3.1E-02	i	2.86E-02	i	1	75-01-4	Vinyl chloride (child/adult)+++	7.9E-02	ca			1.1E-01	ca	2.0E-02	ca	1.0E-02	7.0E-04
7.5E-01	i	3.00E-03	i	1.6E-02	i	2.86E-02	i	1	75-01-4	Vinyl chloride (adult)			7.5E-01	ca						
	3.0E-04	i	3.0E-04	r	0	0.10	81-81-2	Warfarin	1.8E+01	nc	1.8E+02	nc	1.1E+00	nc	1.1E+01	nc				
	7.0E-01	i	2.9E-02	i	1	0.10	1330-20-7	Xylenes	2.7E+02	nc	4.2E+02	sat	1.1E+02	nc	2.1E+02	nc	2.1E+02	1.0E+01		
	3.0E-01	i			0		7440-66-6	Zinc	2.3E+04	nc	1.0E+05	max			1.1E+04	nc	1.2E+04	6.2E+02		
	3.0E-04	i			0		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	0.10	12122-67-7	Zineb	3.1E+03	nc	3.1E+04	nc	1.8E+02	nc	1.8E+03	nc				