

## Appendix A: Hot Spots Unit Risk and Cancer Potency Values

Updated May 2019

Chemical	Chemical Abstract Service (CAS) Number	Source	Unit Risk ( $\mu\text{g}/\text{m}^3$ ) <sup>-1</sup>	Slope Factor ( $\text{mg}/\text{kg}\cdot\text{day}$ ) <sup>-1</sup>	US EPA Class <sup>C</sup>	IARC Class <sup>C</sup>
Acetaldehyde	75-07-0	TAC	2.7 E-6	1.0 E-2	B2	2B
Acetamide	60-35-5	P65-E	2.0 E-5	7.0 E-2	NC	2B
Acrylamide	79-06-1	IRIS	1.3 E-3	4.5 E+0	B2	2A
Acrylonitrile	107-13-1	P65-S	2.9 E-4	1.0 E+0	B1	2A
Allyl chloride	107-05-1	P65-S	6.0 E-6	2.1 E-2	C	3
2-Aminoanthraquinone	117-79-3	P65-E	9.4 E-6	3.3 E-2	NC	3
Aniline	62-53-3	IRIS	1.6 E-6	5.7 E-3	B2	3
Arsenic (inorganic)	7440-38-2	(inhalation) TAC	3.3 E-3	1.2 E+1	A	1
(oral)		IRIS		1.5 E+0		
Asbestos	1332-21-4	TAC	6.3 E-2 1.9 E-4 <sup>#</sup>	2.2 E+2	A	1
Benz[ <i>a</i> ]anthracene <sup>BaP</sup>	56-55-3	(inhalation) TAC	1.1 E-4	3.9 E-1	B2	2A
(oral)				1.2 E+0		
Benzene	71-43-2	TAC	2.9 E-5	1.0 E-1	A	1
Benzidine	92-87-5	P65-S	1.4 E-1	5.0 E+2	A	1
Benzo[ <i>a</i> ]pyrene	50-32-8	(inhalation) TAC	1.1 E-3	3.9 E+0	B2	2A
(oral)				1.2 E+1		
Benzo[ <i>b</i> ]fluoranthrene <sup>BaP</sup>	205-99-2	(inhalation) TAC	1.1 E-4	3.9 E-1	B2	2B
(oral)				1.2 E+0		
Benzo[ <i>j</i> ]fluoranthrene <sup>BaP</sup>	205-82-3	(inhalation) TAC	1.1 E-4	3.9 E-1	NC	2B
(oral)				1.2 E+0		
Benzo[ <i>k</i> ]fluoranthrene <sup>BaP</sup>	207-08-9	(inhalation) TAC	1.1 E-4	3.9 E-1	B2	2B
(oral)				1.2 E+0		
Benzyl chloride	100-44-7	IRIS	4.9 E-5	1.7 E-1	B2	2B
Beryllium	7440-41-7	IRIS	2.4 E-3	8.4 E+0	B2	1
Bis(2-chloroethyl) ether	111-44-4	P65-S	7.1 E-4	2.5 E+0	B2	3
Bis(chloromethyl)ether	542-88-1	P65-S	1.3 E-2	4.6 E+1	A	1
1,3-Butadiene	106-99-0	TAC	1.7 E-4	6.0 E-1	B2	2A
Cadmium (and compounds)	7440-43-9	TAC	4.2 E-3	1.5 E+1	B1	1
Carbon tetrachloride	56-23-5	TAC	4.2 E-5	1.5 E-1	B2	2B
Chlorinated dibenzo- <i>p</i> -dioxins <sup>A</sup>	1746-01-6	TAC			B2	2B
2,3,7,8-Tetrachlorodibenzo- <i>p</i> -dioxin			3.8 E+1	1.3 E+5		
1,2,3,7,8-Pentachlorodibenzo- <i>p</i> -dioxin			3.8 E+1	1.3 E+5		
1,2,3,4,7,8-Hexachlorodibenzo- <i>p</i> -dioxin			3.8 E+0	1.3 E+4		
1,2,3,6,7,8-Hexachlorodibenzo- <i>p</i> -dioxin			3.8 E+0	1.3 E+4		

### Footnotes

- A see Appendix C
- BaP see benzo[*a*]pyrene TAC document
- C see Appendix E
- D Listed by ARB as “Particulate Matter from Diesel-Fueled Engines”; Scientific Review Panel unit risk “reasonable estimate” = 3.0 E-4 ( $\mu\text{g}/\text{m}^3$ )<sup>-1</sup>. Range of unit risks in TAC document was 1.3 E-4 – 2.4 E-3 ( $\mu\text{g}/\text{m}^3$ )<sup>-1</sup>.
- N1 Nickel refinery dust and nickel subsulfide are in Class A; nickel carbonyl is in Class B2
- N2 Nickel compounds are in Class 1; metallic nickel is in Class 2B
- NA not available
- NC not classified
- # [100 PCM fibers/m<sup>3</sup>]-1 ; see Appendix F
- \* can be calculated using PEF factors contained in the benzo[*a*]pyrene TAC document
- P See PCB summary for risk categorization and TEF factors (Appendix C)

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1,2,3,7,8,9-Hexachlorodibenzo- <i>p</i> -dioxin	5120-73-19	TAC	3.8 E+0	1.3 E+4	B2	NC
1,2,3,4,6,7,8-Heptachlorodibenzo- <i>p</i> -dioxin			3.8 E-1	1.3 E+3		
1,2,3,4,6,7,8,9-Octachlorodibenzo- <i>p</i> -dioxin			1.1 E-2	3.9 E+1		
Chlorinated dibenzofurans <sup>A</sup>						
2,3,7,8-Tetrachlorodibenzofuran			3.8 E+0	1.3 E+4		
1,2,3,7,8-Pentachlorodibenzofuran			1.1 E+0	3.9 E+3		
2,3,4,7,8-Pentachlorodibenzofuran			1.1 E+1	3.9 E+4		
1,2,3,4,7,8-Hexachlorodibenzofuran			3.8 E+0	1.3 E+4		
1,2,3,6,7,8-Hexachlorodibenzofuran			3.8 E+0	1.3 E+4		
1,2,3,7,8,9-Hexachlorodibenzofuran			3.8 E+0	1.3 E+4		
2,3,4,6,7,8-Hexachlorodibenzofuran			3.8 E+0	1.3 E+4		
1,2,3,4,6,7,8-Heptachlorodibenzofuran			3.8 E-1	1.3 E+3		
1,2,3,4,7,8,9-Heptachlorodibenzofuran			3.8 E-1	1.3 E+3		
1,2,3,4,6,7,8,9-Octachlorodibenzofuran			1.1 E-2	3.9 E+1		
Chlorinated paraffins	108171-26-2	P65-E	2.5 E-5	8.9 E-2	NC	2B
Chloroform	67-66-3	TAC	5.3 E-6	1.9 E-2	B2	2B
4-Chloro- <i>o</i> -phenylenediamine	95-83-0	P65-E	4.6 E-6	1.6 E-2	NC	2B
<i>p</i> -Chloro- <i>o</i> -toluidine	95-69-2	P65-E	7.7 E-5	2.7 E-1	NC	2A
Chromium (hexavalent)	18540-29-9	TAC	1.5 E-1	5.1 E+2	A	1
(inhalation)						
(oral)		P65-S		4.2 E-1		
Chrysene <sup>BaP</sup>	218-01-9	TAC	1.1 E-5	3.9 E-2	B2	3
(inhalation)						
(oral)				1.2 E-1		
Creosote	8001-58-9	HS	*	*	B1	2A
<i>p</i> -Cresidine	120-71-8	P65-E	4.3 E-5	1.5 E-1	NC	2B
Cupferron	135-20-6	P65-E	6.3 E-5	2.2 E-1	NC	NC
2,4-Diaminoanisole	615-05-4	P65-E	6.6 E-6	2.3 E-2	NC	2B
2,4-Diaminotoluene	95-80-7	P65-E	1.1 E-3	4.0 E+0	NC	2B
Dibenz[ <i>a,h</i> ]acridine <sup>BaP</sup>	226-36-8	TAC	1.1 E-4	3.9 E-1	NC	2B
(inhalation)						
(oral)				1.2 E+0		
Dibenz[ <i>a,j</i> ]acridine <sup>BaP</sup>	224-42-0	TAC	1.1 E-4	3.9 E-1	NC	2B
(inhalation)						
(oral)				1.2 E+0		
Dibenz[ <i>a,h</i> ]anthracene <sup>BaP</sup>	53-70-3	P65-E	1.2 E-3	4.1 E+0	B2	2A
Dibenzo[ <i>a,e</i> ]pyrene <sup>BaP</sup>	192-65-4	TAC	1.1 E-3	3.9 E+0	NC	2B
(inhalation)						
(oral)				1.2 E+1		

### Footnotes

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 N1 Nickel refinery dust and nickel subsulfide are in Class A; nickel carbonyl is in Class B2  
 N2 Nickel compounds are in Class 1; metallic nickel is in Class 2B  
 NA not available  
 NC not classified  
 # [100 PCM fibers/m<sup>3</sup>]-1 ; see Appendix F  
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Dibenzo[ <i>a,h</i> ]pyrene <sup>BaP</sup>	(inhalation)	189-64-0	TAC	1.1 E-2	3.9 E+1	NC	2B
	(oral)				1.2 E+2		
Dibenzo[ <i>a,i</i> ]pyrene <sup>BaP</sup>	(inhalation)	189-55-9	TAC	1.1 E-2	3.9 E+1	NC	2B
	(oral)				1.2 E+2		
Dibenzo[ <i>a,l</i> ]pyrene <sup>BaP</sup>	(inhalation)	191-30-0	TAC	1.1 E-2	3.9 E+1	NC	2B
	(oral)				1.2 E+2		
7H-Dibenzo[ <i>c,g</i> ]carbazole <sup>BaP</sup>	(inhalation)	194-59-2	TAC	1.1 E-3	3.9 E+0	NC	2B
	(oral)				1.2 E+1		
1,2-Dibromo-3-chloropropane		96-12-8	P65-S	2.0 E-3	7.0 E+0	NC	2B
1,4-Dichlorobenzene		106-46-7	P65-S	1.1 E-5	4.0 E-2	NC	2B
3,3'-Dichlorobenzidine		91-94-1	P65-S	3.4 E-4	1.2 E+0	B2	2B
1,1-Dichloroethane		75-34-3	P65-E	1.6 E-6	5.7 E-3	C	NC
Diesel exhaust		NA	TAC	3.0 E-4 <sup>D</sup>	1.1 E+0	NC	2A
Diethylhexylphthalate		117-81-7	PETB	2.4 E-6	8.4 E-3	B2	2B
<i>p</i> -Dimethylaminoazobenzene		60-11-7	P65-E	1.3 E-3	4.6 E+0	NC	2B
7,12-Dimethylbenz[ <i>a</i> ]anthracene <sup>BaP</sup>		57-97-6	P65-E	7.1 E-2	2.5 E+2	NC	NC
1,6-Dinitropyrene <sup>BaP</sup>	(inhalation)	42397-64-8	TAC	1.1 E-2	3.9 E+1	NC	2B
	(oral)				1.2 E+2		
1,8-Dinitropyrene <sup>BaP</sup>	(inhalation)	42397-65-9	TAC	1.1 E-3	3.9 E+0	NC	2B
	(oral)				1.2 E+1		
2,4-Dinitrotoluene		121-14-2	P65-S	8.9 E-5	3.1 E-1	NC	2B
1,4-Dioxane		123-91-1	P65-S	7.7 E-6	2.7 E-2	B2	2B
Epichlorohydrin		106-89-8	P65-S	2.3 E-5	8.0 E-2	B2	2A
Ethylbenzene	(inhalation)	100-41-4	HS	2.5 E-6	8.7 E-3	D	2B
	(oral)				1.1 E-2		
Ethylene dibromide		106-93-4	TAC	7.1 E-5	2.5 E-1	B2	2A
Ethylene dichloride		107-06-2	TAC	2.1 E-5	7.2 E-2	B2	2B
Ethylene oxide		75-21-8	TAC	8.8 E-5	3.1 E-1	NC	1
Ethylene thiourea		96-45-7	P65-E	1.3 E-5	4.5 E-2	UR	2B
Formaldehyde		50-00-0	TAC	6.0 E-6	2.1 E-2	B1	2A
Hexachlorobenzene		118-74-1	P65-S	5.1 E-4	1.8 E+0	B2	2B
Hexachlorocyclohexanes (technical grade)		608-73-1	P65-S	1.1 E-3	4.0 E+0	B2	2B
Hydrazine	(inhalation)	302-01-2	IRIS	4.9 E-3	1.7 E+1	B2	2B
	(oral)				3.0 E+0		

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Indeno[1,2,3- <i>cd</i> ]pyrene <sup>BaP</sup>	(inhalation)	193-39-5	TAC	1.1 E-4	3.9 E-1	B2	2B
	(oral)				1.2 E+0		
Lead and lead compounds	(inhalation)	7439-92-1	TAC	1.2 E-5	4.2 E-2	B2	2B
	(oral)				8.5 E-3		
Lindane		58-89-9	P65-S	3.1 E-4	1.1 E+0	NC	2B
Methyl <i>tert</i> -butyl ether (MTBE)		1634-04-4	HS	2.6 E-7	1.8 E-3	NC	3
3-Methylcholanthrene <sup>BaP</sup>		56-49-5	P65-E	6.3 E-3	2.2 E+1	NC	NC
5-Methylchrysene <sup>BaP</sup>	(inhalation)	3697-24-3	TAC	1.1 E-3	3.9 E+0	NC	2B
	(oral)				1.2 E+1		
4, 4'-Methylene bis(2-chloroaniline) (MOCA)		101-14-4	P65-E	4.3 E-4	1.5 E+0	NC	2A
Methylene chloride		75-09-2	TAC	1.0 E-6	3.5 E-3	B2	2B
4,4'-Methylenedianiline		101-77-9	P65-E	4.6 E-4	1.6 E+0	NC	2B
Michler's ketone		90-94-8	P65-E	2.5 E-4	8.6 E-1	NC	NC
Naphthalene		91-20-3	HS	3.4 E-5	1.2 E-1	NC	2B
Nickel (and compounds)		7440-02-0	TAC	2.6 E-4	9.1 E-1	A, B2 <sup>N1</sup>	1,2B <sup>N2</sup>
5-Nitroacenaphthene <sup>BaP</sup>		602-87-9	P65-E	3.7 E-5	1.3 E-1	NC	2B
6-Nitrochrysene <sup>BaP</sup>	(inhalation)	7496-02-8	TAC	1.1 E-2	3.9 E+1	NC	2B
	(oral)				1.2 E+2		
2-Nitrofluorene <sup>BaP</sup>	(inhalation)	607-57-8	TAC	1.1 E-5	3.9 E-2	NC	2B
	(oral)				1.2 E-1		
1-Nitropyrene <sup>BaP</sup>	(inhalation)	5522-43-0	TAC	1.1 E-4	3.9 E-1	NC	2B
	(oral)				1.2 E+0		
4-Nitropyrene <sup>BaP</sup>	(inhalation)	57835-92-4	TAC	1.1 E-4	3.9 E-1	NC	2B
	(oral)				1.2 E+0		
N-Nitrosodi- <i>n</i> -butylamine		924-16-3	P65-S	3.1 E-3	1.1 E+1	B2	2B
N-Nitroso-N-methylethylamine		10595-95-6	IRIS	6.3 E-3	2.2 E+1	B2	2B
N-Nitrosodi- <i>n</i> -propylamine		621-64-7	IRIS	2.0 E-3	7.0 E+0	B2	2B
N-Nitrosodiethylamine		55-18-5	P65-S	1.0 E-2	3.6 E+1	B2	2A
N-Nitrosodimethylamine		62-75-9	P65-S	4.6 E-3	1.6 E+1	B2	2A
N-Nitrosodiphenylamine		86-30-6	P65-S	2.6 E-6	9.0 E-3	B2	3
<i>p</i> -Nitrosodiphenylamine		156-10-5	P65-E	6.3 E-6	2.2 E-2	NC	3
N-Nitrosomorpholine		59-89-2	P65-E	1.9 E-3	6.7 E+0	NC	2B
N-Nitrosopiperidine		100-75-4	P65-E	2.7 E-3	9.4 E+0	NC	2B
N-Nitrosopyrrolidine		930-55-2	IRIS	6.0 E-4	2.1 E+0	B2	2B
Pentachlorophenol		87-86-5	P65-S	5.1 E-6	1.8 E-2	B2	2B

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Perchloroethylene (inhalation)	127-18-4	TAC	6.1 E-6	2.1 E-2	NC	2A
(oral)		P65-S		5.1 E-2		
Polychlorinated biphenyls (PCBs) (high risk) <sup>P</sup>	1336-36-3	IRIS	5.7 E-4	2.0 E+0	B2	2A
(for use with unspeciated (low risk) <sup>P</sup> )			1.1 E-4	4.0 E-1		
PCB mixtures (lowest risk) <sup>P</sup>			1.1 E-4	4.0 E-1		
			2.0 E-5	7.0 E-2		
(for use where measurements or estimates are available for PCB congeners) <sup>P</sup>						
PCB 77 3,3',4,4'-TCB			3.8 E-3	1.3 E+1		
PCB 81 3,4,4',5-TCB			1.1 E-2	3.9 E+1		
PCB 105 2,3,3',4,4'-PeCB			1.1 E-3	3.9 E+0		
PCB 114 2,3,4,4',5-PeCB			1.1 E-3	3.9 E+0		
PCB 118 2,3',4,4',5-PeCB			1.1 E-3	3.9 E+0		
PCB 123 2',3,4,4',5-PeCB			1.1 E-3	3.9 E+0		
PCB 126 3,3',4,4',5-PeCB			3.8 E+0	1.3 E+4		
PCB 156 2,3,3',4,4',5-HxCB			1.1 E-3	3.9 E+0		
PCB 157 2,3,3',4,4',5'-HxCB			1.1 E-3	3.9 E+0		
PCB 167 2,3',4,4',5,5'-HxCB			1.1 E-3	3.9 E+0		
PCB 169 3,3',4,4',5,5'-HxCB			1.1 E+0	3.9 E+3		
PCB 189 2,3,3',4,4',5,5'-HpCB			1.1 E-3	3.9 E+0		
Potassium bromate	7758-01-2	P65-E	1.4 E-4	4.9 E-1	NC	2B
1,3-Propane sultone	1120-71-4	P65-E	6.9 E-4	2.4 E+0	NC	2B
Propylene oxide (inhalation)	75-56-9	IRIS	3.7 E-6	1.3 E-2	B2	2B
(oral)				2.4 E-1		
<i>Tertiary</i> -butyl acetate (inhalation)	540-88-5	HS	1.3 E-6	4.7 E-3	NC	NC
(oral)				5.0 E-3		
1,1,2,2-Tetrachloroethane	79-34-5	IRIS	5.8 E-5	2.0 E-1	C	3
Thioacetamide	62-55-5	P65-E	1.7 E-3	6.1 E+0	NC	2B
2,4-Toluene diisocyanate	584-84-9	P65-E	1.1 E-5	3.9 E-2	NC	2B
2,6-Toluene diisocyanate	91-08-7	P65-E	1.1 E-5	3.9 E-2	NC	2B
1,1,2-Trichloroethane (vinyl trichloride)	79-00-5	IRIS	1.6 E-5	5.7 E-2	C	3
Trichloroethylene (inhalation)	79-01-6	TAC	2.0 E-6	7.0 E-3	NC	2A
(oral)		P65-S		1.5 E-2		
2,4,6-Trichlorophenol	88-06-2	P65-S	2.0 E-5	7.0 E-2	B2	2B
Urethane	51-79-6	P65-S	2.9 E-4	1.0 E+0	NC	2B
Vinyl chloride	75-01-4	TAC	7.8 E-5	2.7 E-1	NC	1

### Footnotes

- A see Appendix C  
 BaP see benzo[*a*]pyrene TAC document  
 C see Appendix E  
 D Listed by ARB as “Particulate Matter from Diesel-Fueled Engines”; Scientific Review Panel unit risk “reasonable estimate” = 3.0 E-4 ( $\mu\text{g}/\text{m}^3$ )<sup>-1</sup>. Range of unit risks in TAC document was 1.3 E-4 – 2.4 E-3 ( $\mu\text{g}/\text{m}^3$ )<sup>-1</sup>.  
 N1 Nickel refinery dust and nickel subsulfide are in Class A; nickel carbonyl is in Class B2  
 N2 Nickel compounds are in Class 1; metallic nickel is in Class 2B  
 NA not available  
 NC not classified  
 # [100 PCM fibers/m<sup>3</sup>]-1 ; see Appendix F  
 \* can be calculated using PEF factors contained in the benzo[*a*]pyrene TAC document  
 P See chemical summary for risk categorization and TEF factors (Appendix C)

### Source Key

- TAC Toxic Air Contaminant document, Office of Environmental Health Hazard Assessment (OEHHHA)  
 P65-S Standard Proposition 65 document, OEHHHA  
 IRIS Integrated Risk Information System, U.S. Environmental Protection Agency (US EPA)  
 P65-E Expedited Proposition 65 document, OEHHHA  
 HS Air Toxics Hot Spots document, Air and Site Assessment and Climate Indicators Branch, OEHHHA *et al.*  
 PETB Public Health Goal document, Pesticide and Environmental Toxicology Branch, OEHHHA