

Appendix B

Longview Air Toxics Health Risk Screening Analysis Summary - All Compounds

Compound	CAS No.	Longview Frequency Measured >MRL (% of samples)	Longview Annual Average Concentration (ug/m ³)	Screening Values			Potential Health Concern?
				ODEQ Ambient Benchmark Concentration - ABC (ug/m ³)	EPA Final Chronic Screening Value (ug/m ³)	WDOE Acceptable Source Impact Level - ASIL (ug/m ³)	
1,1,2,2-Tetrachloroethylene	127-18-4	4.8%	< 0.95	35	0.17	1.1	Possible
1,2,4-Trimethylbenzene	95-63-6	54.0%	< 1.16	na	na	na	No
1,2-Dimethylbenzene	95-47-6	52.4%	< 0.86	700	10	1500	No
1,3,5-Trimethylbenzene	108-67-8	12.7%	< 0.58	na	na	na	No
1,4/1,3-Dimethylbenzene	108-38-3	74.6%	< 3.14	700	10	1500	No
2,2,4-Trimethylpentane	540-84-1	3.2%	< 0.49	na	na	na	No
2,5-Dimethylbenzaldehyde ^(a)	5779-94-2	1.7%	< 0.084	na	na	na	No
2-Butanone (MEK)	78-93-3	89.8%	0.358	na	500	1000	No
2-Hexanone	591-78-6	23.8%	< 0.49	na	na	67	No
4-Ethyltoluene	622-96-8	4.8%	< 0.52	na	na	na	No
4-Methyl-2-Pentanone (MIBK)	108-10-1	4.8%	< 0.44	na	300	680	No
Acetaldehyde	75-07-0	98.3%	1.362	0.45	0.45	0.45	Yes
Acetone	67-64-1	100.0%	1.44	na	na	5900	No
Arsenic	7440-38-2	100.0%	1.20E-03	0.0002	0.00023	0.00023	Yes
Benzaldehyde	100-52-7	94.9%	0.178	na	na	na	No
Benzene	71-43-2	82.5%	1.27	0.13	0.13	0.12	Yes
Beryllium	7440-41-7	5.3%	< 1.00E-05	0.0004	0.00042	0.00042	No
Butyraldehyde	123-72-8	98.3%	0.217	na	na	na	No
Cadmium	7440-43-9	47.4%	< 2.28E-04	0.0006	0.00056	0.00056	No
Carbon Disulfide ^(a)	75-15-0	1.6%	< 0.33	800	70	100	No
Carbon Tetrachloride	56-23-5	11.1%	< 0.67	0.07	0.067	0.067	Possible
Chloroethane ^(a)	75-00-3	1.6%	< 0.28	na	1000	10000	No
Chloromethane	74-87-3	92.1%	0.84	90	9	340	No
Chromium ^(a)	1606-583-1	1.8%	< 3.29E-03	na	0.00083	0.17 ^(b)	Possible
Cobalt	7440-48-4	82.5%	3.24E-04	0.1	0.01	0.17	No
Crotonaldehyde	4170-30-3	15.3%	< 0.071	na	na	20	No
Cyclohexane	110-82-7	15.9%	< 0.66	na	na	3400	No
Dichlorodifluoromethane	75-71-8	92.1%	1.95	na	na	16000	No
Dichlorotetrafluoroethane ^(a)	76-14-2	1.6%	< 0.73	na	na	23000	No
Ethyl Benzene	100-41-4	38.1%	< 0.71	na	100	1000	No
Formaldehyde	50-00-0	100.0%	0.792	3	0.98	0.077	Yes
n-Heptane	142-82-5	57.1%	< 1.09	na	na	5500	No
Hexaldehyde	66-25-1	98.3%	0.179	na	na	na	No
n-Hexane	110-54-3	12.7%	< 0.4	7000	20	200	No
Isopropanol	67-63-0	84.1%	1.66	na	na	3300	No
Isovaleraldehyde	590-86-3	78.0%	0.176	na	na	na	No
Lead	7439-92-1	100.0%	5.16E-03	0.5	0.15	0.5	No
Manganese	7439-96-5	100.0%	7.80E-03	0.2	0.005	0.4	Yes
Methylene Chloride	75-09-2	55.6%	< 0.67	2.1	2.1	0.56	Possible
Naphthalene	91-20-3	58.1%	< 1.70E-03	0.03	0.029	170	No
Nickel	7440-02-0	52.6%	< 1.84E-03	0.05	0.0021	0.0021	No
PAH							
Benzo[<i>a</i>]pyrene ^(a)	192-97-2	1.6%	< 4.07E-04	.0009 (c.)	0.3	na	No
Coronene	191-07-1	14.5%	< 8.50E-04	na	na	na	No
Dibenzofuran	132-64-9	80.6%	1.69E-03	na	na	na	No
Dibenzothiophene	132-65-0	58.1%	< 5.45E-04	na	na	na	No
Perylene	198-55-0	0.0%	< 4.11E-04	na	na	na	No

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PAH (OR Group) - ABC ^(c)	na	50.0%	< 1.10E-03	0.0009 (c.)	na	na	Possible
Acenaphthene	83-32-9	66.1%	< 1.10E-03		0.3		
Acenaphthylene	208-96-8	53.2%	< 1.10E-03		0.3		
Anthracene	120-12-7	80.6%	1.36E-03		0.3		
Benzo[a]anthracene	56-55-3	29.0%	< 5.27E-04		0.0091		
Benzo[a]pyrene	50-32-8	32.3%	< 4.36E-04		0.00091		
Benzo[b]fluoranthene	205-99-2	45.2%	< 7.23E-04		0.0091		
Benzo[g,h,i]perylene	191-24-2	27.4%	< 4.73E-04		0.3		
Benzo[k]fluoranthene	207-08-9	21.0%	< 3.70E-04		0.0091		
Chrysene	218-01-9	27.4%	< 5.16E-04		0.091		
Dibenz[a,h]anthracene	53-70-3	12.9%	< 4.24E-04		0.00083		
Fluoranthene	206-44-0	82.3%	2.40E-03		0.3		
Fluorene	86-73-7	80.6%	3.00E-03		0.3		
Indeno[1,2,3-cd]pyrene	193-39-5	30.6%	< 5.52E-04		0.0091		
Phenanthrene	85-01-8	80.6%	1.09E-03		0.3		
Pyrene	129-00-0	80.6%	1.80E-03		0.3		
PAH (WA Group) - ABC ^(d)	na	28.3%	< 3.50E-03	na	na	0.00048^(d)	Possible
Benzo[a]anthracene	56-55-3	29.0%	< 5.27E-04		0.0091		
Benzo[a]pyrene	50-32-8	32.3%	< 4.36E-04		0.00091		
Benzo[b]fluoranthene	205-99-2	45.2%	< 7.23E-04		0.0091		
Benzo[k]fluoranthene	207-08-9	21.0%	< 3.70E-04		0.0091		
Chrysene	218-01-9	27.4%	< 5.16E-04		0.091		
Dibenz[a,h]anthracene	53-70-3	12.9%	< 4.24E-04		0.00083		
Indeno[1,2,3-cd]pyrene	193-39-5	30.6%	< 5.52E-04		0.0091		
Propionaldehyde	123-38-6	96.6%	0.232	na	na	na	No
Selenium	7782-49-2	57.9%	< 1.70E-04	na	2	0.67	No
Styrene	100-42-5	6.3%	< 0.49	na	100	1000	No
m-Tolualdehyde	620-23-5	22.0%	< 0.087	na	na	na	No
o-Tolualdehyde	529-20-4	5.1%	< 0.084	na	na	na	No
p-Tolualdehyde	104-87-0	6.8%	< 0.084	na	na	na	No
Toluene	108-88-3	92.1%	3.57	400	40	400	No
Trichloroethylene	79-01-6	9.5%	< 0.6	0.5	0.5	0.59	Possible
Trichlorofluoromethane	75-69-4	44.4%	< 0.85	na	na	19000	No
Trichlorotrifluoroethane	26523-64-8	27.0%	< 0.83	na	na	27000	No
Valeraldehyde	110-62-3	59.3%	< 0.075	na	na	590	No
1,1,1-Trichloroethane	71-55-6	0.0%	< 0.571	1000	100	6400	No
1,1,2,2-Tetrachloroethane	79-34-5	0.0%	< 0.719	na	0.017	23	Possible
1,1,2-Trichloroethane	79-00-5	0.0%	< 0.571	na	0.063	180	Possible
1,1-Dichloroethane	75-34-3	0.0%	< 0.424	na	0.63	2700	No
1,1-Dichloroethylene	75-35-4	0.0%	< 0.415	na	20	67	No
1,2,4-Trichlorobenzene	120-82-1	0.0%	< 0.777	na	20	120	No
1,2-Dibromoethane (EDB)	106-93-4	0.0%	< 0.804	0.002	0.002	0.005	Possible
1,2-Dichlorobenzene	95-50-1	0.0%	< 0.629	na	na	1000	No
1,2-Dichloroethane	107-06-2	0.0%	< 0.424	0.04	0.004	0.038	Possible
1,2-Dichloropropane	78-87-5	0.0%	< 0.484	na	0.053	4	Possible
1,3-Butadiene	106-99-0	0.0%	< 0.232	0.03	0.03	0.004	Possible
1,3-Dichlorobenzene	541-73-1	0.0%	< 0.063	na	na	na	No
1,4-Dichlorobenzene	106-46-7	0.0%	< 0.629	0.09	0.091	1.5	Possible
3-Chloropropene	107-05-1	0.0%	< 0.327	na	0.1	1	Possible
Acrylonitrile	107-13-1	0.0%	< 0.227	0.01	0.015	0.015	Possible
Bromodichloromethane	75-27-4	0.0%	< 0.701	na	na	na	No
Bromoform	75-25-2	0.0%	< 1.082	na	0.91	0.91	Possible

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Bromomethane	74-83-9	0.0%	< 0.407	5	0.5	5	No
Chlorobenzene	108-90-7	0.0%	< 0.482	na	100	150	No
Chloroform	67-66-3	0.0%	< 0.511	98	9.8	0.043	Possible
cis-1,2-Dichloroethylene	156-59-2	0.0%	< 0.415	na	na	na	No
cis-1,3-Dichloropropene	10061-01-5	0.0%	< 0.475	na	na	na	No
Dibromochloromethane	124-48-1	0.0%	< 0.892	na	na	na	No
Hexachloro-1,3-Butadiene	87-68-3	0.0%	< 1.116	na	0.045	0.7	Possible
Hexachloro-1,3-Butadiene	87-68-3	0.0%	< 1.116	na	0.045	0.7	Possible
Methyl tert-butyl ether	1634-04-4	0.0%	< 0.377	na	3.8	500	No
Tetrahydrofuran	109-99-9	0.0%	< 0.309	na	na	2000	No
trans-1,2-Dichloroethene	156-60-5	0.0%	< 0.415	na	na	na	No
trans-1,3-Dichloropropene	10061-02-6	0.0%	< 0.475	na	na	na	No
Vinyl bromide	593-60-2	0.0%	< 0.458	na	0.031	73	Possible
Vinyl Chloride	75-01-4	0.0%	< 0.268	0.1	0.11	0.012	Possible

- Compound was reported above the MRL only once over the term of the study.
- ASIL is for chromium compounds not including hexavalent chromium.
- ABC is compared to the toxicity equivalency factor weighted sum of concentrations for up to 15 individual PAH compounds (Oregon DEQ PAH Group).
- ASIL is compared to the sum of up to 7 PAH compounds as one TAP equivalent in potency to BaP (Washington DOE PAH Group).