



Attachment D

**Regulatory and Human Health Benchmarks
Used for SSL Development**



Attachment D

Regulatory and Human Health Benchmarks for SSL Development

This attachment provides regulatory and human health benchmarks necessary to calculate SSLs for 110 chemicals commonly found at National Priority List (NPL) sites. The sources of these values (shown in the following table) are regularly updated by EPA. Prior to calculating SSLs at a site, check all relevant chemical-specific values in this attachment against the most recent version of their sources to ensure that they are up-to-date.

Attachment D. Regulatory and Human Health Benchmarks Used for SSL Development

| CAS Number | Chemical Name | Maximum Contaminant Level Goal (mg/L) | | Maximum Contaminant Level (mg/L) | | Water Health Based Limits (mg/L) | | Cancer Slope Factor (mg/kg-d) ⁻¹ | | | Unit Risk Factor (µg/m ³) ⁻¹ | | | Reference Dose (mg/kg-d) | | Reference Concentration (mg/m ³) | |
|------------|-----------------------------|---------------------------------------|-------------------|----------------------------------|-------------------|----------------------------------|-----------------|---|-----------------|-------------------|---|---------|-------------------|--------------------------|-------------------|--|-------------------|
| | | MCLG (PMCLG) | Ref. ^a | MCL (PMCL) | Ref. ^a | HBL ^b | Basis | Carc. Class ^c | SF ₀ | Ref. ^a | Carc. Class ^c | URF | Ref. ^a | RfD | Ref. ^a | RfC | Ref. ^a |
| 83-32-9 | Acenaphthene | | | | | 2E+00 | RfD | | | | | | | 6.0E-02 | 1 | | |
| 67-64-1 | Acetone (2-Propanone) | | | | | 4E+00 | RfD | D | | | D | | | 1.0E-01 | 1 | | |
| 309-00-2 | Aldrin | | | | | 5E-06 | SF ₀ | B2 | 1.7E+01 | 1 | B2 | 4.9E-03 | 1 | 3.0E-05 | 1 | | |
| 120-12-7 | Anthracene | | | | | 1E+01 | RfD | D | | | D | | | 3.0E-01 | 1 | | |
| 7440-36-0 | Antimony | 6.0E-03 | 3 | 6.0E-03 | 3 | | | | | | | | | 4.0E-04 | 1 | | |
| 7440-38-2 | Arsenic | | | 5.0E-02 | 3 | | | A | 1.5E+00 | 1 | A | 4.3E-03 | 1 | 3.0E-04 | 1 | | |
| 7440-39-3 | Barium | 2.0E+00 | 3 | 2.0E+00 | 3 | | | | | | | | | 7.0E-02 | 1 | 5.0E-04 | 2 |
| 56-55-3 | Benz(a)anthracene | | | | | 1E-04 | SF ₀ | B2 | 7.3E-01 | 4 | B2 | | | | | | |
| 71-43-2 | Benzene | | | 5.0E-03 | 3 | | | A | 2.9E-02 | 1 | A | 8.3E-06 | 1 | | | | |
| 205-99-2 | Benzo(b)fluoranthene | | | | | 1E-04 | SF ₀ | B2 | 7.3E-01 | 4 | B2 | | | | | | |
| 207-08-9 | Benzo(k)fluoranthene | | | | | 1E-03 | SF ₀ | B2 | 7.3E-02 | 4 | B2 | | | | | | |
| 65-85-0 | Benzoic acid | | | | | 1E+02 | RfD | | | | | | | 4.0E+00 | 1 | | |
| 50-32-8 | Benzo(a)pyrene | | | 2.0E-04 | 3 | | | B2 | 7.3E+00 | 1 | B2 | | | | | | |
| 7440-41-7 | Beryllium | 4.0E-03 | 3 | 4.0E-03 | 3 | | | B2 | 4.3E+00 | 1 | B2 | 2.4E-03 | 1 | 5.0E-03 | 1 | | |
| 111-44-4 | Bis(2-chloroethyl)ether | | | | | 8E-05 | SF ₀ | B2 | 1.1E+00 | 1 | B2 | 3.3E-04 | 1 | | | | |
| 117-81-7 | Bis(2-ethylhexyl)phthalate | | | 6.0E-03 | 3 | | | B2 | 1.4E-02 | 1 | B2 | | | 2.0E-02 | 1 | | |
| 75-27-4 | Bromodichloromethane | | | 1.0E-01 * | 3 | | | B2 | 6.2E-02 | 1 | B2 | | | 2.0E-02 | 1 | | |
| 75-25-2 | Bromoform (tribromomethane) | | | 1.0E-01 * | 3 | | | B2 | 7.9E-03 | 1 | B2 | 1.1E-06 | 1 | 2.0E-02 | 1 | | |
| 71-36-3 | Butanol | | | | | 4E+00 | RfD | D | | | D | | | 1.0E-01 | 1 | | |
| 85-68-7 | Butyl benzyl phthalate | | | | | 7E+00 | RfD | C | | | C | | | 2.0E-01 | 1 | | |
| 7440-43-9 | Cadmium | 5.0E-03 | 3 | 5.0E-03 | 3 | | | B1 | | | B1 | 1.8E-03 | 1 | 1.0E-03** | 1 | | |
| 86-74-8 | Carbazole | | | | | 4E-03 | SF ₀ | B2 | 2.0E-02 | 2 | | | | | | | |
| 75-15-0 | Carbon disulfide | | | | | 4E+00 | RfD | | | | | | | 1.0E-01 | 1 | 7.0E-01 | 1 |
| 56-23-5 | Carbon tetrachloride | | | 5.0E-03 | 3 | | | B2 | 1.3E-01 | 1 | B2 | 1.5E-05 | 1 | 7.0E-04 | 1 | | |
| 57-74-9 | Chlordane | | | 2.0E-03 | 3 | | | B2 | 1.3E+00 | 1 | B2 | 3.7E-04 | 1 | 6.0E-05 | 1 | | |
| 106-47-8 | p-Chloroaniline | | | | | 1E-01 | RfD | | | | | | | 4.0E-03 | 1 | | |
| 108-90-7 | Chlorobenzene | 1.0E-01 | 3 | 1.0E-01 | 3 | | | D | | | D | | | 2.0E-02 | 1 | 2.0E-02 | 2 |
| 124-48-1 | Chlorodibromomethane | 6.0E-02 | 3 | 1.0E-01 * | 3 | | | C | 8.4E-02 | 1 | C | | | 2.0E-02 | 1 | | |
| 67-66-3 | Chloroform | | | 1.0E-01 * | 3 | | | B2 | 6.1E-03 | 1 | B2 | 2.3E-05 | 1 | 1.0E-02 | 1 | | |
| 95-57-8 | 2-Chlorophenol | | | | | 2E-01 | RfD | | | | | | | 5.0E-03 | 1 | | |

* Proposed MCL = 0.08 mg/L, *Drinking Water Regulations and Health Advisories*, U.S. EPA (1995).

** Cadmium RfD is based on dietary exposure.

Attachment D (continued)

| CAS Number | Chemical Name | Maximum Contaminant Level Goal (mg/L) | | Maximum Contaminant Level (mg/L) | | Water Health Based Limits (mg/L) | | Cancer Slope Factor (mg/kg-d) ⁻¹ | | | Unit Risk Factor (µg/m ³) ⁻¹ | | | Reference Dose (mg/kg-d) | | Reference Concentration (mg/m ³) | |
|------------|----------------------------|---------------------------------------|-------------------|----------------------------------|-------------------|----------------------------------|-----------------|---|-----------------|-------------------|---|---------|-------------------|--------------------------|-------------------|--|-------------------|
| | | MCLG (PMCLG) | Ref. ^a | MCL (PMCL) | Ref. ^a | HBL ^b | Basis | Carc. Class ^c | SF _o | Ref. ^a | Carc. Class ^c | URF | Ref. ^a | RfD | Ref. ^a | RfC | Ref. ^a |
| 7440-47-3 | Chromium | 1.0E-01 | 3 | 1.0E-01 | 3 | | | A | | | A | 1.2E-02 | 1 | 5.0E-03 | 1 | | |
| 16065-83-1 | Chromium (III) | | | | | 4E+01 | RfD | | | | | | | 1.0E+00 | 1 | | |
| 18540-29-9 | Chromium (VI) | | | 1.0E-01 | 3 * | | | A | | | A | 1.2E-02 | 1 | 5.0E-03 | 1 | | |
| 218-01-9 | Chrysene | | | | | 1E-02 | SF _o | B2 | 7.3E-03 | 4 | | | | | | | |
| 57-12-5 | Cyanide (amenable) | (2.0E-01) | 3 | (2.0E-01) | 3 | | | D | | | D | | | 2.0E-02 | 1 | | |
| 72-54-8 | DDD | | | | | 4E-04 | SF _o | B2 | 2.4E-01 | 1 | B2 | | | | | | |
| 72-55-9 | DDE | | | | | 3E-04 | SF _o | B2 | 3.4E-01 | 1 | B2 | | | | | | |
| 50-29-3 | DDT | | | | | 3E-04 | SF _o | B2 | 3.4E-01 | 1 | B2 | 9.7E-05 | 1 | 5.0E-04 | 1 | | |
| 53-70-3 | Dibenz(a,h)anthracene | | | | | 1E-05 | SF _o | B2 | 7.3E+00 | 4 | B2 | | | | | | |
| 84-74-2 | Di-n-butyl phthalate | | | | | 4E+00 | RfD | D | | | D | | | 1.0E-01 | 1 | | |
| 95-50-1 | 1,2-Dichlorobenzene | 6.0E-01 | 3 | 6.0E-01 | 3 | | | D | | | D | | | 9.0E-02 | 1 | 2.0E-01 | 2 |
| 106-46-7 | 1,4-Dichlorobenzene | 7.5E-02 | 3 | 7.5E-02 | 3 | | | B2 | 2.4E-02 | 2 | B2 | | | | | 8.0E-01 | 1 |
| 91-94-1 | 3,3-Dichlorobenzidine | | | | | 2E-04 | SF _o | B2 | 4.5E-01 | 1 | B2 | | | | | | |
| 75-34-3 | 1,1-Dichloroethane | | | | | 4E+00 | RfD | C | | | C | | | 1.0E-01 | 7 | 5.0E-01 | 2 |
| 107-06-2 | 1,2-Dichloroethane | | | 5.0E-03 | 3 | | | B2 | 9.1E-02 | 1 | B2 | 2.6E-05 | 1 | | | | |
| 75-35-4 | 1,1-Dichloroethylene | 7.0E-03 | 3 | 7.0E-03 | 3 | | | C | 6.0E-01 | 1 | C | 5.0E-05 | 1 | 9.0E-03 | 1 | | |
| 156-59-2 | cis-1,2-Dichloroethylene | 7.0E-02 | 3 | 7.0E-02 | 3 | | | D | | | D | | | 1.0E-02 | 2 | | |
| 156-60-5 | trans-1,2-Dichloroethylene | 1.0E-01 | 3 | 1.0E-01 | 3 | | | | | | | | | 2.0E-02 | 1 | | |
| 120-83-2 | 2,4-Dichlorophenol | | | | | 1E-01 | RfD | | | | | | | 3.0E-03 | 1 | | |
| 78-87-5 | 1,2-Dichloropropane | | | 5.0E-03 | 3 | | | B2 | 6.8E-02 | 2 | B2 | | | | | 4.0E-03 | 1 |
| 542-75-6 | 1,3-Dichloropropene | | | | | 5E-04 | SF _o | B2 | 1.8E-01 | 2 | B2 | 3.7E-05 | 2 | 3.0E-04 | 1 | 2.0E-02 | 1 |
| 60-57-1 | Dieldrin | | | | | 5E-06 | SF _o | B2 | 1.6E+01 | 1 | B2 | 4.6E-03 | 1 | 5.0E-05 | 1 | | |
| 84-66-2 | Diethylphthalate | | | | | 3E+01 | RfD | D | | | D | | | 8.0E-01 | 1 | | |
| 105-67-9 | 2,4-Dimethylphenol | | | | | 7E-01 | RfD | | | | | | | 2.0E-02 | 1 | | |
| 51-28-5 | 2,4-Dinitrophenol | | | | | 4E-02 | RfD | | | | | | | 2.0E-03 | 1 | | |
| 121-14-2 | 2,4-Dinitrotoluene** | | | | | 1E-04 | SF _o | B2 | 6.8E-01 | 1 | | | | 2.0E-03 | 1 | | |
| 606-20-2 | 2,6-Dinitrotoluene** | | | | | 1E-04 | SF _o | B2 | 6.8E-01 | 1 | | | | 1.0E-03 | 2 | | |
| 117-84-0 | Di-n-octyl phthalate | | | | | 7E-01 | RfD | | | | | | | 2.0E-02 | 2 | | |
| 115-29-7 | Endosulfan | | | | | 2E-01 | RfD | | | | | | | 6.0E-03 | 2 | | |
| 72-20-8 | Endrin | 2.0E-03 | 3 | 2.0E-03 | 3 | | | D | | | D | | | 3.0E-04 | 1 | | |

* MCL for total chromium is based on Cr (VI) toxicity.

** Cancer Slope Factor is for 2,4-, 2,6-Dinitrotoluene mixture.

Attachment D (continued)

| CAS Number | Chemical Name | Maximum Contaminant Level Goal (mg/L) | | Maximum Contaminant Level (mg/L) | | Water Health Based Limits (mg/L) | | Cancer Slope Factor (mg/kg-d) ⁻¹ | | | Unit Risk Factor (µg/m ³) ⁻¹ | | | Reference Dose (mg/kg-d) | | Reference Concentration (mg/m ³) | |
|------------|--|---------------------------------------|-------------------|----------------------------------|-------------------|----------------------------------|-----------------|---|-----------------|-------------------|---|---------|-------------------|--------------------------|-------------------|--|-------------------|
| | | MCLG (PMCLG) | Ref. ^a | MCL (PMCL) | Ref. ^a | HBL ^b | Basis | Carc. Class ^c | SF _o | Ref. ^a | Carc. Class ^c | URF | Ref. ^a | RfD | Ref. ^a | RfC | Ref. ^a |
| 100-41-4 | Ethylbenzene | 7.0E-01 | 3 | 7.0E-01 | 3 | | | D | | | D | | 1.0E-01 | 1 | 1.0E+00 | 1 | |
| 206-44-0 | Fluoranthene | | | | | 1E+00 | RfD | D | | | D | | 4.0E-02 | 1 | | | |
| 86-73-7 | Fluorene | | | | | 1E+00 | RfD | D | | | | | 4.0E-02 | 1 | | | |
| 76-44-8 | Heptachlor | | | 4.0E-04 | 3 | | | B2 | 4.5E+00 | 1 | B2 | 1.3E-03 | 1 | 5.0E-04 | 1 | | |
| 1024-57-3 | Heptachlor epoxide | | | 2.0E-04 | 3 | | | B2 | 9.1E+00 | 1 | B2 | 2.6E-03 | 1 | 1.3E-05 | 1 | | |
| 118-74-1 | Hexachlorobenzene | | | 1.0E-03 | 3 | | | B2 | 1.6E+00 | 1 | B2 | 4.6E-04 | 1 | 8.0E-04 | 1 | | |
| 87-68-3 | Hexachloro-1,3-butadiene | 1.0E-03 | 3 | | | 1E-03 | SF _o | C | 7.8E-02 | 1 | C | 2.2E-05 | 1 | 2.0E-04 | 2 | | |
| 319-84-6 | -HCH (-BHC) | | | | | 1E-05 | SF _o | B2 | 6.3E+00 | 1 | B2 | 1.8E-03 | 1 | | | | |
| 319-85-7 | -HCH (-BHC) | | | | | 5E-05 | SF _o | C | 1.8E+00 | 1 | C | 5.3E-04 | 1 | | | | |
| 58-89-9 | -HCH (Lindane) | 2.0E-04 | 3 | 2.0E-04 | 3 | | | B2 | 1.3E+00 | 2 | C | | | 3.0E-04 | 1 | | |
| 77-47-4 | Hexachlorocyclopentadiene | 5.0E-02 | 3 | 5.0E-02 | 3 | | | D | | | D | | | 7.0E-03 | 1 | 7.0E-05 | 2 |
| 67-72-1 | Hexachloroethane | | | | | 6E-03 | SF _o | C | 1.4E-02 | 1 | C | 4.0E-06 | 1 | 1.0E-03 | 1 | | |
| 193-39-5 | Indeno(1,2,3- <i>cd</i>)pyrene | | | | | 1E-04 | SF _o | B2 | 7.3E-01 | 4 | B2 | | | | | | |
| 78-59-1 | Isophorone | | | | | 9E-02 | SF _o | C | 9.5E-04 | 1 | C | | | 2.0E-01 | 1 | | |
| 7439-97-6 | Mercury | 2.0E-03 | 3 | 2.0E-03 | 3 | | | D | | | D | | | 3.0E-04 | 2 | 3.0E-04 | 2 |
| 72-43-5 | Methoxychlor | 4.0E-02 | 3 | 4.0E-02 | 3 | | | D | | | D | | | 5.0E-03 | 1 | | |
| 74-83-9 | Methyl bromide | | | | | 5E-02 | RfD | D | | | D | | | 1.4E-03 | 1 | 5.0E-03 | 1 |
| 75-09-2 | Methylene chloride | | | 5.0E-03 | 3 | | | B2 | 7.5E-03 | 1 | B2 | 4.7E-07 | 1 | 6.0E-02 | 1 | 3.0E+00 | 2 |
| 95-48-7 | 2-Methylphenol (<i>o</i> -cresol) | | | | | 2E+00 | RfD | C | | | C | | | 5.0E-02 | 1 | | |
| 91-20-3 | Naphthalene | | | | | 1E+00 | RfD | D | | | D | | | 4.0E-02 | 6 | | |
| 7440-02-0 | Nickel | | | | | 1E-01 | HA * | A | | | A | 2.4E-04 | 1 | 2.0E-02 | 1 | | |
| 98-95-3 | Nitrobenzene | | | | | 2E-02 | RfD | D | | | D | | | 5.0E-04 | 1 | 2.0E-03 | 2 |
| 86-30-6 | <i>N</i> -Nitrosodiphenylamine | | | | | 2E-02 | SF _o | B2 | 4.9E-03 | 1 | B2 | | | | | | |
| 621-64-7 | <i>N</i> -Nitrosodi- <i>n</i> -propylamine | | | | | 1E-05 | SF _o | B2 | 7.0E+00 | 1 | B2 | | | | | | |
| 87-86-5 | Pentachlorophenol | | | 1.0E-03 | 3 | | | B2 | 1.2E-01 | 1 | B2 | | | 3.0E-02 | 1 | | |
| 108-95-2 | Phenol | | | | | 2E+01 | RfD | D | | | D | | | 6.0E-01 | 1 | | |
| 129-00-0 | Pyrene | | | | | 1E+00 | RfD | D | | | D | | | 3.0E-02 | 1 | | |
| 7782-49-2 | Selenium | 5.0E-02 | 3 | 5.0E-02 | 3 | | | D | | | D | | | 5.0E-03 | 1 | | |
| 7440-22-4 | Silver | | | | | 2E-01 | RfD | D | | | D | | | 5.0E-03 | 1 | | |
| 100-42-5 | Styrene | 1.0E-01 | 3 | 1.0E-01 | 3 | | | D | | | D | | | 2.0E-01 | 1 | 1.0E+00 | 1 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | | | | 4E-04 | SF _o | C | 2.0E-01 | 1 | C | 5.8E-05 | 1 | | | | |

* Health advisory for nickel (MCL is currently remanded); EPA Office of Science and Technology, 7/10/95.

Attachment D (continued)

| CAS Number | Chemical Name | Maximum Contaminant Level Goal (mg/L) | | Maximum Contaminant Level (mg/L) | | Water Health Based Limits (mg/L) | | Cancer Slope Factor (mg/kg-d) ⁻¹ | | | Unit Risk Factor (µg/m ³) ⁻¹ | | | Reference Dose (mg/kg-d) | | Reference Concentration (mg/m ³) | |
|------------|-------------------------------|---------------------------------------|-------------------|----------------------------------|-------------------|----------------------------------|-----------------|---|-----------------|-------------------|---|---------|-------------------|--------------------------|-------------------|--|-------------------|
| | | MCLG (PMCLG) | Ref. ^a | MCL (PMCL) | Ref. ^a | HBL ^b | Basis | Carc. Class ^c | SF _o | Ref. ^a | Carc. Class ^c | URF | Ref. ^a | RfD | Ref. ^a | RfC | Ref. ^a |
| 127-18-4 | Tetrachloroethylene | | | 5.0E-03 | 3 | | | | 5.2E-02 | 5 | | 5.8E-07 | 5 | 1.0E-02 | 1 | | |
| 7440-28-0 | Thallium | 5.0E-04 | 3 | 2.0E-03 | 3 | | | | | | | | | | | | |
| 108-88-3 | Toluene | 1.0E+00 | 3 | 1.0E+00 | 3 | | | | | | | | | 2.0E-01 | 1 | 4.0E-01 | 1 |
| 8001-35-2 | Toxaphene | | | 3.0E-03 | 3 | | | | B2 | 1.1E+00 | 1 | B2 | 3.2E-04 | 1 | | | |
| 120-82-1 | 1,2,4-Trichlorobenzene | 7.0E-02 | 3 | 7.0E-02 | 3 | | | | D | | | D | | 1.0E-02 | 1 | 2.0E-01 | 2 |
| 71-55-6 | 1,1,1-Trichloroethane | 2.0E-01 | 3 | 2.0E-01 | 3 | | | | D | | | D | | | | 1.0E+00 | 5 |
| 79-00-5 | 1,1,2-Trichloroethane | 3.0E-03 | 3 | 5.0E-03 | 3 | | | | C | 5.7E-02 | 1 | C | 1.6E-05 | 1 | 4.0E-03 | 1 | |
| 79-01-6 | Trichloroethylene | zero | 3 | 5.0E-03 | 3 | | | | | 1.1E-02 | 5 | | 1.7E-06 | 5 | | | |
| 95-95-4 | 2,4,5-Trichlorophenol | | | | | 4E+00 | RfD | | | | | | | 1.0E-01 | 1 | | |
| 88-06-2 | 2,4,6-Trichlorophenol | | | | | 8E-03 | SF _o | B2 | 1.1E-02 | 1 | B2 | 3.1E-06 | 1 | | | | |
| 7440-62-2 | Vanadium | | | | | 3E-01 | RfD | | | | | | | 7.0E-03 | 2 | | |
| 108-05-4 | Vinyl acetate | | | | | 4E+01 | RfD | | | | | | | 1.0E+00 | 1 | 2.0E-01 | 1 |
| 75-01-4 | Vinyl chloride (chloroethene) | | | 2.0E-03 | 3 | | | A | 1.9E+00 | 2 | A | 8.4E-05 | 2 | | | | |
| 108-38-3 | <i>m</i> -Xylene | 1.0E+01 | 3 * | 1.0E+01 | 3 * | | | D | | | D | | 2.0E+00 | 2 | | | |
| 95-47-6 | <i>o</i> -Xylene | 1.0E+01 | 3 * | 1.0E+01 | 3 * | | | D | | | D | | 2.0E+00 | 2 | | | |
| 106-42-3 | <i>p</i> -Xylene | 1.0E+01 | 3 * | 1.0E+01 | 3 * | | | D | | | D | | 2.0E+00 | 1 ** | | | |
| 7440-66-6 | Zinc | | | | | 1E+01 | RfD | D | | | D | | 3.0E-01 | 1 | | | |

* MCL for total xylenes [1330-20-7] is 10 mg/L.

** RfD for total xylenes is 2 mg/kg-day.

- ^a References:
- 1 = IRIS, U.S. EPA (1995)
 - 2 = HEAST, U.S. EPA (1995)
 - 3 = U.S. EPA (1995)
 - 4 = OHEA, U.S. EPA (1993)
 - 5 = Interim toxicity criteria provided by Superfund Health Risk Technical Support Center, Environmental Criteria Assessment Office (ECAO), Cincinnati, OH (1994)
 - 6 = ECAO, U.S. EPA (1994i)
 - 7 = ECAO, U.S. EPA (1994h)

^c Categorization of overall weight of evidence for human carcinogenicity:

- Group A: human carcinogen
- Group B: probable human carcinogen
 - B1: limited evidence from epidemiologic studies
 - B2: "sufficient" evidence from animal studies and "inadequate" evidence or "no data" from epidemiologic studies
- Group C: possible human carcinogen
- Group D: not classifiable as to health carcinogenicity
- Group E: evidence of noncarcinogenicity for humans

^b Health Based Limits calculated for 30-year exposure duration, 10⁻⁶ risk or hazard quotient = 1.