

ILLINOIS POLLUTION CONTROL BOARD
January 24, 2002

IN THE MATTER OF:)
)
PROPOSED AMENDMENTS TO TIERED) R00-19(C)
APPROACH TO CORRECTIVE ACTION) (Rulemaking – Land)
OBJECTIVES (TACO) (MTBE): 35 ILL.)
ADM. CODE 742)

Adopted Rule. Final Notice.

OPINION AND ORDER OF THE BOARD (by C.A. Manning, N.J. Melas, R.C. Flemal):

By today's order, the Board adopts amendments to the Tiered Approach to Corrective Action Objectives (TACO) found at Part 742 of the Board's land regulations (35 Ill. Adm. Code 742). The TACO rules were originally adopted by the Board on June 5, 1997. *See Tiered Approach to Corrective Action Objectives (TACO): 35 Ill. Adm. Code 742, R97-12(A)* (June 5, 1997). Part 742 contains procedures for developing remediation objectives based on risks to human health and the environment posed by environmental conditions at sites undergoing remediation in the Site Remediation Program, the Leaking Underground Storage Tank Program, and pursuant to the Resource Conservation and Recovery Act (RCRA) Part B permits and closures.¹

PROCEDURAL HISTORY

Background

On May 15, 2000, the Illinois Environmental Protection Agency (Agency) submitted proposed amendments to the TACO regulations. The Board accepted this matter for hearing on May 18, 2000. On July 27, 2000, the Board moved the Agency's proposed rulemaking to first notice. In doing so, the Board divided the proposal into two subdockets, A and B.² On

¹ The TACO regulations provide for a three-tiered approach to cleanup objectives. Under a Tier 1 analysis, an applicant compares levels of contaminants of concern at the remediation site to pre-determined remediation objectives. *See Tiered Approach to Corrective Action Objectives (TACO): 35 Ill. Adm. Code 742, R97-12(A)* (June 5, 1997). For a Tier 2 analysis, an applicant uses site-specific information and equations set forth in the rules to develop alternative remediation objectives for contaminants of concern. *Id.* Finally, a Tier 3 analysis provides greater flexibility by allowing an applicant to develop site-specific remediation objectives using alternative parameters not found in Tier 1 or Tier 2. *Id.*

² The Subocket A amendments were adopted on December 21, 2000 (Proposed Amendments to Tiered Approach to Corrective Action Objectives (TACO): 35 Ill. Adm. Code 742, R00-19(A) (Dec. 21, 2000)), and the Subocket B amendments were adopted on July 26, 2001

June 7, 2001, the Board opened this Subocket C for the purpose of addressing the proposed cleanup standards for methyl tertiary-butyl ether (MTBE) that were originally contained in Subocket B. The Board adopted its first-notice opinion and order in this matter on September 6, 2001, and it was published on September 21, 2001. 25 Ill. Reg. 11994. The Board's second-notice opinion and order was adopted on December 6, 2001. Following adoption of the second-notice opinion and order, the Board submitted the proposal to the Joint Committee on Administrative Rules (JCAR) for consideration. JCAR considered the proposal at its January 9, 2002 meeting and voted "no objection" to the Board proceeding with the amendments. The Board today adopts the proposed MTBE cleanup standards which, with only minor exceptions, are identical to those amendments that were originally proposed by the Board at first notice.

The Board has coordinated this rulemaking with another pending Agency proposal that will add groundwater quality standards for MTBE. *See generally Proposed MTBE Groundwater Quality Standards Amendments: 35 Ill. Adm. Code 620, R01-14.* Today the Board also adopts the proposed amendments in R01-14.

Subocket C Amendments

There have been only a few minor, non-substantive changes to the rule from that proposed by the Board in its first-notice opinion and order. These minor changes amount to basically typographical changes prompted by comments from JCAR. The Board received two public comments during the public comment period: one comment from the Agency and one from the Illinois Petroleum Council. The public comments did not seek to change the substance of the rule proposed at first notice. Rather, both public comments supported the substance of the Board's proposed addition of MTBE. As a result, no changes were made to the substance of the rule in response to the public comments. For a more detailed discussion of the public comments, please see the Board's second-notice opinion and order. Proposed Amendments to Tiered Approach to Corrective Action Objectives (TACO) (MTBE): 35 Ill. Adm. Code 742, R00-19(C) (Dec. 6, 2001).

Based on the record developed through hearing and public comments, the Board has determined that MTBE should be added to the TACO regulations. Accordingly, the Board adopts these amendments to 35 Ill. Adm. Code 742.

ORDER

The Board hereby adopts these amendments to the TACO regulations and directs the Clerk to file the following adopted amendments with the Secretary of State.

(Proposed Amendments to Tiered Approach to Corrective Action Objectives (TACO): 35 Ill. Adm. Code 742, R00-19(B) (July 26, 2001)).

**TITLE 35: ENVIRONMENTAL PROTECTION
SUBTITLE G: WASTE DISPOSAL
CHAPTER I: POLLUTION CONTROL BOARD
SUBCHAPTER f: RISK BASED CLEANUP OBJECTIVES
PART 742**

TIERED APPROACH TO CORRECTIVE ACTION OBJECTIVES

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SUBPART E: TIER 1 EVALUATION

| | |
|---------|--|
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| 742.505 | Tier 1 Soil and Groundwater Remediation Objectives |
| 742.510 | Tier 1 Remediation Objectives Tables |

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| | |
|---------|------------------------------|
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| 742.605 | Land Use |
| 742.610 | Chemical and Site Properties |

SUBPART G: TIER 2 SOIL EVALUATION

| | |
|---------|---|
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| 742.705 | Parameters for Soil Remediation Objective Equations |
| 742.710 | SSL Soil Equations |
| 742.715 | RBCA Soil Equations |
| 742.720 | Chemicals with Cumulative Noncarcinogenic Effects |

SUBPART H: TIER 2 GROUNDWATER EVALUATION

| | |
|---------|--|
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SUBPART I: TIER 3 EVALUATION

| | |
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| 742.905 | Modifications of Parameters |
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| 742.925 | Exposure Routes |
| 742.930 | Derivation of Toxicological Data |

SUBPART J: INSTITUTIONAL CONTROLS

| | |
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SUBPART K: ENGINEERED BARRIERS

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ILLUSTRATION B Developing Groundwater Remediation Objectives Under the Tiered Approach

TABLE A Soil Saturation Limits (C_{sat}) for Chemicals Whose Melting Point is Less than 30°C

TABLE B Tolerance Factor (K)

TABLE C Coefficients $\{A_{N-i+1}\}$ for W Test of Normality, for N=2(1)50

TABLE D Percentage Points of the W Test for n=3(1)50

TABLE E Similar-Acting Noncarcinogenic Chemicals

TABLE F Similar-Acting Carcinogenic Chemicals

TABLE G Concentrations of Inorganic Chemicals in Background Soils

TABLE H Chemicals Whose Tier 1 Class I Groundwater Remediation Objective Exceeds the 1 in 1,000,000 Cancer Risk Concentration

APPENDIX B Tier 1 Tables and Illustrations

ILLUSTRATION A Tier 1 Evaluation

TABLE A Tier 1 Soil Remediation Objectives for Residential Properties

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TABLE C pH Specific Soil Remediation Objectives for Inorganics and Ionizing Organics for the Soil Component of the Groundwater Ingestion Route (Class I Groundwater)

TABLE D pH Specific Soil Remediation Objectives for Inorganics and Ionizing Organics for the Soil Component of the Groundwater Ingestion Route (Class II Groundwater)

TABLE E Tier 1 Groundwater Remediation Objectives for the Groundwater Component of the Groundwater Ingestion Route

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APPENDIX C Tier 2 Tables and Illustrations

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 TABLE J Values to be Substituted for k_d or k_s when Evaluating Inorganics as a Function of pH (cm³/g or L/kg or cm³_{water}/g_{soil})
 TABLE K Parameter Estimates for Calculating Water-Filled Soil Porosity (θ_w)

AUTHORITY: Implementing Sections 22.4, 22.12, Title XVI, and Title XVII and authorized by Sections 27 and 58.5 of the Environmental Protection Act [415 ILCS 5/22.4, 22.12, 27, and 58.5 and Title XVI and Title XVII].

SOURCE: Adopted in R97-12(A) at 21 Ill. Reg. 7942, effective July 1, 1997; amended in R97-12(B) at 21 Ill. Reg. 16391, effective December 8, 1997; amended in R97-12(C) at 22 Ill. Reg. 10847, effective June 8, 1998; amended in R00-19(A) at 25 Ill. Reg. 651, effective January 6, 2001; amended in R00-19(B) at 25 Ill. Reg. 10374, effective August 15, 2001; amended in R00-19(C) at 26 Ill. Reg. ___, effective ____.

NOTE: Capitalization indicates statutory language.

Section 742.APPENDIX A: General

Section 742.TABLE A: Soil Saturation Limits (C_{sat})for Chemicals Whose Melting Point is Less than 30° C

| CAS No. | Chemical Name | C_{sat} (mg/kg) |
|----------|---|-------------------|
| 67-64-1 | Acetone | 100,000 |
| 71-43-2 | Benzene | 870 |
| 111-44-4 | Bis(2-chloroethyl)ether | 3,300 |
| 117-81-7 | Bis(2-ethylhexyl)phthalate | 31,000 |
| 75-27-4 | Bromodichloromethane (Dichlorobromomethane) | 3,000 |
| 75-25-2 | Bromoform | 1,900 |
| 71-36-3 | Butanol | 10,000 |
| 85-68-7 | Butyl benzyl phthalate | 930 |
| 75-15-0 | Carbon disulfide | 720 |
| 56-23-5 | Carbon tetrachloride | 1,100 |
| 108-90-7 | Chlorobenzene (Monochlorobenzene) | 680 |
| 124-48-1 | Chlorodibromomethane (Dibromochloromethane) | 1,300 |
| 67-66-3 | Chloroform | 2,900 |

| | | |
|------------------|---|--------------|
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 1,400 |
| 106-93-4 | 1,2-Dibromoethane (Ethylene dibromide) | 2,800 |
| 84-74-2 | Di- <i>n</i> -butyl phthalate | 2,300 |
| 95-50-1 | 1,2-Dichlorobenzene (o-Dichlorobenzene) | 560 |
| 75-34-3 | 1,1-Dichloroethane | 1,700 |
| 107-06-2 | 1,2-Dichloroethane (Ethylene dichloride) | 1,800 |
| 75-35-4 | 1,1-Dichloroethylene | 1,500 |
| 156-59-2 | <i>cis</i> -1,2-Dichloroethylene | 1,200 |
| 156-60-5 | <i>trans</i> -1,2-Dichloroethylene | 3,100 |
| 78-87-5 | 1,2-Dichloropropane | 1,100 |
| 542-75-6 | 1,3-Dichloropropene (1,3-Dichloropropylene, <i>cis</i> + <i>trans</i>) | 1,400 |
| 84-66-2 | Diethyl phthalate | 2,000 |
| 117-84-0 | Di- <i>n</i> -octyl phthalate | 10,000 |
| 100-41-4 | Ethylbenzene | 400 |
| 77-47-4 | Hexachlorocyclopentadiene | 2,200 |
| 78-59-1 | Isophorone | 4,600 |
| 74-83-9 | Methyl bromide (Bromomethane) | 3,200 |
| <u>1634-04-4</u> | <u>Methyl tert-butyl ether</u> <u>Methyl tertiary-butyl ether</u> | <u>8,800</u> |
| 75-09-2 | Methylene chloride (Dichloromethane) | 2,400 |
| 98-95-3 | Nitrobenzene | 1,000 |
| 100-42-5 | Styrene | 1,500 |
| 127-18-4 | Tetrachloroethylene (Perchloroethylene) | 240 |
| 108-88-3 | Toluene | 650 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 3,200 |
| 71-55-6 | 1,1,1-Trichloroethane | 1,200 |
| 79-00-5 | 1,1,2-Trichloroethane | 1,800 |
| 79-01-6 | Trichloroethylene | 1,300 |
| 108-05-4 | Vinyl acetate | 2,700 |
| 75-01-4 | Vinyl chloride | 1,200 |
| 108-38-3 | m-Xylene | 420 |

| | | |
|-----------|---------------------------|--------|
| 95-47-6 | o-Xylene | 410 |
| 106-42-3 | p-Xylene | 460 |
| 1330-20-7 | Xylenes (total) | 320 |
| | Ionizable Organics | |
| 95-57-8 | 2-Chlorophenol | 53,000 |
| | | |

(Source: Amended at 26 Ill. Reg. _____, effective _____)

Section 742.APPENDIX A: General

Section 742.TABLE E: Similar-Acting Noncarcinogenic Chemicals

| | |
|--|---|
| <u>Kidney</u> | <u>Central Nervous System</u> |
| Acetone | Butanol (Ingestion only) |
| Cadmium (Ingestion only) | Cyanide (amenable) |
| Chlorobenzene | 2,4-Dimethylphenol |
| Dalapon | Endrin |
| 1,1-Dichloroethane | Manganese |
| Di-n-octyl phthalate (Ingestion only) | 2-Methylphenol |
| Endosulfan | Mercury (Inhalation only) |
| Ethylbenzene | Styrene (Inhalation only) |
| Fluoranthene | Toluene (Inhalation only) |
| Nitrobenzene | Xylenes (Ingestion only) |
| <u>Pyrene</u> | |
| Toluene (Ingestion only) | <u>Circulatory System</u> |
| 2,4,5-Trichlorophenol | Antimony |
| Vinyl Acetate (Ingestion only) | Barium (Ingestion only) |
| <u>Liver</u> | <u>2,4-D</u> |
| Aceanaphthene | cis-1,2-Dichloroethylene (Ingestion only) |
| Acetone (Ingestion only) | Nitrobenzene |
| Butylbenzyl phthalate (Ingestion only) | trans-1,2-Dichloroethylene (Ingestion only) |
| Chlorobenzene (Ingestion only) | 2,4-Dimethylphenol |
| 1,1-Dichloroethylene (Ingestion only) | Fluoranthene |
| Di-n-octyl phthalate (Ingestion only) | Fluorene |
| Endrin | Styrene (Ingestion only) |
| Ethylbenzene | Zinc |
| Fluoranthene | <u>Gastrointestinal System</u> |
| Nitrobenzene | Beryllium (Ingestion only) |
| Picloram | Endothall |
| Styrene (Ingestion only) | Hexachlorocyclopentadiene (Ingestion only) |
| 2,4,5-TP (Silvex) | Methyl bromide (Ingestion only) |
| Toluene (Ingestion only) | |
| 1,2,4-Trichlorobenzene (Inhalation only) | |
| 2,4,5-Trichlorophenol | |

Reproductive System

Barium (Inhalation only)

Boron (Ingestion only)

Carbon disulfide

2-Chlorophenol (Ingestion only)

1,2 Dibromo 3-Chloropropane (Inhalation only)

Dinoseb

Ethylbenzene (Inhalation only)

Methoxychlor

Phenol

Cholinesterase Inhibition

Aldicarb

Carbofuran

Decreased Body Weight Gainsand Circulatory System Effects

Atrazine

Simazine

Adrenal GlandNitrobenzene

1,2,4 Trichlorobenzene (ingestion only)

Respiratory System

1,2 Dichloropropane (Inhalation only)

1,3 Dichloropropylene (Inhalation only)

Hexachlorocyclopentadiene (Inhalation only)

Methyl bromide (Inhalation only)

Naphthalene (Inhalation only)

Toluene (Inhalation only)

Vinyl acetate (Inhalation only)

Immune System

2,4 Dichlorophenol

p-Chloroaniline

Mercury (Ingestion only)

Adrenal GlandNitrobenzene1,2,4-Trichlorobenzene (Ingestion only)KidneyAcetone (Ingestion only)Cadmium (Ingestion only)ChlorobenzeneDalapon1,1-DichloroethaneDi-n-octyl phthalate (Ingestion only)EndosulfanEthylbenzeneFluorantheneMethyl tert butyl ether (Inhalation only)Methyl tertiary-butyl ether (Inhalation only)NitrobenzenePyreneToluene (Ingestion only)2,4,5-TrichlorophenolVinyl acetate (Ingestion only)LiverAcenaphtheneAcetone (Ingestion only)Butylbenzyl phthalate (Ingestion only)Chlorobenzene (Ingestion only)1,1-Dichloroethylene (Ingestion only)Di-n-octyl phthalate (Ingestion only)EndrinEthylbenzeneFluorantheneMethyl tert butyl ether (Inhalation only)Methyl tertiary-butyl ether (Inhalation only)NitrobenzenePicloramStyrene (Ingestion only)2,4,5-TP (Silvex)Toluene (Ingestion only)1,2,4-Trichlorobenzene (Inhalation only)2,4,5-TrichlorophenolCentral Nervous SystemButanol (Ingestion only)Cyanide (amenable)2,4-Demethylphenol2,4-DimethylphenolEndrinManganese2,Methylphenol2-MethylphenolMercury (Inhalation only)Styrene (Inhalation only)Toluene (Inhalation only)Xylenes (Ingestion only)Circulatory SystemAntimonyBarium (Ingestion only)2,4-Dcis-1,2-Dichloroethylene (Ingestion only)Nitrobenzenetrans-1,2-Dichloroethylene (Ingestion only)2,4-DimethylphenolFluorantheneFluoreneStyrene (Ingestion only)ZincGastrointestinal SystemBeryllium (Ingestion only)EndothallHexachlorocyclopentadiene (Ingestion only)Methyl bromide (Ingestion only)Methyl tert butyl ether (Ingestion only)Methyl tertiary-butyl ether (Ingestion only)

Immune System2,4-Dichlorophenolp-ChloroanilineMercury (Ingestion only)Reproductive SystemBarium (Inhalation only)Boron (Ingestion only)Carbon disulfide2-Chlorophenol (Ingestion only)1,2 Dibromo-3-Chloropropane (Inhalation only)DinosebEthylbenzene (Inhalation only)MethoxychlorPhenolRespiratory System1,2-Dichloropropane (Inhalation only)1,3-Dichloropropylene (Inhalation only)Hexachlorocyclopentadiene (Inhalation Inhalation only)Methyl bromide (Inhalation only)Naphthalene (Inhalation only)Toluene (Inhalation only)Vinyl acetate (Inhalation only)Cholinesterase InhibitionAldicarbCarbofuranDecreased Body Weight Gainsand Circulatory System EffectsAtrazineSimazine

(Source: Amended at 26 Ill. Reg. ____, effective ____)

Section 742.APPENDIX B: Tier 1 Tables and Illustrations

Section 742.TABLE A: Tier 1 Soil Remediation Objectives^a for Residential Properties

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | ADL (mg/kg) |
|------------|-------------------------------|--|-----------------------|---|---------------------|----------------|
| | | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | |
| 83-32-9 | Acenaphthene | 4,700 ^b | --- ^c | 570 ^b | 2,900 | * |
| 67-64-1 | Acetone | 7,800 ^b | 100,000 ^d | 16 ^b | 16 | * |
| 15972-60-8 | Alachlor ^e | 8 ^e | --- ^c | 0.04 | 0.2 | NA |
| 116-06-3 | Aldicarb ^e | 78 ^b | --- ^c | 0.013 | 0.07 | NA |
| 309-00-2 | Aldrin | 0.04 ^e | 3 ^e | 0.5 ^e | 2.5 | 0.94 |
| 120-12-7 | Anthracene | 23,000 ^b | --- ^c | 12,000 ^b | 59,000 | * |
| 1912-24-9 | Atrazine ^e | 2700 ^b | --- ^c | 0.066 | 0.33 | NA |
| 71-43-2 | Benzene | 12 ^e | 0.8 ^e | 0.03 | 0.17 | * |
| 56-55-3 | Benzo(<i>a</i>)anthracene | 0.9 ^e | --- ^c | 2 | 8 | * |
| 205-99-2 | Benzo(<i>b</i>)fluoranthene | 0.9 ^e | --- ^c | 5 | 25 | * |

| | | Exposure Route-Specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | | |
|-----------|--|--|---------------------|---|---------------------|-------------|--|
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | ADL (mg/kg) | |
| 207-08-9 | Benzo(<i>k</i>)fluroanthene | 9 ^e | --- ^c | 49 | 250 | * | |
| 50-32-8 | Benzo(<i>a</i>)pyrene | 0.09 ^{e,f} | --- ^c | 8 | 82 | * | |
| 111-44-4 | Bis(2-chloroethyl)ether | 0.6 ^e | 0.2 ^{e,f} | 0.0004 ^{e,f} | 0.0004 | 0.66 | |
| 117-81-7 | Bis(2-ethylhexyl)phthalate | 46 ^e | 31,000 ^d | 3,600 | 31,000 ^d | * | |
| 75-27-4 | Bromodichloromethane (Dichlorobromomethane) | 10 ^e | 3,000 ^d | 0.6 | 0.6 | * | |
| 75-25-2 | Bromoform | 81 ^e | 53 ^e | 0.8 | 0.8 | * | |
| 71-36-3 | Butanol | 7,800 ^b | 10,000 ^d | 17 ^b | 17 | NA | |
| 85-68-7 | Butyl benzyl phthalate | 16,000 ^b | 930 ^d | 930 ^d | 930 ^d | * | |
| 86-74-8 | Carbazole | 32 ^e | --- ^c | 0.6 ^e | 2.8 | NA | |
| 1563-66-2 | Carbofuran ^o | 390 ^b | --- ^c | 0.22 | 1.1 | NA | |
| 75-15-0 | Carbon disulfide | 7,800 ^b | 720 ^d | 32 ^b | 160 | * | |

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | ADL (mg/kg) |
|----------|--|--|--------------------|---|------------------|-------------|
| | | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | |
| 56-23-5 | Carbon tetrachloride | 5 ^e | 0.3 ^e | 0.07 | 0.33 | * |
| 57-74-9 | Chlordane | 1.8 ^e | 72 ^e | 10 | 48 | * |
| 106-47-8 | 4-Chloroaniline (<i>p</i> -Chloroaniline) | 310 ^b | --- | 0.7 ^b | 0.7 | * |
| 108-90-7 | Chlorobenzene (Monochlorobenzene) | 1,600 ^b | 130 ^b | 1 | 6.5 | * |
| 124-48-1 | Chlorodibromomethane (Dibromochloromethane) | 1,600 ^b | 1,300 ^d | 0.4 | 0.4 | * |
| 67-66-3 | Chloroform | 100 ^e | 0.3 ^e | 0.6 | 2.9 | * |
| 218-01-9 | Chrysene | 88 ^e | --- | 160 | 800 | * |
| 94-75-7 | 2,4-D ^a | 780 ^b | --- | 1.5 | 7.7 | * |
| 75-99-0 | Dalapon ^o | 2,300 ^b | --- | 0.85 | 8.5 | * |
| 72-54-8 | DDD | 3 ^e | --- | 16 ^e | 80 | * |
| 72-55-9 | DDE | 2 ^e | --- | 54 ^e | 270 | * |

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| | | Exposure Route-Specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | | |
|----------|--|--|---------------------|---|--------------------|-------------|--|
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | ADL (mg/kg) | |
| 50-29-3 | DDT | 2 ^e | --- ^g | 32 ^e | 160 | * | |
| 53-70-3 | Dibenzo(<i>a,h</i>)anthracene | 0.09 ^{e,f} | --- ^c | 2 | 7.6 | * | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 0.46 ^e | 11 ^b | 0.002 | 0.002 | * | |
| 106-93-4 | 1,2-Dibromoethane (Ethylene dibromide) | 0.0075 ^e | 0.17 ^e | 0.0004 | 0.004 | 0.005 | |
| 84-74-2 | Di- <i>n</i> -butyl phthalate | 7,800 ^b | 2,300 ^d | 2,300 ^d | 2,300 ^d | * | |
| 95-50-1 | 1,2-Dichlorobenzene (<i>o</i> - Dichlorobenzene) | 7,000 ^b | 560 ^d | 17 | 43 | * | |
| 106-46-7 | 1,4-Dichlorobenzene (<i>p</i> - Dichlorobenzene) | --- ^c | 11,000 ^b | 2 | 11 | * | |
| 91-94-1 | 3,3'-Dichlorobenzidine | 1 ^e | --- ^c | 0.007 ^{e,f} | 0.033 | 1.3 | |
| 75-34-3 | 1,1-Dichloroethane | 7,800 ^b | 1,300 ^b | 23 ^b | 110 | * | |

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | ADL (mg/kg) |
|----------|--|--|--------------------|---|------------------|-------------|
| | | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | |
| 107-06-2 | 1,2-Dichloroethane (Ethylene dichloride) | 7 ^e | 0.4 ^e | 0.02 | 0.1 | * |
| 75-35-4 | 1,1-Dichloroethylene | 700 ^b | 1,500 ^d | 0.06 | 0.3 | * |
| 156-59-2 | <i>cis</i> -1,2-Dichloroethylene | 780 ^b | 1,200 ^d | 0.4 | 1.1 | * |
| 156-60-5 | <i>trans</i> -1,2-Dichloroethylene | 1,600 ^b | 3,100 ^d | 0.7 | 3.4 | * |
| 78-87-5 | 1,2-Dichloropropane | 9 ^e | 15 ^b | 0.03 | 0.15 | * |
| 542-75-6 | 1,3-Dichloropropene (1,3-Dichloropropylene, <i>cis</i> + <i>trans</i>) | 6.4 ^e | 1.1 ^e | 0.004 ^e | 0.02 | 0.005 |
| 60-57-1 | Dieldrin ⁿ | 0.04 ^e | 1 ^e | 0.004 ^e | 0.02 | 0.603 |
| 84-66-2 | Diethyl phthalate | 63,000 ^b | 2,000 ^d | 470 ^b | 470 | * |
| 105-67-9 | 2,4-Dimethylphenol | 1,600 ^b | --- ^c | 9 ^b | 9 | * |
| 121-14-2 | 2,4-Dinitrotoluene | 0.9 ^e | --- ^c | 0.0008 ^{e,f} | 0.0008 | 0.250 |

| | | Exposure Route-Specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | |
|-----------|--|--|---------------------|---|---------------------|-------------|
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | ADL (mg/kg) |
| 606-20-2 | 2,6-Dinitrotoluene | 0.9 ^e | --- ^c | 0.0007 ^{e,f} | 0.0007 | 0.260 |
| 117-84-0 | Di- <i>n</i> -octyl phthalate | 1,600 ^b | 10,000 ^d | 10,000 ^d | 10,000 ^d | * |
| 115-29-7 | Endosulfan ^o | 470 ^b | --- ^c | 18 ^b | 90 | * |
| 145-73-3 | Endothall ^o | 1,600 ^b | --- ^c | 0.4 | 0.4 | NA |
| 72-20-8 | Endrin | 23 ^b | --- ^c | 1 | 5 | * |
| 100-41-4 | Ethylbenzene | 7,800 ^b | 400 ^d | 13 | 19 | * |
| 206-44-0 | Fluoranthene | 3,100 ^b | --- ^c | 4,300 ^b | 21,000 | * |
| 86-73-7 | Fluorene | 3,100 ^b | --- ^c | 560 ^b | 2,800 | * |
| 76-44-8 | Heptachlor | 0.1 ^e | 0.1 ^e | 23 | 110 | 0.871 |
| 1024-57-3 | Heptachlor epoxide | 0.07 ^e | 5 ^e | 0.7 | 3.3 | 1.005 |
| 118-74-1 | Hexachlorobenzene | 0.4 ^e | 1 ^e | 2 | 11 | * |
| 319-84-6 | <i>Alpha</i> -HCH (<i>alpha</i> -BHC) | 0.1 ^e | 0.8 ^e | 0.0005 ^{e,f} | 0.003 | 0.0074 |

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | ADL (mg/kg) |
|------------------|--|--|--------------------------|---|--------------------|-------------|
| | | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | |
| 58-89-9 | <i>Gamma-HCH</i> (Lindane) ⁿ | 0.5 ^e | --- ^c | 0.009 | 0.047 | * |
| 77-47-4 | Hexachlorocyclopentadiene | 550 ^b | 10 ^b | 400 | 2,200 ^d | * |
| 67-72-1 | Hexachloroethane | 78 ^b | --- ^c | 0.5 ^b | 2.6 | * |
| 193-39-5 | Indeno(1,2,3- <i>c,d</i>)pyrene | 0.9 ^e | --- ^c | 14 | 69 | * |
| 78-59-1 | Isophorone | 15,600 ^b | 4,600 ^d | 8 ^b | 8 | * |
| 72-43-5 | Methoxychlor ^o | 390 ^b | --- ^c | 160 | 780 | * |
| 74-83-9 | Methyl bromide (Bromomethane) | 110 ^b | 10 ^b | 0.2 ^b | 1.2 | * |
| <u>1634-04-4</u> | <u>Methyl tert-butyl ether</u> <u>Methyl tertiary-butyl ether</u> | <u>780^b</u> | <u>8,800^d</u> | <u>0.32</u> | <u>0.32</u> | * |
| 75-09-2 | Methylene chloride (Dichloromethane) | 85 ^e | 13 ^e | 0.02 ^e | 0.2 | * |
| 95-48-7 | 2-Methylphenol (<i>o</i> -Cresol) | 3,900 ^b | --- ^c | 15 ^b | 15 | * |
| 91-20-3 | Naphthalene | 1,600 ^b | 170 ^b | 12 ^b | 18 | * |
| 98-95-3 | Nitrobenzene | 39 ^b | 92 ^b | 0.1 ^{b,f} | 0.1 | 0.26 |

| | | Exposure Route-Specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | |
|-----------|---|--|--------------------|---|------------------|-------------|
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | ADL (mg/kg) |
| 86-30-6 | <i>N</i> -Nitrosodiphenylamine | 130 ^e | --- ^c | 1 ^e | 5.6 | * |
| 621-64-7 | <i>N</i> -Nitrosodi- <i>n</i> -propylamine | 0.09 ^{e,f} | --- ^c | 0.00005 ^{e,f} | 0.00005 | 0.0018 |
| 108-95-2 | Phenol | 47,000 ^b | --- ^c | 100 ^b | 100 | * |
| 1918-02-1 | Picloram ^o | 5,500 ^b | --- ^c | 2 | 20 | NA |
| 1336-36-3 | Polychlorinated biphenyls (PCBs) ⁿ | 1 ^h | --- ^{c,h} | --- ^h | --- ^h | * |
| 129-00-0 | Pyrene | 2,300 ^b | --- ^c | 4,200 ^b | 21,000 | * |
| 122-34-9 | Simazine ^o | 390 ^b | --- ^c | 0.04 | 0.37 | NA |
| 100-42-5 | Styrene | 16,000 ^b | 1,500 ^d | 4 | 18 | * |
| 127-18-4 | Tetrachloroethylene (Perchloroethylene) | 12 ^e | 11 ^e | 0.06 | 0.3 | * |
| 108-88-3 | Toluene | 16,000 ^b | 650 ^d | 12 | 29 | * |

| | | Exposure Route-Specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | |
|-----------|------------------------|--|--------------------|---|------------------|-------------|
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | ADL (mg/kg) |
| 8001-35-2 | Toxaphene ^a | 0.6 ^e | 89 ^e | 31 | 150 | * |
| 120-82-1 | 1,2,4-Trichlorobenzene | 780 ^b | 3,200 ^b | 5 | 53 | * |
| 71-55-6 | 1,1,1-Trichloroethane | --- ^c | 1,200 ^d | 2 | 9.6 | * |
| 79-00-5 | 1,1,2-Trichloroethane | 310 ^b | 1,800 ^d | 0.02 | 0.3 | * |
| 79-01-6 | Trichloroethylene | 58 ^e | 5 ^e | 0.06 | 0.3 | * |
| 108-05-4 | Vinyl acetate | 78,000 ^b | 1,000 ^b | 170 ^b | 170 | * |
| 75-01-4 | Vinyl chloride | 0.46 ^e | 0.28 ^e | 0.01 ^f | 0.07 | * |
| 108-38-3 | m-Xylene | 160,000 ^b | 420 ^d | 210 | 210 | * |
| 95-47-6 | o-Xylene | 160,000 ^b | 410 ^d | 190 | 190 | * |
| 106-42-3 | p-Xylene | 160,000 ^b | 460 ^d | 200 | 200 | * |

| | | Exposure Route-Specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | |
|-----------|---------------------------|--|-----------------------|---|---------------------|----------------|
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | ADL (mg/kg) |
| 1330-20-7 | Xylenes (total) | 160,000 ^b | 320 ^d | 150 | 150 | * |
| | Ionizable Organics | | | | | |
| 65-85-0 | Benzoic Acid | 310,000 ^b | --- ^c | 400 ^{b,i} | 400 ⁱ | * |
| 95-57-8 | 2-Chlorophenol | 390 ^b | 53,000 ^d | 4 ^{b,i} | 4 ⁱ | * |
| 120-83-2 | 2,4-Dichlorophenol | 230 ^b | --- ^c | 1 ^{b,i} | 1 ⁱ | * |
| 51-28-5 | 2,4-Dinitrophenol | 160 ^b | --- ^c | 0.2 ^{b,f} | 0.2 | 3.3 |
| 88-85-7 | Dinoseb ^o | 78 ^b | --- ^c | 0.34 ^{b,i} | 3.4 ⁱ | * |
| 87-86-5 | Pentachlorophenol | 3 ^{e,j} | --- ^c | 0.03 ^{f,i} | 0.14 ⁱ | * |
| 93-72-1 | 2,4,5-TP (Silvex) | 630 ^b | --- ^c | 11 ⁱ | 55 ⁱ | * |
| 95-95-4 | 2,4,5-Trichlorophenol | 7,800 ^b | --- ^c | 270 ^{b,i} | 1,400 ⁱ | * |
| 88-06-2 | 2,4,6 Trichlorophenol | 58 ^e | 200 ^e | 0.2 ^{e,f,i} | 0.77 ⁱ | 0.66 |

| CAS No. | Chemical Name | Exposure Route-specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | ADL (mg/kg) |
|-------------------|---------------------------|--|----------------------|---|--------------------|-------------|
| | | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/L) | Class II (mg/L) | |
| Inorganics | | | | | | |
| 7440-36-0 | Antimony | 31 ^b | --- | 0.006 ^m | 0.024 ^m | * |
| 7440-38-2 | Arsenic ^{l,n} | t | 750 ^e | 0.05 ^m | 0.2 ^m | * |
| 7440-39-3 | Barium | 5,500 ^b | 690,000 ^b | 2.0 ^m | 2.0 ^m | * |
| 7440-41-7 | Beryllium | 160 ^b | 1,300 ^e | 0.004 ^m | 0.5 ^m | * |
| 7440-42-8 | Boron | 7,000 ^b | --- | 2.0 ^m | 2.0 ^m | * |
| 7440-43-9 | Cadmium ^{l,n} | 78 ^{b,r} | 1,800 ^e | 0.005 ^m | 0.05 ^m | * |
| 16887-00-6 | Chloride | --- | --- | 200 ^m | 200 ^m | * |
| 7440-47-3 | Chromium, total | 230 ^b | 270 ^e | 0.1 ^m | 1.0 ^m | * |
| 16065-83-1 | Chromium, ion, trivalent | 120,000 ^b | --- | --- | --- | * |
| 18540-29-9 | Chromium, ion, hexavalent | 230 ^b | 270 ^e | --- | --- | * |
| 7440-48-4 | Cobalt | 4,700 ^b | --- | 1.0 ^m | 1.0 ^m | * |

| | | Exposure Route-specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | |
|------------|---------------------------|--|---------------------|---|--------------------|-------------|
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/L) | Class II (mg/L) | ADL (mg/kg) |
| 7440-50-8 | Copper ⁿ | 2,900 ^b | --- ^c | 0.65 ^m | 0.65 ^m | * |
| 57-12-5 | Cyanide (amenable) | 1,600 ^b | --- ^c | 0.2 ^{q,m} | 0.6 ^{q,m} | * |
| 7782-41-4 | Fluoride | 4,700 ^b | --- ^c | 4.0 ^m | 4.0 ^m | * |
| 15438-31-0 | Iron | --- ^c | --- ^c | 5.0 ^m | 5.0 ^m | * |
| 7439-92-1 | Lead | 400 ^k | --- ^c | 0.0075 ^m | 0.1 ^m | * |
| 7439-96-5 | Manganese | 3,700 ^b | 69,000 ^b | 0.15 ^m | 10.0 ^m | * |
| 7439-97-6 | Mercury ^{l,n,s} | 23 ^b | 10 ^b | 0.002 ^m | 0.01 ^m | * |
| 7440-02-0 | Nickel ^l | 1,600 ^b | 13,000 ^e | 0.1 ^m | 2.0 ^m | * |
| 14797-55-8 | Nitrate as N ^p | 130,000 ^b | --- ^c | 10.0 ^q | 100 ^q | * |
| 7782-49-2 | Selenium ^{l,n} | 390 ^b | --- ^c | 0.05 ^m | 0.05 ^m | * |

| | | Exposure Route-specific Values for Soils | | Soil Component of the Groundwater Ingestion Exposure Route Values | | |
|------------|-------------------|--|--------------------|---|-------------------|-------------|
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/L) | Class II (mg/L) | ADL (mg/kg) |
| 7440-22-4 | Silver | 390 ^b | --- ^c | 0.05 ^m | --- | * |
| 14808-79-8 | Sulfate | --- ^c | --- ^c | 400 ^m | 400 ^m | * |
| 7440-28-0 | Thallium | 6.3 ^{b,u} | --- ^c | 0.002 ^m | 0.02 ^m | * |
| 7440-62-2 | Vanadium | 550 ^b | --- ^c | 0.049 ^m | 0.1 ^m | * |
| 7440-66-6 | Zinc ^l | 23,000 ^b | --- ^c | 5.0 ^m | 10 ^m | * |

"*" indicates that the ADL is less than or equal to the specified remediation objective.
NA means not available; no PQL or EQL available in USEPA analytical methods.

Chemical Name and Soil Remediation Objective Notations

- ^a Soil remediation objectives based on human health criteria only.
- ^b Calculated values correspond to a target hazard quotient of 1.
- ^c No toxicity criteria available for the route of exposure.
- ^d Soil saturation concentration (C_{sat}) = the concentration at which the absorptive limits of the soil particles, the solubility limits of the available soil moisture, and saturation of soil pore air have been reached. Above the soil saturation concentration, the assumptions regarding vapor transport to air and/or dissolved phase transport to groundwater (for chemicals which are liquid at ambient soil temperatures) have been violated, and alternative modeling approaches are required.
- ^e Calculated values correspond to a cancer risk level of 1 in 1,000,000.
- ^f Level is at or below Contract Laboratory Program required quantitation limit for Regular Analytical Services (RAS).
- ^g Chemical-specific properties are such that this route is not of concern at any soil contaminant concentration.
- ^h 40 CFR 761 contains applicability requirements and methodologies for the development of PCB remediation objectives. Requests for approval of a Tier 3 evaluation must address the applicability of 40 CFR 761.
- ⁱ Soil remediation objective for pH of 6.8. If soil pH is other than 6.8, refer to Appendix B, Tables C and D of this Part.
- ^j Ingestion soil remediation objective adjusted by a factor of 0.5 to account for dermal route.
- ^k A preliminary remediation goal of 400 mg/kg has been set for lead based on *Revised Interim Soil Lead Guidance for CERCLA Sites and RCRA Corrective Action Facilities*, OSWER Directive #9355.4-12.
- ^l Potential for soil-plant-human exposure.
- ^m The person conducting the remediation has the option to use: 1) TCLP or SPLP test results to compare with the remediation objectives listed in this Table; or 2) the total amount of contaminant in the soil sample results to compare with pH specific remediation objectives listed in Appendix B, Table C or D of this Part. (See Section 742.510.) If the person conducting the remediation wishes to calculate soil remediation objectives based on background concentrations, this should be done in accordance with Subpart D of this Part.
- ⁿ The Agency reserves the right to evaluate the potential for remaining contaminant concentrations to pose significant threats to crops, livestock, or wildlife.
- ^o For agrichemical facilities, remediation objectives for surficial soils which are based on field application rates may be more appropriate for currently registered pesticides. Consult the Agency for further information.
- ^p For agrichemical facilities, soil remediation objectives based on site-specific background concentrations of Nitrate as N may be more appropriate. Such determinations shall be conducted in accordance with the procedures set forth in Subparts D and I of this Part.
- ^q The TCLP extraction must be done using water at a pH of 7.0.
- ^r Value based on dietary Reference Dose.
- ^s Value for Ingestion based on Reference Dose for Mercuric chloride (CAS No. 7487-94-7); value for Inhalation based on Reference Concentration for elemental Mercury (CAS No. 7439-97-6).
- ^t For the ingestion route for arsenic, see 742.Appendix A, Table G.
- ^u Value based on Reference Dose for Thallium sulfate (CAS No. 7446-18-6).

(Source: Amended at 26 Ill. Reg. ___, effective ____)

Section 742.APPENDIX B: Tier 1 Tables and Illustrations

Section 742.Table B: Tier 1 Soil Remediation Objectives^a for Industrial/Commercial Properties

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route | | Values |
|------------|-----------------------|--|----------------------|----------------------|----------------------|--|--------------------|--------|
| | | Industrial-Commercial | Construction Worker | Ingestion (mg/kg) | Inhalation (mg/kg) | Ingestion (mg/kg) | Inhalation (mg/kg) | |
| 83-32-9 | Acenaphthene | 120,000 ^b | ----- ^c | 120,000 ^b | ----- ^c | 570 ^b | 2,900 | * |
| 67-64-1 | Acetone | 200,000 ^b | 100,000 ^d | 200,000 ^b | 100,000 ^d | 16 ^b | 16 | * |
| 15972-60-8 | Alachlor ^o | 72 ^e | ----- ^c | 1,600 ^e | ----- ^c | 0.04 | 0.2 | NA |
| 116-06-3 | Aldicarb ^o | 2,000 ^b | ----- ^c | 200 ^b | ----- ^c | 0.013 | 0.07 | NA |
| 309-00-2 | Aldrin | 0.3 ^e | 6.6 ^e | 6.1 ^b | 9.3 ^e | 0.5 ^e | 2.5 | 0.94 |
| 120-12-7 | Anthracene | 610,000 ^b | ----- ^c | 610,000 ^b | ----- ^c | 12,000 ^b | 59,000 | * |
| 1912-24-9 | Atrazine ^o | 72,000 ^b | ----- ^c | 7,100 ^b | ----- ^c | 0.066 | 0.33 | NA |
| 71-43-2 | Benzene | 100 ^e | 1.6 ^e | 2,300 ^e | 2.2 ^e | 0.03 | 0.17 | * |

| | | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route | | |
|----------|--|--|---------------------|----------------------|---------------------|--|---------------------|-------------|
| | | Industrial-Commercial | | Construction Worker | | Values | | |
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | ADL (mg/kg) |
| 56-55-3 | Benzo(<i>a</i>)anthracene | 8 ^e | ---- ^c | 170 ^e | ---- ^c | 2 | 8 | * |
| 205-99-2 | Benzo(<i>b</i>)fluoranthene | 8 ^e | ---- ^c | 170 ^e | ---- ^c | 5 | 25 | * |
| 207-08-9 | Benzo(<i>k</i>)fluroanthene | 78 ^e | ---- ^c | 1,700 ^e | ---- ^c | 49 | 250 | * |
| 50-32-8 | Benzo(<i>a</i>)pyrene | 0.8 ^e | ---- ^c | 17 ^e | ---- ^c | 8 | 82 | * |
| 111-44-4 | Bis(2-chloroethyl)ether | 5 ^e | 0.47 ^e | 75 ^e | 0.66 ^e | 0.0004 ^{e,f} | 0.0004 | 0.66 |
| 117-81-7 | Bis(2-ethylhexyl)phthalate | 410 ^e | 31,000 ^d | 4,100 ^b | 31,000 ^d | 3,600 | 31,000 ^d | * |
| 75-27-4 | Bromodichloromethane (Dichlorobromomethane) | 92 ^e | 3,000 ^d | 2,000 ^e | 3,000 ^d | 0.6 | 0.6 | * |
| 75-25-2 | Bromoform | 720 ^e | 100 ^e | 16,000 ^e | 140 ^e | 0.8 | 0.8 | * |
| 71-36-3 | Butanol | 200,000 ^b | 10,000 ^d | 200,000 ^b | 10,000 ^d | 17 ^b | 17 | NA |
| 85-68-7 | Butyl benzyl phthalate | 410,000 ^b | 930 ^d | 410,000 ^b | 930 ^d | 930 ^d | 930 ^d | * |
| 86-74-8 | Carbazole | 290 ^e | ---- ^c | 6,200 ^e | ---- ^c | 0.6 ^e | 2.8 | NA |

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route | Values | |
|-----------|---|--|---------------------|---------------------|--------------------|--|--------|----|
| | | Industrial-Commercial | Construction Worker | Ingestion (mg/kg) | Inhalation (mg/kg) | | | |
| 1563-66-2 | Carbofuran ^o | 10,000 ^b | ---- ^c | 1,000 ^b | ---- ^c | 0.22 | 1.1 | NA |
| 75-15-0 | Carbon disulfide | 200,000 ^b | 720 ^d | 20,000 ^b | 9.0 ^b | 32 ^b | 160 | * |
| 56-23-5 | Carbon tetrachloride | 44 ^e | 0.64 ^e | 410 ^b | 0.90 ^e | 0.07 | 0.33 | * |
| 57-74-9 | Chlordane | 1.6 ^e | 140 ^e | 100 ^b | 22 ^b | 10 | 48 | * |
| 106-47-8 | 4 - Chloroaniline (<i>p</i> -Chloroaniline) | 8,200 ^b | ---- ^c | 820 ^b | ---- ^c | 0.7 ^b | 0.7 | * |
| 108-90-7 | Chlorobenzene (Monochlorobenzene) | 41,000 ^b | 210 ^b | 4,100 ^b | 1.3 ^b | 1 | 6.5 | * |
| 124-48-1 | Chlorodibromomethane (Dibromochloromethane) | 41,000 ^b | 1,300 ^d | 41,000 ^b | 1,300 ^d | 0.4 | 0.4 | * |
| 67-66-3 | Chloroform | 940 ^e | 0.54 ^e | 2,000 ^b | 0.76 ^e | 0.6 | 2.9 | * |
| 218-01-9 | Chrysene | 780 ^e | ---- ^c | 17,000 ^e | ---- ^e | 160 | 800 | * |
| 94-75-7 | 2,4-D ^o | 20,000 ^b | ---- ^c | 2,000 ^b | ---- ^c | 1.5 | 7.7 | * |

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| | | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route | Values | |
|----------|--|--|---------------------|----------------------|--------------------|--|--------------------|-------------|
| | | Industrial-Commercial | Construction Worker | Ingestion (mg/kg) | Inhalation (mg/kg) | | | |
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | ADL (mg/kg) |
| 75-99-0 | Dalapon ^o | 61,000 ^b | ----- ^c | 6,100 ^b | ----- ^c | 0.85 | 8.5 | * |
| 72-54-8 | DDD | 24 ^e | ----- ^c | 520 ^e | ----- ^c | 16 ^e | 80 | * |
| 72-55-9 | DDE | 17 ^e | ----- ^c | 370 ^e | ----- ^c | 54 ^e | 270 | * |
| 50-29-3 | DDT | 17 ^e | 1,500 ^e | 100 ^b | 2,100 ^e | 32 ^e | 160 | * |
| 53-70-3 | Dibenzo(<i>a,h</i>)anthracene | 0.8 ^e | ----- ^c | 17 ^e | ----- ^c | 2 | 7.6 | * |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 4 ^e | 17 ^b | 89 ^e | 0.11 ^b | 0.002 | 0.002 | * |
| 106-93-4 | 1,2-Dibromoethane (Ethylene dibromide) | 0.07 ^e | 0.32 ^e | 1.5 ^e | 0.45 ^e | 0.0004 | 0.004 | 0.005 |
| 84-74-2 | Di- <i>n</i> -butyl phthalate | 200,000 ^b | 2,300 ^d | 200,000 ^b | 2,300 ^d | 2,300 ^d | 2,300 ^d | * |
| 95-50-1 | 1,2-Dichlorobenzene (<i>o</i> - Dichlorobenzene) | 180,000 ^b | 560 ^d | 18,000 ^b | 310 ^b | 17 | 43 | * |
| 106-46-7 | 1,4-Dichlorobenzene (<i>p</i> - Dichlorobenzene) | ----- ^c | 17,000 ^b | ----- ^c | 340 ^b | 2 | 11 | * |

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| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route Values | | |
|----------|--|--|---------------------|------------------------|--------------------|---|-------|-------|
| | | Industrial-Commercial | Construction Worker | Ingestion (mg/kg) | Inhalation (mg/kg) | | | |
| 91-94-1 | 3,3'-Dichlorobenzidine | 13 ^e | ----- ^c | 280 ^e | ----- ^c | 0.007 ^{e,f} | 0.033 | 1.3 |
| 75-34-3 | 1,1-Dichloroethane | 200,000 ^b | 1,700 ^d | 200,000 ^b | 130 ^b | 23 ^b | 110 | * |
| 107-06-2 | 1,2-Dichloroethane (Ethylene dichloride) | 63 ^e | 0.70 ^e | 1,400 ^e | 0.99 ^e | 0.02 | 0.1 | * |
| 75-35-4 | 1,1-Dichloroethylene | 18,000 ^b | 1,500 ^d | 1,800 ^b | 300 ^v | 0.06 | 0.3 | * |
| 156-59-2 | <i>cis</i> -1,2-Dichloroethylene | 20,000 ^b | 1,200 ^d | 20,000 ^b | 1,200 ^d | 0.4 | 1.1 | * |
| 156-60-5 | <i>Trans</i> -1,2-Dichloroethylene | 41,000 ^b | 3,100 ^d | 41,000 ^b | 3,100 ^d | 0.7 | 3.4 | * |
| 78-87-5 | 1,2-Dichloropropane | 84 ^e | 23 ^b | 1,800 ^e | 0.50 ^b | 0.03 | 0.15 | * |
| 542-75-6 | 1,3-Dichloropropene (1,3-Dichloropropylene, <i>cis</i> + <i>trans</i>) | 57 ^e | 2.1 ^e | 1,200 ^e | 0.39 ^b | 0.004 ^e | 0.02 | 0.005 |
| 60-57-1 | Dieldrin ⁿ | 0.4 ^e | 2.2 ^e | 7.8 ^e | 3.1 ^e | 0.004 ^e | 0.02 | 0.603 |
| 84-66-2 | Diethyl phthalate | 1,000,000 ^b | 2,000 ^d | 1,000,000 ^b | 2,000 ^d | 470 ^b | 470 | * |

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route | | Values |
|----------|-------------------------------|--|---------------------|---------------------|---------------------|--|---------------------|--------|
| | | Industrial-Commercial | Construction Worker | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | |
| 105-67-9 | 2,4-Dimethylphenol | 41,000 ^b | ---- ^c | 41,000 ^b | ---- ^c | 9 ^b | 9 | * |
| 121-14-2 | 2,4-Dinitrotoluene | 8.4 ^e | ---- ^c | 180 ^e | ---- ^c | 0.0008 ^{e,f} | 0.0008 | 0.250 |
| 606-20-2 | 2,6-Dinitrotoluene | 8.4 ^e | ---- ^c | 180 ^e | ---- ^c | 0.0007 ^{e,f} | 0.0007 | 0.260 |
| 117-84-0 | Di- <i>n</i> -octyl phthalate | 41,000 ^e | 10,000 ^d | 4,100 ^b | 10,000 ^d | 10,000 ^d | 10,000 ^d | * |
| 115-29-7 | Endosulfan ^g | 12,000 ^b | ---- ^c | 1,200 ^b | ---- ^c | 18 ^b | 90 | * |
| 145-73-3 | Endothall ^h | 41,000 ^c | ---- ^c | 4,100 ^b | ---- ^c | 0.4 | 0.4 | NA |
| 72-20-8 | Endrin | 610 ^b | ---- ^c | 61 ^b | ---- ^c | 1 | 5 | * |
| 100-41-4 | Ethylbenzene | 200,000 ^b | 400 ^d | 20,000 ^b | 58 ^b | 13 | 19 | * |
| 206-44-0 | Fluoranthene | 82,000 ^b | ---- ^c | 82,000 ^b | ---- ^c | 4,300 ^b | 21,000 | * |
| 86-73-7 | Fluorene | 82,000 ^b | ---- ^c | 82,000 ^b | ---- ^c | 560 ^b | 2,800 | * |
| 76-44-8 | Heptachlor | 1 ^e | 11 ^e | 28 ^e | 16 ^e | 23 | 110 | * |

f

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route | | Values |
|-----------|--|--|---------------------|----------------------|--------------------|--|--------------------|--------|
| | | Industrial-Commercial | Construction Worker | Ingestion (mg/kg) | Inhalation (mg/kg) | | | |
| 1024-57-3 | Heptachlor epoxide | 0.6 ^e | 9.2 ^e | 2.7 ^b | 13 ^e | 0.7 | 3.3 | 1.005 |
| 118-74-1 | Hexachlorobenzene | 4 ^e | 1.8 ^e | 78 ^e | 2.6 ^e | 2 | 11 | * |
| 319-84-6 | <i>Alpha</i> -HCH (<i>alpha</i> -BHC) | 0.9 ^e | 1.5 ^e | 20 ^e | 2.1 ^e | 0.0005 ^{e,f} | 0.003 | 0.0074 |
| 58-89-9 | <i>Gamma</i> -HCH (Lindane) ^g | 4 ^e | ---- ^c | 96 ^e | ---- ^c | 0.009 | 0.047 | * |
| 77-47-4 | Hexachlorocyclopentadiene | 14,000 ^b | 16 ^b | 14,000 ^b | 1.1 ^b | 400 | 2,200 ^d | * |
| 67-72-1 | Hexachloroethane | 2,000 ^b | ---- ^c | 2,000 ^b | ---- ^c | 0.5 ^b | 2.6 | * |
| 193-39-5 | Indeno(1,2,3- <i>c,d</i>)pyrene | 8 ^e | ---- ^c | 170 ^e | ---- ^c | 14 | 69 | * |
| 78-59-1 | Isophorone | 410,000 ^b | 4,600 ^d | 410,000 ^b | 4,600 ^d | 8 ^b | 8 | * |
| 72-43-5 | Methoxychlor ^o | 10,000 ^b | ---- ^c | 1,000 ^b | ---- ^c | 160 | 780 | * |
| 74-83-9 | Methyl bromide (Bromomethane) | 2,900 ^b | 15 ^b | 1,000 ^b | 3.9 ^b | 0.2 ^b | 1.2 | * |

| | | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route | Values | |
|-----------|--|--|---------------------|----------------------|---------------------|--|-------------------|-------------|
| | | Industrial-Commercial | Construction Worker | Ingestion (mg/kg) | Inhalation (mg/kg) | | | |
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | ADL (mg/kg) |
| 1634-04-4 | Methyl tert-butyl ether Methyl tertiary-butyl ether | 20,000 ^b | 8,800 ^d | 2,000 ^b | 140 ^b | 0.32 | 0.32 | * |
| 75-09-2 | Methylene chloride (Dichloromethane) | 760 ^e | 24 ^e | 12,000 ^b | 34 ^e | 0.02 ^e | 0.2 | * |
| 95-48-7 | 2-Methylphenol (o - Cresol) | 100,000 ^b | ---- ^c | 100,000 ^b | ---- ^c | 15 ^b | 15 | * |
| 86-30-6 | N-Nitrosodiphenylamine | 1,200 ^e | ---- ^c | 25,000 ^e | ---- ^c | 1 ^e | 5.6 | * |
| 621-64-7 | N-Nitrosodi-n-propylamine | 0.8 ^e | ---- ^c | 18 ^e | ---- ^c | 0.00005 ^{e,f} | 0.00005 | 0.0018 |
| 91-20-3 | Naphthalene | 41,000 ^b | 270 ^b | 4,100 ^b | 1.8 ^b | 12 ^b | 18 | * |
| 98-95-3 | Nitrobenzene | 1,000 ^b | 140 ^b | 1,000 ^b | 9.4 ^b | 0.1 ^{b,f} | 0.1 | 0.26 |
| 108-95-2 | Phenol | 1,000,000 ^b | ---- ^c | 120,000 ^b | ---- ^c | 100 ^b | 100 | * |
| 1918-02-1 | Picloram ^o | 140,000 ^b | ---- ^c | 14,000 ^b | ---- ^c | 2 | 20 | NA |
| 1336-36-3 | Polychlorinated biphenyls (PCBs) ⁿ | 1 ^h | ---- ^{c,h} | 1 ^h | ---- ^{c,h} | ---- ^h | ---- ^h | * |
| 129-00-0 | Pyrene | 61,000 ^b | ---- ^c | 61,000 ^b | ---- ^c | 4,200 ^b | 21,000 | * |

| | | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route | Values | |
|-----------|--|--|---------------------|----------------------|--------------------|--|------------------|-------------|
| | | Industrial-Commercial | Construction Worker | Ingestion (mg/kg) | Inhalation (mg/kg) | | | |
| CAS No. | Chemical Name | Ingestion (mg/kg) | Inhalation (mg/kg) | Ingestion (mg/kg) | Inhalation (mg/kg) | Class I (mg/kg) | Class II (mg/kg) | ADL (mg/kg) |
| 122-34-9 | Simazine ^o | 10,000 ^b | ----- ^c | 1,000 ^b | ----- ^c | 0.04 | 0.37 | NA |
| 100-42-5 | Styrene | 410,000 ^b | 1,500 ^d | 41,000 ^b | 430 ^b | 4 | 18 | * |
| 127-18-4 | Tetrachloroethylene (Perchloroethylene) | 110 ^e | 20 ^e | 2,400 ^e | 28 ^e | 0.06 | 0.3 | * |
| 108-88-3 | Toluene | 410,000 ^b | 650 ^d | 410,000 ^b | 42 ^b | 12 | 29 | * |
| 8001-35-2 | Toxaphene ⁿ | 5.2 ^e | 170 ^e | 110 ^e | 240 ^e | 31 | 150 | * |
| 120-82-1 | 1,2,4-Trichlorobenzene | 20,000 ^b | 3,200 ^d | 2,000 ^b | 920 ^b | 5 | 53 | * |
| 71-55-6 | 1,1,1-Trichloroethane | ----- ^c | 1,200 ^d | ----- ^c | 1,200 ^d | 2 | 9.6 | * |
| 79-00-5 | 1,1,2-Trichloroethane | 8,200 ^b | 1,800 ^d | 8,200 ^b | 1,800 ^d | 0.02 | 0.3 | * |
| 79-01-6 | Trichloroethylene | 520 ^e | 8.9 ^e | 1,200 ^b | 12 ^e | 0.06 | 0.3 | * |
| 108-05-4 | Vinyl acetate | 1,000,000 ^b | 1,600 ^b | 200,000 ^b | 10 ^b | 170 ^b | 170 | * |

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route | Values | |
|-----------|---------------------------|--|---------------------|----------------------|---------------------|--|------------------|-----|
| | | Industrial-Commercial | Construction Worker | Ingestion (mg/kg) | Inhalation (mg/kg) | | | |
| 75-01-4 | Vinyl chloride | 7.9 ^e | 1.1 ^e | 170 ^e | 1.1 ^b | 0.01 ^f | 0.07 | * |
| 108-38-3 | m-Xylene | 1,000,000 | 420 ^d | 410,000 ^b | 420 ^d | 210 | 210 | * |
| 95-47-6 | o-Xylene | 1,000,000 | 410 ^d | 410,000 ^b | 410 ^d | 190 | 190 | * |
| 106-42-3 | p-Xylene | 1,000,000 | 460 ^d | 410,000 ^b | 460 ^d | 200 | 200 | * |
| 1330-20-7 | Xylenes (total) | 1,000,000 ^b | 320 ^d | 410,000 ^b | 320 ^d | 150 | 150 | * |
| | Ionizable Organics | | | | | | | |
| 65-85-0 | Benzoic Acid | 1,000,000 ^b | ---- ^c | 820,000 ^b | ---- ^c | 400 ^{b,i} | 400 ⁱ | * |
| 95-57-8 | 2-Chlorophenol | 10,000 ^b | 53,000 ^d | 10,000 ^b | 53,000 ^d | 4 ^{b,i} | 20 ⁱ | * |
| 120-83-2 | 2,4-Dichlorophenol | 6,100 ^b | ---- ^c | 610 ^b | ---- ^c | 1 ^{b,i} | 1 ⁱ | * |
| 51-28-5 | 2,4-Dinitrophenol | 4,100 ^b | ---- ^c | 410 ^b | ---- ^c | 0.2 ^{b,f,i} | 0.2 ⁱ | 3.3 |
| 88-85-7 | Dinoseb ^o | 2,000 ^b | ---- ^c | 200 ^b | ---- ^c | 0.34 ^{b,i} | 3.4 ⁱ | * |

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route | | Values |
|---------|------------------------|--|---------------------|----------------------|--------------------|--|--------------------|--------|
| | | Industrial-Commercial | Construction Worker | Ingestion (mg/kg) | Inhalation (mg/kg) | | | |
| 87-86-5 | Pentachlorophenol | 24 ^{e,j} | ---- ^c | 520 ^{e,j} | ---- ^c | 0.03 ^{f,i} | 0.14 ⁱ | * |
| 93-72-1 | 2,4,5-TP (Silvex) | 16,000 ^b | ---- ^c | 1,600 ^b | ---- ^c | 11 ⁱ | 55 ⁱ | * |
| 95-95-4 | 2,4,5-Trichlorophenol | 200,000 ^b | ---- ^c | 200,000 ^b | ---- ^c | 270 ^{b,i} | 1,400 ⁱ | * |
| 88-06-2 | 2,4,6- Trichlorophenol | 520 ^e | 390 ^e | 11,000 ^e | 540 ^e | 0.2 ^{e,f,i} | 0.77 ⁱ | 0.66 |

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route Values | | |
|------------|---------------------------|--|----------------------|----------------------|----------------------|---|--------------------|---|
| | | Industrial-Commercial | Construction Worker | Ingestion (mg/kg) | Inhalation (mg/kg) | | | |
| | Inorganics | | | | | | | |
| 7440-36-0 | Antimony | 820 ^b | ---- ^c | 82 ^b | ---- ^c | 0.006 ^m | 0.024 ^m | * |
| 7440-38-2 | Arsenic ^{l,n} | t | 1,200 ^e | 61 ^b | 25,000 ^e | 0.05 ^m | 0.2 ^m | * |
| 7440-39-3 | Barium | 140,000 ^b | 910,000 ^b | 14,000 ^b | 870,000 ^b | 2.0 ^m | 2.0 ^m | * |
| 7440-41-7 | Beryllium | 4,100 ^b | 2,100 ^e | 410 ^b | 44,000 ^e | 0.004 ^m | 0.5 ^m | * |
| 7440-42-8 | Boron | 180,000 ^b | 1,000,000 | 18,000 ^b | 1,000,000 | 2.0 ^m | 2.0 ^m | * |
| 7440-43-9 | Cadmium ^{l,n} | 2,000 ^{b,r} | 2,800 ^e | 200 ^{b,r} | 59,000 ^e | 0.005 ^m | 0.05 ^m | * |
| 16887-00-6 | Chloride | ----- ^c | ----- ^c | ----- ^c | ----- ^c | 200 ^m | 200 ^m | * |
| 7440-47-3 | Chromium, total | 6,100 ^b | 420 ^e | 4,100 ^b | 690 ^b | 0.1 ^m | 1.0 ^m | * |
| 16065-83-1 | Chromium, ion, trivalent | 1,000,000 ^b | ---- ^c | 310,000 ^b | ---- ^c | ---- ^g | ---- ^g | * |
| 18540-29-9 | Chromium, ion, hexavalent | 6,100 ^b | 420 ^e | 4,100 ^b | 690 ^b | ---- | ---- | * |

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route Values | | | |
|------------|---------------------------|--|----------------------|----------------------|----------------------|---|--------------------|---|--|
| | | Industrial-Commercial | | Construction Worker | | | | | |
| | | Ingestion (mg/kg) | Inhalation (mg/kg) | Ingestion (mg/kg) | Inhalation (mg/kg) | | | | |
| 7440-48-4 | Cobalt | 120,000 ^b | ---- ^c | 12,000 ^b | ---- ^c | 1.0 ^m | 1.0 ^m | * | |
| 7440-50-8 | Copper ⁿ | 82,000 ^b | ---- ^c | 8,200 ^b | ---- ^c | 0.65 ^m | 0.65 ^m | * | |
| 57-12-5 | Cyanide (amenable) | 41,000 ^b | ---- ^c | 4,100 ^b | ---- ^c | 0.2 ^{q,m} | 0.6 ^{q,m} | * | |
| 7782-41-4 | Fluoride | 120,000 ^b | ---- ^c | 12,000 ^b | ---- ^c | 4.0 ^m | 4.0 ^m | * | |
| 15438-31-0 | Iron | ---- ^c | ---- ^c | ---- ^c | ---- ^c | 5.0 ^m | 5.0 ^m | * | |
| 7439-92-1 | Lead | 400 ^k | ---- ^c | 400 ^k | ---- ^c | 0.0075 ^m | 0.1 ^m | * | |
| 7439-96-5 | Manganese | 96,000 ^b | 91,000 ^b | 9,600 ^b | 8,700 ^b | 0.15 ^m | 10.0 ^m | * | |
| 7439-97-6 | Mercury ^{l,n,s} | 610 ^b | 540,000 ^b | 61 ^b | 52,000 ^b | 0.002 ^m | 0.01 ^m | * | |
| 7440-02-0 | Nickel ^l | 41,000 ^b | 21,000 ^e | 4,100 ^b | 440,000 ^e | 0.1 ^m | 2.0 ^m | * | |
| 14797-55-8 | Nitrate as N ^p | 1,000,000 ^b | ---- ^c | 330,000 ^b | ---- ^c | 10.0 ^q | 100 ^q | * | |
| 7782-49-2 | Selenium ^{l,n} | 10,000 ^b | ---- ^c | 1,000 ^b | ---- ^c | 0.05 ^m | 0.05 ^m | * | |

| CAS No. | Chemical Name | Exposure Route-Specific Values for Soils | | | | Soil Component of the Groundwater Ingestion Exposure Route Values | | | |
|------------|-------------------|--|--------------------|---------------------|--------------------|---|-------------------|---|--|
| | | Industrial-Commercial | | Construction Worker | | | | | |
| | | Ingestion (mg/kg) | Inhalation (mg/kg) | Ingestion (mg/kg) | Inhalation (mg/kg) | | | | |
| 7440-22-4 | Silver | 10,000 ^b * | ---- ^c | 1,000 ^b | ---- ^c | 0.05 ^m | ---- | * | |
| 14808-79-8 | Sulfate | ---- ^c | ---- ^c | ---- ^c | ---- ^c | 400 ^m | 400 ^m | * | |
| 7440-28-0 | Thallium | 160 ^{b,u} | ---- ^c | 160 ^{b,u} | ---- ^c | 0.002 ^m | 0.02 ^m | * | |
| 7440-62-2 | Vanadium | 14,000 ^b | ---- ^c | 1,400 ^b | ---- ^c | 0.049 ^m | 0.1 ^m | * | |
| 7440-66-6 | Zinc ^j | 610,000 ^b | ---- ^c | 61,000 ^b | ---- ^c | 5.0 ^m | 10 ^m | * | |

"*" indicates that the ADL is less than or equal to the specified remediation objective.

NA means Not Available; no PQL or EQL available in USEPA analytical methods.

Chemical Name and Soil Remediation Objective Notations (2nd, 5th thru 8th Columns)

- ^a Soil remediation objectives based on human health criteria only.
- ^b Calculated values correspond to a target hazard quotient of 1.
- ^c No toxicity criteria available for this route of exposure.
- ^d Soil saturation concentration ($C_{[sat]}$) = the concentration at which the absorptive limits of the soil particles, the solubility limits of the available soil moisture, and saturation of soil pore air have been reached. Above the soil saturation concentration, the assumptions regarding vapor transport to air and/or dissolved phase transport to groundwater (for chemicals which are liquid at ambient soil temperatures) have been violated, and alternative modeling approaches are required.
- ^e Calculated values correspond to a cancer risk level of 1 in 1,000,000.
- ^f Level is at or below Contract Laboratory Program required quantitation limit for Regular Analytical Services (RAS).
- ^g Chemical-specific properties are such that this route is not of concern at any soil contaminant concentration.
- ^h 40 CFR 761 contains applicability requirements and methodologies for the development of PCB remediation objectives. Requests for approval of a Tier 3 evaluation must address the applicability of 40 CFR 761.
- ⁱ Soil remediation objective for pH of 6.8. If soil pH is other than 6.8, refer to Appendix B, Tables C and D in this Part.
- ^j Ingestion soil remediation objective adjusted by a factor of 0.5 to account for dermal route.
- ^k A preliminary remediation goal of 400 mg/kg has been set for lead based on *Revised Interim Soil Lead Guidance for CERCLA Sites and RCRA Corrective Action Facilities*, OSWER Directive #9355.4-12.
- ^l Potential for soil-plant-human exposure.
- ^m The person conducting the remediation has the option to use: (1) TCLP or SPLP test results to compare with the remediation objectives listed in this Table; or (2) the total amount of contaminant in the soil sample results to compare with pH specific remediation objectives listed in Appendix B, Table C or D of this Part. (See Section 742.510.) If the person conducting the remediation wishes to calculate soil remediation objectives based on background concentrations, this should be done in accordance with Subpart D of this Part.
- ⁿ The Agency reserves the right to evaluate the potential for remaining contaminant concentrations to pose significant threats to crops, livestock, or wildlife.
- ^o For agrichemical facilities, remediation objectives for surficial soils which are based on field application rates may be more appropriate for currently registered pesticides. Consult the Agency for further information.
- ^p For agrichemical facilities, soil remediation objectives based on site-specific background concentrations of Nitrate as N may be more appropriate. Such determinations shall be conducted in accordance with the procedures set forth in Subparts D and I of this Part.
- ^q The TCLP extraction must be done using water at a pH of 7.0.
- ^r Value based on dietary Reference Dose.
- ^s Value for Ingestion based on Reference Dose for Mercuric chloride (CAS No. 7487-94-7); value for Inhalation based on Reference Concentration for elemental Mercury (CAS No. 7439-97-6).
- ^t For the ingestion route for arsenic for industrial/commercial, see 742.Appendix A, Table G.
- ^u Value based on Reference Dose for Thallium sulfate (CAS No. 7446-18-6).
- ^v Calculated values correspond to soil concentrations that should not result in air concentrations that exceed criteria for workplace air.

(Source: Amended at 26 Ill. Reg. ____, effective ____)

Section 742.APPENDIX B: Tier 1 Tables and Illustrations

Section 742.TABLE E: Tier 1 Groundwater Remediation Objectives for the Groundwater Component of the Groundwater Ingestion Route

| | | Groundwater Remediation Objective | |
|------------|--|-----------------------------------|--------------------|
| CAS No. | Chemical Name | Class I (mg/L) | Class II (mg/L) |
| 83-32-9 | Acenaphthene | 0.42 | 2.1 |
| 67-64-1 | Acetone | 0.7 | 0.7 |
| 15972-60-8 | Alachlor | 0.002 ^c | 0.01 ^c |
| 116-06-3 | Aldicarb | 0.003 ^c | 0.015 ^c |
| 309-00-2 | Aldrin | 0.014 ^a | 0.07 |
| 120-12-7 | Anthracene | 2.1 | 10.5 |
| 1912-24-9 | Atrazine | 0.003 ^c | 0.015 ^c |
| 71-43-2 | Benzene | 0.005 ^c | 0.025 ^c |
| 56-55-3 | Benzo(<i>a</i>)anthracene | 0.00013 ^a | 0.00065 |
| 205-99-2 | Benzo(<i>b</i>)fluoranthene | 0.00018 ^a | 0.0009 |
| 207-08-9 | Benzo(<i>k</i>)fluroanthene | 0.00017 ^a | 0.00085 |
| 50-32-8 | Benzo(<i>a</i>)pyrene | 0.0002 ^{a,c} | 0.002 ^c |
| 111-44-4 | Bis(2-chloroethyl)ether | 0.01 ^a | 0.01 |
| 117-81-7 | Bis(2-ethylhexyl)phthalate (Di(2-ethylhexyl)phthalate) | 0.006 ^c | 0.06 ^c |
| 75-27-4 | Bromodichloromethane (Dichlorobromomethane) | 0.0002 ^a | 0.0002 |
| 75-25-2 | Bromoform | 0.001 ^a | 0.001 |
| 71-36-3 | Butanol | 0.7 | 0.7 |
| 85-68-7 | Butyl benzyl phthalate | 1.4 | 7.0 |
| 86-74-8 | Carbazole | --- | --- |
| 1563-66-2 | Carbofuran | 0.04 ^c | 0.2 ^c |
| 75-15-0 | Carbon disulfide | 0.7 | 3.5 |
| 56-23-5 | Carbon tetrachloride | 0.005 ^c | 0.025 ^c |
| 57-74-9 | Chlordane | 0.002 ^c | 0.01 ^c |

| Groundwater Remediation Objective | | | |
|-----------------------------------|--|----------------------|---------------------|
| CAS No. | Chemical Name | Class I (mg/L) | Class II (mg/L) |
| 108-90-7 | Chlorobenzene (Monochlorobenzene) | 0.1 ^c | 0.5 ^c |
| 124-48-1 | Chlorodibromomethane (Dibromochloromethane) | 0.14 | 0.14 |
| 67-66-3 | Chloroform | 0.0002 ^a | 0.001 |
| 218-01-9 | Chrysene | 0.0015 ^a | 0.0075 |
| 94-75-7 | 2,4-D | 0.07 ^c | 0.35 ^c |
| 75-99-0 | Dalapon | 0.2 ^c | 2.0 ^c |
| 72-54-8 | DDD | 0.014 ^a | 0.07 |
| 72-55-9 | DDE | 0.01 ^a | 0.05 |
| 50-29-3 | DDT | 0.006 ^a | 0.03 |
| 53-70-3 | Dibenzo(<i>a,h</i>)anthracene | 0.0003 ^a | 0.0015 |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 0.0002 ^c | 0.0002 ^c |
| 106-93-4 | 1,2-Dibromoethane (Ethylene dibromide) | 0.00005 ^c | 0.0005 ^c |
| 84-74-2 | Di- <i>n</i> -butyl phthalate | 0.7 | 3.5 |
| 95-50-1 | 1,2-Dichlorobenzene (<i>o</i> – Dichlorobenzene) | 0.6 ^c | 1.5 ^c |
| 106-46-7 | 1,4-Dichlorobenzene (<i>p</i> – Dichlorobenzene) | 0.075 ^c | 0.375 ^c |
| 91-94-1 | 3,3'-Dichlorobenzidine | 0.02 ^a | 0.1 |
| 75-34-3 | 1,1-Dichloroethane | 0.7 | 3.5 |
| 107-06-2 | 1,2-Dichloroethane (Ethylene dichloride) | 0.005 ^c | 0.025 ^c |
| 75-35-4 | 1,1-Dichloroethylene ^b | 0.007 ^c | 0.035 ^c |
| 156-59-2 | <i>cis</i> -1,2-Dichloroethylene | 0.07 ^c | 0.2 ^c |
| 156-60-5 | <i>trans</i> -1,2-Dichloroethylene | 0.1 ^c | 0.5 ^c |
| 78-87-5 | 1,2-Dichloropropane | 0.005 ^c | 0.025 ^c |
| 542-75-6 | 1,3-Dichloropropene (1,3-Dichloropropylene, <i>cis</i> + <i>trans</i>) | 0.001 ^a | 0.005 |

| Groundwater Remediation Objective | | | |
|-----------------------------------|--|----------------------|--------------------|
| CAS No. | Chemical Name | Class I (mg/L) | Class II (mg/L) |
| 60-57-1 | Dieldrin | 0.009 ^a | 0.045 |
| 84-66-2 | Diethyl phthalate | 5.6 | 5.6 |
| 121-14-2 | 2,4-Dinitrotoluene ^a | 0.00002 ^a | 0.00002 |
| 606-20-2 | 2,6-Dinitrotoluene ^a | 0.00031 ^a | 0.00031 |
| 88-85-7 | Dinoseb | 0.007 ^c | 0.07 ^c |
| 117-84-0 | Di-n-octyl phthalate | 0.14 | 0.7 |
| 115-29-7 | Endosulfan | 0.042 | 0.21 |
| 145-73-3 | Endothall | 0.1 ^c | 0.1 ^c |
| 72-20-8 | Endrin | 0.002 ^c | 0.01 ^c |
| 100-41-4 | Ethylbenzene | 0.7 ^c | 1.0 ^c |
| 206-44-0 | Fluoranthene | 0.28 | 1.4 |
| 86-73-7 | Fluorene | 0.28 | 1.4 |
| 76-44-8 | Heptachlor | 0.0004 ^c | 0.002 ^c |
| 1024-57-3 | Heptachlor epoxide | 0.0002 ^c | 0.001 ^c |
| 118-74-1 | Hexachlorobenzene | 0.00006 ^a | 0.0003 |
| 319-84-6 | <i>alpha</i> -HCH (<i>alpha</i> -BHC) | 0.00011 ^a | 0.00055 |
| 58-89-9 | <i>Gamma</i> -HCH (Lindane) | 0.0002 ^c | 0.001 ^c |
| 77-47-4 | Hexachlorocyclopentadiene | 0.05 ^c | 0.5 ^c |
| 67-72-1 | Hexachloroethane | 0.007 | 0.035 |
| 193-39-5 | Indeno(1,2,3-c,d)pyrene | 0.00043 ^a | 0.00215 |
| 78-59-1 | Isophorone | 1.4 | 1.4 |
| 72-43-5 | Methoxychlor | 0.04 ^c | 0.2 ^c |
| 74-83-9 | Methyl bromide (Bromomethane) | 0.0098 | 0.049 |
| 1634-04-4 | <u>Methyl tert-butyl ether</u> <u>Methyl tertiary-butyl ether</u> | 0.07 | 0.07 |
| 75-09-2 | Methylene chloride (Dichloromethane) | 0.005 ^c | 0.05 ^c |
| 91-20-3 | Naphthalene | 0.14 | 0.22 |
| 98-95-3 | Nitrobenzene ^b | 0.0035 | 0.0035 |

| Groundwater Remediation Objective | | | |
|-----------------------------------|---|---------------------|---------------------|
| CAS No. | Chemical Name | Class I (mg/L) | Class II (mg/L) |
| 86-30-6 | <i>N</i> -Nitrosodiphenylamine | 0.0032 ^a | 0.016 |
| 621-64-7 | <i>N</i> -Nitrosodi- <i>n</i> -propylamine | 0.0018 ^a | 0.0018 |
| 87-86-5 | Pentachlorophenol | 0.001 ^c | 0.005 ^c |
| 108-95-2 | Phenol | 0.1 ^c | 0.1 ^c |
| 1918-02-1 | Picloram | 0.5 ^c | 5.0 ^c |
| 1336-36-3 | Polychlorinated biphenyls (PCBs) | 0.0005 ^c | 0.0025 ^c |
| 129-00-0 | Pyrene | 0.21 | 1.05 |
| 122-34-9 | Simazine | 0.004 ^c | 0.04 ^c |
| 100-42-5 | Styrene | 0.1 ^c | 0.5 ^c |
| 93-72-1 | 2,4,5-TP (Silvex) | 0.05 ^c | 0.25 ^c |
| 127-18-4 | Tetrachloroethylene (Perchloroethylene) | 0.005 ^c | 0.025 ^c |
| 108-88-3 | Toluene | 1.0 ^c | 2.5 ^c |
| 8001-35-2 | Toxaphene | 0.003 ^c | 0.015 ^c |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.07 ^c | 0.7 ^c |
| 71-55-6 | 1,1,1-Trichloroethane ^b | 0.2 ^c | 1.0 ^c |
| 79-00-5 | 1,1,2-Trichloroethane | 0.005 ^c | 0.05 ^c |
| 79-01-6 | Trichloroethylene | 0.005 ^c | 0.025 ^c |
| 108-05-4 | Vinyl acetate | 7.0 | 7.0 |
| 75-01-4 | Vinyl chloride | 0.002 ^c | 0.01 ^c |
| 1330-20-7 | Xylenes (total) | 10.0 ^c | 10.0 ^c |
| Ionizable Organics | | | |
| 65-85-0 | Benzoic Acid | 28 | 28 |
| 106-47-8 | 4-Chloroaniline (<i>p</i> -Chloroaniline) | 0.028 | 0.028 |
| 95-57-8 | 2-Chlorophenol | 0.035 | 0.175 |
| 120-83-2 | 2,4-Dichlorophenol | 0.021 | 0.021 |
| 105-67-9 | 2,4-Dimethylphenol | 0.14 | 0.14 |
| 51-28-5 | 2,4-Dinitrophenol | 0.014 | 0.014 |
| 95-48-7 | 2-Methylphenol (<i>o</i> -Cresol) | 0.35 | 0.35 |

Groundwater Remediation Objective

| CAS No. | Chemical Name | Class I (mg/L) | Class II (mg/L) |
|-------------------|---------------------------|---------------------|--------------------|
| 95-95-4 | 2,4,5-Trichlorophenol | 0.7 | 3.5 |
| 88-06-2 | 2,4,6 Trichlorophenol | 0.01 ^a | 0.05 |
| Inorganics | | | |
| 7440-36-0 | Antimony | 0.006 ^c | 0.024 ^c |
| 7440-38-2 | Arsenic | 0.05 ^c | 0.2 ^c |
| 7440-39-3 | Barium | 2.0 ^c | 2.0 ^c |
| 7440-41-7 | Beryllium | 0.004 ^c | 0.5 ^c |
| 7440-42-8 | Boron | 2.0 ^c | 2.0 ^c |
| 7440-43-9 | Cadmium | 0.005 ^c | 0.05 ^c |
| 16887-00-6 | Chloride | 200 ^c | 200 ^c |
| 7440-47-3 | Chromium, total | 0.1 ^c | 1.0 ^c |
| 18540-29-9 | Chromium, ion, hexavalent | --- | --- |
| 7440-48-4 | Cobalt | 1.0 ^c | 1.0 ^c |
| 7440-50-8 | Copper | 0.65 ^c | 0.65 ^c |
| 57-12-5 | Cyanide | 0.2 ^c | 0.6 ^c |
| 7782-41-4 | Fluoride | 4.0 ^c | 4.0 ^c |
| 15438-31-0 | Iron | 5.0 ^c | 5.0 ^c |
| 7439-92-1 | Lead | 0.0075 ^c | 0.1 ^c |
| 7439-96-5 | Manganese | 0.15 ^c | 10.0 ^c |
| 7439-97-6 | Mercury | 0.002 ^c | 0.01 ^c |
| 7440-02-0 | Nickel | 0.1 ^c | 2.0 ^c |
| 14797-55-8 | Nitrate as N | 10.0 ^c | 100 ^c |
| 7782-49-2 | Selenium | 0.05 ^c | 0.05 ^c |
| 7440-22-4 | Silver | 0.05 ^c | --- |
| 14808-79-8 | Sulfate | 400 ^c | 400 ^c |

| Groundwater Remediation Objective | | | |
|-----------------------------------|-----------------------|--------------------|--------------------|
| CAS No. | Chemical Name | Class I (mg/L) | Class II (mg/L) |
| 7440-28-0 | Thallium | 0.002 ^c | 0.02 ^c |
| 7440-62-2 | Vanadium ^b | 0.049 | 0.1 |
| 7440-66-6 | Zinc | 5.0 ^c | 10 ^c |

Chemical Name and Groundwater Remediation Objective Notations

^a The groundwater remediation objective is equal to the ADL for carcinogens according to the procedures specified in 35 Ill. Adm. Code 620.

^b Oral Reference Dose and/or Reference Concentration under review by USEPA. Listed values subject to change.

^c Value listed is also the Groundwater Quality Standard for this chemical pursuant to 35 Ill. Adm. Code 620.410 for Class I Groundwater or 35 Ill. Adm. Code 620.420 for Class II Groundwater.

(Source: Amended at 26 Ill. Reg. ____, effective ____)

Section 742.APPENDIX B: Tier 1 Tables and Illustrations

Section 742.TABLE F: Values Used to Calculate the Tier 1 Soil Remediation Objectives for the Soil Component of the Groundwater Ingestion Route

| CAS No. | Chemical Name | GW _{obj} Concentration used to Calculate Tier 1 Soil Remediation Objectives ^a | |
|------------|--|---|--------------------|
| | | Class I (mg/L) | Class II (mg/L) |
| 83-32-9 | Acenaphthene | 2.0 ^b | 10 |
| 67-64-1 | Acetone | 4.0 ^b | 4.0 |
| 15972-60-8 | Alachlor | 0.002 ^c | 0.01 ^c |
| 116-06-3 | Aldicarb | 0.003 ^c | 0.015 ^c |
| 309-00-2 | Aldrin | 5.0E-6 ^b | 2.5E-5 |
| 120-12-7 | Anthracene | 10 ^b | 50 |
| 1912-24-9 | Atrazine | 0.003 ^c | 0.015 ^c |
| 71-43-2 | Benzene | 0.005 ^c | 0.025 ^c |
| 56-55-3 | Benzo(<i>a</i>)anthracene | 0.0001 ^b | 0.0005 |
| 205-99-2 | Benzo(<i>b</i>)fluoranthene | 0.0001 ^b | 0.0005 |
| 207-08-9 | Benzo(<i>k</i>)fluroanthene | 0.001 ^b | 0.005 |
| 50-32-8 | Benzo(<i>a</i>)pyrene | 0.0002 ^{a,c} | 0.002 ^c |
| 111-44-4 | Bis(2-chloroethyl)ether | 8.0E-5 ^b | 8.0E-5 |
| 117-81-7 | Bis(2-ethylhexyl)phthalate (Di(2-ethylhexyl)phthalate) | 0.006 ^{a,c} | 0.06 ^c |
| 75-27-4 | Bromodichloromethane (Dichlorobromomethane) | 0.1 ^b | 0.1 |
| 75-25-2 | Bromoform | 0.1 ^b | 0.01 |
| 71-36-3 | Butanol | 4.0 ^b | 4.0 |
| 85-68-7 | Butyl benzyl phthalate | 7.0 ^b | 35 |
| 86-74-8 | Carbazole | 0.004 ^b | 0.02 |
| 1563-66-2 | Carbofuran | 0.04 ^c | 0.2 ^c |
| 75-15-0 | Carbon disulfide | 4.0 ^b | 20 |
| 56-23-5 | Carbon tetrachloride | 0.005 ^c | 0.025 ^c |
| 57-74-9 | Chlordane | 0.002 ^c | 0.01 ^c |

| GW _{obj} Concentration used to Calculate Tier 1 Soil Remediation Objectives ^a | | | |
|--|--|------------------------|---------------------|
| CAS No. | Chemical Name | Class I (mg/L) | Class II (mg/L) |
| 108-90-7 | Chlorobenzene (Monochlorobenzene) | 0.1 ^c | 0.5 ^c |
| 124-48-1 | Chlorodibromomethane (Dibromochloromethane) | 0.06 ^b | 0.06 |
| 67-66-3 | Chloroform | 0.1 ^b | 0.5 |
| 218-01-9 | Chrysene | 0.1 ^b | 0.05 |
| 94-75-7 | 2,4-D | 0.07 ^c | 0.35 ^c |
| 75-99-0 | Dalapon | 0.2 ^c | 2.0 ^c |
| 72-54-8 | DDD | 0.0004 ^b | 0.002 |
| 72-55-9 | DDE | 0.0003 ^b | 0.0015 |
| 50-29-3 | DDT | 0.0003 ^b | 0.0015 |
| 53-70-3 | Dibenzo(<i>a,h</i>)anthracene | 1.0E-5 ^b | 5.0E-5 |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 0.0002 ^c | 0.0002 ^c |
| 106-93-4 | 1,2-Dibromoethane (Ethylene dibromide) | 0.00005 ^{a,c} | 0.0005 ^c |
| 84-74-2 | Di- <i>n</i> -butyl phthalate | 4.0 ^b | 20 |
| 95-50-1 | 1,2-Dichlorobenzene (<i>o</i> – Dichlorobenzene) | 0.6 ^c | 1.5 ^c |
| 106-46-7 | 1,4-Dichlorobenzene (<i>p</i> – Dichlorobenzene) | 0.075 ^c | 0.375 ^c |
| 91-94-1 | 3,3'-Dichlorobenzidine | 0.0002 ^b | 0.001 |
| 75-34-3 | 1,1-Dichloroethane | 4.0 ^b | 20 |
| 107-06-2 | 1,2-Dichloroethane (Ethylene dichloride) | 0.005 ^c | 0.025 ^c |
| 75-35-4 | 1,1-Dichloroethylene | 0.007 ^c | 0.035 ^c |
| 156-59-2 | <i>cis</i> -1,2-Dichloroethylene | 0.07 ^c | 0.2 ^c |
| 156-60-5 | <i>trans</i> -1,2-Dichloroethylene | 0.1 ^c | 0.5 ^c |
| 78-97-5 | 1,2-Dichloropropane | 0.005 ^c | 0.025 ^c |
| 542-75-6 | 1,3-Dichloropropene (1,3-Dichloropropylene, <i>cis</i> + <i>trans</i>) | 0.0005 ^b | 0.0025 |

| GW _{obj} Concentration used to Calculate Tier 1 Soil Remediation Objectives ^a | | | |
|---|--|---------------------|--------------------|
| CAS No. | Chemical Name | Class I (mg/L) | Class II (mg/L) |
| 60-57-1 | Dieldrin | 5.0E-6 ^b | 2.5E-5 |
| 84-66-2 | Diethyl phthalate | 30 ^b | 30 |
| 121-14-2 | 2,4-Dinitrotoluene | 0.0001 ^b | 0.0001 |
| 606-20-2 | 2,6-Dinitrotoluene | 0.0001 | 0.0001 |
| 88-85-7 | Dinoseb | 0.007 ^c | 0.07 ^c |
| 117-84-0 | Di-n-octyl phthalate | 0.7 ^b | 3.5 |
| 115-29-7 | Endosulfan | 0.2 ^b | 1.0 |
| 145-73-3 | Endothall | 0.1 ^c | 0.1 ^c |
| 72-20-8 | Endrin | 0.002 ^c | 0.01 ^c |
| 100-41-4 | Ethylbenzene | 0.7 ^c | 1.0 ^c |
| 206-44-0 | Fluoranthene | 1.0 ^b | 5.0 |
| 86-73-7 | Fluorene | 1.0 ^b | 5.0 |
| 76-44-8 | Heptachlor | 0.0004 ^c | 0.002 ^c |
| 1024-57-3 | Heptachlor epoxide | 0.0002 ^c | 0.001 ^c |
| 118-74-1 | Hexachlorobenzene | 0.001 ^b | 0.005 |
| 319-84-6 | <i>alpha</i> -HCH (<i>alpha</i> -BHC) | 1.0E-5 ^b | 5.0E-5 |
| 58-89-9 | <i>Gamma</i> -HCH (Lindane) | 0.0002 ^c | 0.001 ^c |
| 77-47-4 | Hexachlorocyclopentadiene | 0.05 ^c | 0.5 ^c |
| 67-72-1 | Hexachloroethane | 0.007 | 0.035 |
| 193-39-5 | Indeno(1,2,3- <i>c,d</i>)pyrene | 0.0001 ^b | 0.0005 |
| 78-59-1 | Isophorone | 1.4 | 1.4 |
| 72-43-5 | Methoxychlor | 0.04 ^c | 0.2 ^c |
| 74-83-9 | Methyl bromide (Bromomethane) | 0.05 ^b | 0.25 |
| 1634-04-4 | <u>Methyl tert butyl ether</u> <u>Methyl tertiary-butyl ether</u> | 0.07 | 0.07 |
| 75-09-2 | Methylene chloride (Dichloromethane) | 0.005 ^c | 0.05 ^c |
| 91-20-3 | Naphthalene | 0.14 | 0.22 |
| 98-95-3 | Nitrobenzene | 0.02 ^b | 0.02 |

| | | GW _{obj} Concentration used to Calculate Tier 1 Soil Remediation Objectives ^a | |
|---------------------------|---|--|--------------------|
| CAS No. | Chemical Name | Class I (mg/L) | Class II (mg/L) |
| 1918-02-1 | Picloram | 0.5 ^c | 5.0 ^c |
| 1336-36-3 | Polychlorinated biphenyls (PCBs) | --- | --- |
| 129-00-0 | Pyrene | 1.0 ^b | 5.0 |
| 122-34-9 | Simazine | 0.004 ^c | 0.04 ^c |
| 100-42-5 | Styrene | 0.1 ^c | 0.5 ^c |
| 93-72-1 | 2,4,5-TP (Silvex) | 0.05 ^c | 0.25 ^c |
| 127-18-4 | Tetrachloroethylene (Perchloroethylene) | 0.005 ^c | 0.025 ^c |
| 108-88-3 | Toluene | 1.0 ^c | 2.5 ^c |
| 8001-35-2 | Toxaphene | 0.003 ^c | 0.015 ^c |
| 120-82-1 | 1,2,4-Trichlorobenzene | 0.07 ^c | 0.7 ^c |
| 71-55-6 | 1,1,1-Trichloroethane ² | 0.2 ^c | 1.0 ^c |
| 79-00-5 | 1,1,2-Trichloroethane | 0.005 ^c | 0.05 ^c |
| 79-01-6 | Trichloroethylene | 0.005 ^c | 0.025 ^c |
| 108-05-4 | Vinyl acetate | 40 ^b | 40 |
| 75-01-4 | Vinyl chloride | 0.002 ^c | 0.01 ^c |
| 1330-20-7 | Xylenes (total) | 10.0 ^c | 10.0 ^c |
| Ionizable Organics | | | |
| 65-85-0 | Benzoic Acid | 100 ^b | 100 |
| 106-47-8 | 4-Chloroaniline (<i>p</i> -Chloroaniline) | 0.1 ^b | 0.1 |
| 95-57-8 | 2-Chlorophenol | 0.2 ^b | 1.0 |
| 120-83-2 | 2,4-Dichlorophenol | 0.1 ^b | 0.1 |
| 105-67-9 | 2,4-Dimethylphenol | 0.7 ^b | 0.7 |
| 51-28-5 | 2,4-Dinitrophenol | 0.04 ^b | 0.04 |
| 95-48-7 | 2-Methylphenol (<i>o</i> - Cresol) | 2.0 ^b | 2.0 |
| 86-30-6 | <i>N</i> -Nitrosodiphenylamine | 0.02 ^b | 0.1 |

| CAS No. | Chemical Name | GW _{obj} Concentration used to Calculate Tier 1 Soil Remediation Objectives ^a | |
|-------------------|--|---|--------------------|
| | | Class I (mg/L) | Class II (mg/L) |
| 621-64-7 | <i>N</i> -Nitrosodi- <i>n</i> -propylamine | 1.0E-5 ^b | 1.0E-5 |
| 87-86-5 | Pentachlorophenol | 0.001 ^{a,c} | 0.005 ^c |
| 108-95-2 | Phenol | 0.1 ^c | 0.1 ^c |
| 95-95-4 | 2,4,5-Trichlorophenol | 4.0 ^b | 20 |
| 88-06-2 | 2,4,6-Trichlorophenol | 0.008 ^b | 0.04 |
| Inorganics | | | |
| 7440-36-0 | Antimony | 0.006 ^c | 0.024 ^c |
| 7440-38-2 | Arsenic | 0.05 ^c | 0.2 ^c |
| 7440-39-3 | Barium | 2.0 ^c | 2.0 ^c |
| 7440-41-7 | Beryllium | 0.004 ^c | 0.5 ^c |
| 7440-42-8 | Boron | 2.0 ^c | 2.0 ^c |
| 7440-43-9 | Cadmium | 0.005 ^c | 0.05 ^c |
| 16887-00-6 | Chloride | 200 ^c | 200 ^c |
| 7440-47-3 | Chromium, total | 0.1 ^c | 1.0 ^c |
| 18540-29-9 | Chromium, ion, hexavalent | --- | --- |
| 7440-48-4 | Cobalt | 1.0 ^c | 1.0 ^c |
| 7440-50-8 | Copper | 0.65 ^c | 0.65 ^c |
| 57-12-5 | Cyanide | 0.2 ^c | 0.6 ^c |
| 7782-41-4 | Fluoride | 4.0 ^c | 4.0 ^c |
| 15438-31-0 | Iron | 5.0 ^c | 5.0 ^c |
| 7439-92-1 | Lead | 0.0075 ^c | 0.1 ^c |
| 7439-96-5 | Manganese | 0.15 ^c | 10.0 ^c |
| 7439-97-6 | Mercury | 0.002 ^c | 0.01 ^c |
| 7440-02-0 | Nickel | 0.1 ^c | 2.0 ^c |
| 14797-55-8 | Nitrate as N | 10.0 ^c | 100 ^c |
| 7782-49-2 | Selenium | 0.05 ^c | 0.05 ^c |
| 7440-22-4 | Silver | 0.05 ^c | --- |
| 14808-79-8 | Sulfate | 400 ^c | 400 ^c |

| | | GW _{obj} Concentration used to Calculate Tier 1 Soil Remediation Objectives ^a | |
|-----------|---------------|---|--------------------|
| CAS No. | Chemical Name | Class I (mg/L) | Class II (mg/L) |
| 7440-28-0 | Thallium | 0.002 ^c | 0.02 ^c |
| 7440-62-2 | Vanadium | 0.049 | 0.1 |
| 7440-66-6 | Zinc | 5.0 ^c | 10 ^c |

Chemical Name and Groundwater Remediation Objective Notations

^a The Equation S17 is used to calculate the Soil Remediation Objective for the Soil Component of the Groundwater Ingestion Route; this equation requires calculation of the Target Soil Leachate Concentration (C_w) from Equation S18: C_w = DF x GW_{obj}.

^b Value listed is the Water Health Based Limit (HBL) for this chemical from Soil Screening Guidance: User's Guide, incorporated by reference at Section 742.210. The HBL is equal to the non-zero MCLG (if available); the MCL (if available); or, for carcinogens, a cancer risk of 1.0E-6, and for noncarcinogens is equal to a Hazard Quotient of 1.0. NOTE: These GW_{obj} concentrations are not equal to the Tier 1 Groundwater Remediation Objectives for the Direct Ingestion of Groundwater Component of the Groundwater Ingestion Route, listed in Section 742.Appendix B, Table E.

^c Value listed is also the Groundwater Quality Standard for this chemical pursuant to 35 Ill. Adm. Code 620.410 for Class I Groundwater or 35 Ill. Adm. Code 620.420 for Class II Groundwater.

(Source: Amended at 26 Ill. Reg. ___, effective ____)

Section 742.APPENDIX C: Tier 2 Tables and Illustrations

Section 742.Table E: Default Physical and Chemical Parameters

| CAS No. | Chemical | Solubility in Water (S) (mg/L) | Diffusivity in Air (D _i) (cm ² /s) | Diffusivity in Water (D _w) (cm ² /s) | Dimensionless Henry's Law Constant (H') (25°C) | Organic Carbon Partition Coefficient (K _{oc}) (L/kg) | First Order Degradation Constant (λ) (d ⁻¹) |
|------------------|--------------|--------------------------------|---|---|--|--|---|
| Neutral Organics | | | | | | | |
| 83-32-9 | Acenaphthene | 4.24 | 0.0421 | 7.69E-6 | 0.00636 | 7,080 | 0.0034 |
| 67-64-1 | Acetone | 1,000,000 | 0.124 | 1.14E-5 | 0.00159 | 0.575 | 0.0495 |
| 15972-60-8 | Alachlor | 242 | 0.0198 | 5.69E-6 | 0.00000132 | 394 | No Data |
| 116-06-3 | Aldicarb | 6,000 | 0.0305 | 7.19E-6 | 0.000000574 | 12 | 0.00109 |
| 309-00-2 | Aldrin | 0.18 | 0.0132 | 4.86E-6 | 0.00697 | 2,450,000 | 0.00059 |
| 120-12-7 | Anthracene | 0.0434 | 0.0324 | 7.74E-6 | 0.00267 | 29,500 | 0.00075 |
| 1912-24-9 | Atrazine | 70 | 0.0258 | 6.69E-6 | 0.00000005 | 451 | No Data |
| 71-43-2 | Benzene | 1,750 | 0.088 | 9.80E-6 | 0.228 | 58.9 | 0.0009 |

| CAS No. | Chemical | Solubility in Water (S) (mg/L) | Diffusivity in Air (D _i) (cm ² /s) | Diffusivity in Water (D _w) (cm ² /s) | Dimensionless Henry's Law Constant (H') (25°C) | Organic Carbon Partition Coefficient (K _{oc}) (L/kg) | First Order Degradation Constant (λ) (d ⁻¹) |
|----------|----------------------------|--------------------------------|---|---|--|--|---|
| 56-55-3 | Benzo(a)anthracene | 0.0094 | 0.0510 | 9.00E-6 | 0.000137 | 398,000 | 0.00051 |
| 205-99-2 | Benzo(b)fluoranthene | 0.0015 | 0.0226 | 5.56E-6 | 0.00455 | 1,230,000 | 0.00057 |
| 207-08-9 | Benzo(k)fluoranthene | 0.0008 | 0.0226 | 5.56E-6 | 0.000034 | 1,230,000 | 0.00016 |
| 65-85-0 | Benzoic Acid | 3,500 | 0.0536 | 7.97E-6 | 0.0000631 | 0.600 | No Data |
| 50-32-8 | Benzo(a)pyrene | 0.00162 | 0.043 | 9.00E-6 | 0.0000463 | 1,020,000 | 0.00065 |
| 111-44-4 | Bis(2-chloroethyl)ether | 17,200 | 0.0692 | 7.53E-6 | 0.000738 | 15.5 | 0.0019 |
| 117-81-7 | Bis(2-ethylhexyl)phthalate | 0.34 | 0.0351 | 3.66E-6 | 0.00000418 | 15,100,000 | 0.0018 |
| 75-27-4 | Bromodichloromethane | 6,740 | 0.0298 | 1.06E-5 | 0.0656 | 55.0 | No Data |
| 75-25-2 | Bromoform | 3,100 | 0.0149 | 1.03E-5 | 0.0219 | 87.1 | 0.0019 |
| 71-36-3 | Butanol | 74,000 | 0.0800 | 9.30E-6 | 0.000361 | 6.92 | 0.01283 |
| 85-68-7 | Butyl Benzyl Phthalate | 2.69 | 0.0174 | 4.83E-6 | 0.0000517 | 57,500 | 0.00385 |
| 86-74-8 | Carbazole | 7.48 | 0.0390 | 7.03E-6 | 0.000000626 | 3,390 | No Data |

| CAS No. | Chemical | Solubility in Water (S) (mg/L) | Diffusivity in Air (D _i) (cm ² /s) | Diffusivity in Water (D _w) (cm ² /s) | Dimensionless Henry's Law Constant (H') (25°C) | Organic Carbon Partition Coefficient (K _{oc}) (L/kg) | First Order Degradation Constant (λ) (d ⁻¹) |
|-----------|----------------------|--------------------------------|---|---|--|--|---|
| 1563-66-2 | Carbofuran | 320 | 0.0249 | 6.63E-6 | .00377 | 37 | No Data |
| 75-15-0 | Carbon Disulfide | 1,190 | 0.104 | 1.00E-5 | 1.24 | 45.7 | No Data |
| 56-23-5 | Carbon Tetrachloride | 793 | 0.0780 | 8.80E-6 | 1.25 | 174 | 0.0019 |
| 57-74-9 | Chlordane | 0.056 | 0.0118 | 4.37E-6 | 0.00199 | 120,000 | 0.00025 |
| 106-47-8 | p-Chloroaniline | 5,300 | 0.0483 | 1.01E-5 | 0.0000136 | 66.1 | No Data |
| 108-09-7 | Chlorobenzene | 472 | 0.0730 | 8.70E-6 | 0.152 | 219 | 0.0023 |
| 124-48-1 | Chlorodibromomethane | 2,600 | 0.0196 | 1.05E-5 | 0.0321 | 63.1 | 0.00385 |
| 67-66-3 | Chloroform | 7,920 | 0.104 | 1.00E-5 | 0.15 | 39.8 | 0.00039 |
| 95-57-8 | 2-Chlorophenol | 22,000 | 0.0501 | 9.46E-6 | 0.016 | 388 | No Data |
| 218-01-9 | Chrysene | 0.0016 | 0.0248 | 6.21E-6 | 0.00388 | 398,000 | 0.00035 |
| 94-75-7 | 2,4-D | 680 | 0.0231 | 7.31E-6 | 0.00000041 | 451 | 0.00385 |
| 72-54-8 | 4,4'-DDD | 0.09 | 0.0169 | 4.76E-6 | 0.000164 | 1,000,000 | 0.000062 |

| CAS No. | Chemical | Solubility in Water (S) (mg/L) | Diffusivity in Air (D _i) (cm ² /s) | Diffusivity in Water (D _w) (cm ² /s) | Dimensionless Henry's Law Constant (H') (25°C) | Organic Carbon Partition Coefficient (K _{oc}) (L/kg) | First Order Degradation Constant (λ) (d ⁻¹) |
|----------|-----------------------------|--------------------------------|---|---|--|--|---|
| 72-55-9 | 4,4'-DDE | 0.12 | 0.0144 | 5.87E-6 | 0.000861 | 4,470,000 | 0.000062 |
| 50-29-3 | 4,4'-DDT | 0.025 | 0.0137 | 4.95E-6 | 0.000332 | 2,630,000 | 0.000062 |
| 75-99-0 | Dalapon | 900,000 | 0.0414 | 9.46E-6 | 0.00000264 | 5.8 | 0.005775 |
| 53-70-3 | Dibenzo(a,h)anthracene | 0.00249 | 0.0202 | 5.18E-6 | 0.000000603 | 3,800,000 | 0.00037 |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 1,200 | 0.0212 | 7.02E-6 | 0.00615 | 182 | 0.001925 |
| 106-93-4 | 1,2-Dibromoethane | 4,200 | 0.0287 | 8.06E-6 | 0.0303 | 93 | 0.005775 |
| 84-74-2 | Di-n-butyl Phthalate | 11.2 | 0.0438 | 7.86E-6 | 0.0000000385 | 33,900 | 0.03013 |
| 95-50-1 | 1,2-Dichlorobenzene | 156 | 0.0690 | 7.90E-6 | 0.0779 | 617 | 0.0019 |
| 106-46-7 | 1,4-Dichlorobenzene | 73.8 | 0.0690 | 7.90E-6 | 0.0996 | 617 | 0.0019 |
| 91-94-1 | 3,3-Dichlorobenzidine | 3.11 | 0.0194 | 6.74E-6 | 0.000000164 | 724 | 0.0019 |

| CAS No. | Chemical | Solubility in Water (S) (mg/L) | Diffusivity in Air (D _i) (cm ² /s) | Diffusivity in Water (D _w) (cm ² /s) | Dimensionless Henry's Law Constant (H') (25°C) | Organic Carbon Partition Coefficient (K _{oc}) (L/kg) | First Order Degradation Constant (λ) (d ⁻¹) |
|----------|-------------------------------------|--------------------------------|---|---|--|--|---|
| 75-34-3 | 1,1-Dichloroethane | 5,060 | 0.0742 | 1.05E-5 | 0.23 | 31.6 | 0.0019 |
| 107-06-2 | 1,2-Dichloroethane | 8,520 | 0.104 | 9.90E-6 | 0.0401 | 17.4 | 0.0019 |
| 75-35-4 | 1,1-Dichloroethylene | 2,250 | 0.0900 | 1.04E-5 | 1.07 | 58.9 | 0.0053 |
| 156-59-2 | Cis-1,2-Dichloroethylene | 3,500 | 0.0736 | 1.13E-5 | 0.167 | 35.5 | 0.00024 |
| 156-60-5 | Trans-1,2-Dichloroethylene | 6,300 | 0.0707 | 1.19E-5 | 0.385 | 52.5 | 0.00024 |
| 120-83-2 | 2,4-Dichlorophenol | 4,500 | 0.0346 | 8.77E-6 | 0.00013 | 147 | 0.00027 |
| 78-87-5 | 1,2-Dichloropropane | 2,800 | 0.0782 | 8.73E-6 | 0.115 | 43.7 | 0.00027 |
| 542-75-6 | 1,3-Dichloropropylene (cis + trans) | 2,800 | 0.0626 | 1.00E-5 | 0.726 | 45.7 | 0.061 |
| 60-57-1 | Dieldrin | 0.195 | 0.0125 | 4.74E-6 | 0.000619 | 21,400 | 0.00032 |
| 84-66-2 | Diethyl Phthalate | 1,080 | 0.0256 | 6.35E-6 | 0.0000185 | 288 | 0.00619 |
| 105-67-9 | 2,4-Dimethylphenol | 7,870 | 0.0584 | 8.69E-6 | 0.000082 | 209 | 0.0495 |
| 51-28-5 | 2,4-Dinitrophenol | 2,790 | 0.0273 | 9.06E-6 | 0.0000182 | 0.01 | 0.00132 |

| CAS No. | Chemical | Solubility in Water (S) (mg/L) | Diffusivity in Air (D _i) (cm ² /s) | Diffusivity in Water (D _w) (cm ² /s) | Dimensionless Henry's Law Constant (H') (25°C) | Organic Carbon Partition Coefficient (K _{oc}) (L/kg) | First Order Degradation Constant (λ) (d ⁻¹) |
|-----------|----------------------|--------------------------------|---|---|--|--|---|
| 121-14-2 | 2,4-Dinitrotoluene | 270 | 0.203 | 7.06E-6 | 0.0000038 | 95.5 | 0.00192 |
| 606-20-2 | 2,6-Dinitrotoluene | 182 | 0.0327 | 7.26E-6 | 0.0000306 | 69.2 | 0.00192 |
| 88-85-7 | Dinoseb | 52 | 0.0215 | 6.62E-6 | 0.0000189 | 1,120 | 0.002817 |
| 117-84-0 | Di-n-octyl Phthalate | 0.02 | 0.0151 | 3.58E-6 | 0.00274 | 83,200,000 | 0.0019 |
| 115-29-7 | Endosulfan | 0.51 | 0.0115 | 4.55E-6 | 0.000459 | 2,140 | 0.07629 |
| 145-73-3 | Endothall | 21,000 | 0.0291 | 8.07E-6 | 0.0000000107 | 0.29 | No Data |
| 72-20-8 | Endrin | 0.25 | 0.0125 | 4.74E-6 | 0.000308 | 12,300 | 0.00032 |
| 100-41-4 | Ethylbenzene | 169 | 0.0750 | 7.80E-6 | 0.323 | 363 | 0.003 |
| 206-44-0 | Fluoranthene | 0.206 | 0.0302 | 6.35E-6 | 0.00066 | 107,000 | 0.00019 |
| 86-73-7 | Fluorene | 1.98 | 0.0363 | 7.88E-6 | 0.00261 | 13,800 | 0.000691 |
| 76-44-8 | Heptachlor | 0.18 | 0.0112 | 5.69E-6 | 60.7 | 1,410,000 | 0.13 |
| 1024-57-3 | Heptachlor epoxide | 0.2 | 0.0132 | 4.23E-6 | 0.00039 | 83,200 | 0.00063 |

| CAS No. | Chemical | Solubility in Water (S) (mg/L) | Diffusivity in Air (D _i) (cm ² /s) | Diffusivity in Water (D _w) (cm ² /s) | Dimensionless Henry's Law Constant (H') (25°C) | Organic Carbon Partition Coefficient (K _{oc}) (L/kg) | First Order Degradation Constant (λ) (d ⁻¹) |
|-----------|--|--------------------------------|---|---|--|--|---|
| 118-74-1 | Hexachlorobenzene | 6.2 | 0.0542 | 5.91E-6 | 0.0541 | 55,000 | 0.00017 |
| 319-84-6 | Alpha-HCH (alpha-BHC) | 2.0 | 0.0142 | 7.34E-6 | 0.000435 | 1,230 | 0.0025 |
| 58-89-9 | Gamma-HCH (Lindane) | 6.8 | 0.0142 | 7.34E-6 | 0.000574 | 1,070 | 0.0029 |
| 77-47-4 | Hexachlorocyclo-Pentadiene | 1.8 | 0.0161 | 7.21E-6 | 1.11 | 200,000 | 0.012 |
| 67-72-1 | Hexachloroethane | 50 | 0.0025 | 6.80E-6 | 0.159 | 1,780 | 0.00192 |
| 193-39-5 | Indeno(1,2,3-c,d)pyrene | 0.000022 | 0.0190 | 5.66E-6 | 0.0000656 | 3,470,000 | 0.00047 |
| 78-59-1 | Isophorone | 12,000 | 0.0623 | 6.76E-6 | 0.000272 | 46.8 | 0.01238 |
| 7439-97-6 | Mercury | --- | 0.0307 | 6.30E-6 | 0.467 | --- | No Data |
| 72-43-5 | Methoxychlor | 0.045 | 0.0156 | 4.46E-6 | 0.000648 | 97,700 | 0.0019 |
| 74-83-9 | Methyl Bromide | 15,200 | 0.0728 | 1.21E-5 | 0.256 | 10.5 | 0.01824 |
| 1634-04-4 | <u>Methyl tert-butyl ether</u> <u>Methyl tertiary-butyl ether</u> | <u>51,000</u> | <u>0.102</u> | <u>1.10E-5</u> | <u>0.0241</u> | <u>11.5</u> | <u>No Data</u> |
| 75-09-2 | Methylene Chloride | 13,000 | 0.101 | 1.17E-5 | 0.0898 | 11.7 | 0.012 |
| 65-48-7 | 2-Methylphenol | 26,000 | 0.0740 | 8.30E-6 | 0.0000492 | 91.2 | 0.0495 |

| CAS No. | Chemical | Solubility in Water (S) (mg/L) | Diffusivity in Air (D _i) (cm ² /s) | Diffusivity in Water (D _w) (cm ² /s) | Dimensionless Henry's Law Constant (H') (25°C) | Organic Carbon Partition Coefficient (K _{oc}) (L/kg) | First Order Degradation Constant (λ) (d ⁻¹) |
|-----------|----------------------------------|--------------------------------|---|---|--|--|---|
| 91-20-3 | Naphthalene | 31.0 | 0.0590 | 7.50E-6 | 0.0198 | 2,000 | 0.0027 |
| 98-95-3 | Nitrobenzene | 2,090 | 0.0760 | 8.60E-6 | 0.000984 | 64.6 | 0.00176 |
| 86-30-6 | N-Nitrosodiphenylamine | 35.1 | 0.0312 | 6.35E-6 | 0.000205 | 1,290 | 0.01 |
| 621-64-7 | N-Nitrosodi-n-propylamine | 9,890 | 0.0545 | 8.17E-6 | 0.0000923 | 24.0 | 0.0019 |
| 87-86-5 | Pentachlorophenol | 1,950 | 0.0560 | 6.10E-6 | 0.000001 | 592 | 0.00045 |
| 108-95-2 | Phenol | 82,800 | 0.0820 | 9.10E-6 | 0.0000163 | 28.8 | 0.099 |
| 1918-02-1 | Picloram | 430 | 0.0255 | 5.28E-6 | 0.00000000166 | 1.98 | No Data |
| 1336-36-3 | Polychlorinated biphenyls (PCBs) | 0.7 | ----- ^a | ----- ^a | ----- ^a | 309,000 | No Data |
| 129-00-0 | Pyrene | 0.135 | 0.0272 | 7.24E-6 | 0.000451 | 105,000 | 0.00018 |
| 122-34-9 | Simazine | 5 | 0.027 | 7.36E-6 | 0.0000000133 | 133 | No Data |
| 100-42-5 | Styrene | 310 | 0.0710 | 8.00E-6 | 0.113 | 776 | 0.0033 |
| 93-72-1 | 2,4,5-TP (Silvex) | 31 | 0.0194 | 5.83E-6 | 0.0000000032 | 5,440 | No Data |

| CAS No. | Chemical | Solubility in Water (S) (mg/L) | Diffusivity in Air (D _i) (cm ² /s) | Diffusivity in Water (D _w) (cm ² /s) | Dimensionless Henry's Law Constant (H') (25°C) | Organic Carbon Partition Coefficient (K _{oc}) (L/kg) | First Order Degradation Constant (λ) (d ⁻¹) |
|-----------|------------------------|--------------------------------|---|---|--|--|---|
| 127-18-4 | Tetrachloroethylene | 200 | 0.0720 | 8.20E-6 | 0.754 | 155 | 0.00096 |
| 108-88-3 | Toluene | 526 | 0.0870 | 8.60E-6 | 0.272 | 182 | 0.011 |
| 8001-35-2 | Toxaphene | 0.74 | 0.0116 | 4.34E-6 | 0.000246 | 257,000 | No Data |
| 120-82-1 | 1,2,4-Trichlorobenzene | 300 | 0.0300 | 8.23E-6 | 0.0582 | 1,780 | 0.0019 |
| 71-55-6 | 1,1,1-Trichloroethane | 1,330 | 0.0780 | 8.80E-6 | 0.705 | 110 | 0.0013 |
| 79-00-5 | 1,1,2-Trichloroethane | 4,420 | 0.0780 | 8.80E-6 | 0.0374 | 50.1 | 0.00095 |
| 79-01-6 | Trichloroethylene | 1,100 | 0.0790 | 9.10E-6 | 0.422 | 166 | 0.00042 |
| 95-95-4 | 2,4,5-Trichlorophenol | 1,200 | 0.0291 | 7.03E-6 | 0.000178 | 1,600 | 0.00038 |
| 88-06-2 | 2,4,6-Trichlorophenol | 800 | 0.0318 | 6.25E-6 | 0.000319 | 381 | 0.00038 |
| 108-05-4 | Vinyl Acetate | 20,000 | 0.0850 | 9.20E-6 | 0.021 | 5.25 | No Data |
| 57-01-4 | Vinyl Chloride | 2,760 | 0.106 | 1.23E-6 | 1.11 | 18.6 | 0.00024 |
| 108-38-3 | m-Xylene | 161 | 0.070 | 7.80E-6 | 0.301 | 407 | 0.0019 |

| CAS No. | Chemical | Solubility in Water (S) (mg/L) | Diffusivity in Air (D _i) (cm ² /s) | Diffusivity in Water (D _w) (cm ² /s) | Dimensionless Henry's Law Constant (H') (25°C) | Organic Carbon Partition Coefficient (K _{oc}) (L/kg) | First Order Degradation Constant (λ) (d ⁻¹) |
|-----------|-----------------|--------------------------------|---|---|--|--|---|
| 95-47-6 | o-Xylene | 178 | 0.087 | 1.00E-5 | 0.213 | 363 | 0.0019 |
| 106-42-3 | p-Xylene | 185 | 0.0769 | 8.44E-6 | 0.314 | 389 | 0.0019 |
| 1330-20-7 | Xylenes (total) | 186 | 0.0720 | 9.34E-6 | 0.25 | 260 | 0.0019 |

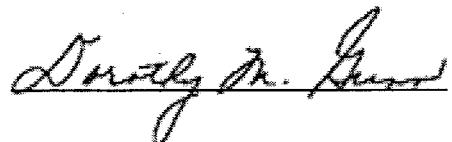
Chemical Abstracts Service (CAS) registry number. This number in the format xxx-xx-x, is unique for each chemical and allows efficient searching on computerized data bases.

^aSoil Remediation objectives are determined pursuant to 40 CFR 761, as incorporated by reference at Section 732.104 (the USEPA "PCB Spill Cleanup Policy"), for most sites; persons remediating sites should consult with BOL if calculation of Tier 2 soil remediation objectives is desired.

(Source: Amended at 26 Ill. Reg. ___, effective ____)

IT IS SO ORDERED.

I, Dorothy M. Gunn, Clerk of the Illinois Pollution Control Board, certify that the Board adopted the above opinion and order on January 24, 2002, by a vote of 7-0.

A handwritten signature in black ink, appearing to read "Dorothy M. Gunn". The signature is fluid and cursive, with a horizontal line drawn through it for emphasis.

Dorothy M. Gunn, Clerk
Illinois Pollution Control Board