

4. CHEMICAL AND PHYSICAL INFORMATION

4.1 CHEMICAL IDENTITY

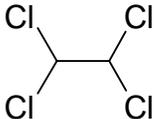
Information regarding the chemical identity of 1,1,2,2-tetrachloroethane is located in Table 4-1.

4.2 PHYSICAL AND CHEMICAL PROPERTIES

Information regarding the physical and chemical properties of 1,1,2,2-tetrachloroethane is located in Table 4-2.

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Table 4-1. Chemical Identity of 1,1,2,2-Tetrachloroethane

| Characteristics | Information | References |
|-------------------------|--|------------------------|
| Chemical name | 1,1,2,2-Tetrachloroethane | Lide 2005 |
| Synonyms | Tetrachloroethane; acetylene tetrachloride; dichloro-2,2-dichloroethane; s-tetrachloroethane | ChemID 2004 |
| Trade names | Bonoform, Cellon, Westron, Acetosol | ChemID 2004; HSDB 2006 |
| Chemical formula | C ₂ H ₂ Cl ₄ | O'Neil et al. 2001 |
| Chemical structure |  | Lide 2005 |
| Identification numbers: | | |
| CAS registry | 79-34-5 | ChemID 2004 |
| NIOSH RTECS | KI8575000 | RTECS 2006 |
| EPA hazardous waste | U209 | RTECS 2006 |
| OHM/TADS | 8100014 | HSDB 1995 |
| DOT/UN/NA/IMCO shipping | UN 1702; IMO 6.1 | HSDB 2006 |
| HSDB | 123 | HSDB 2006 |
| NCI | NCI-C03554 | RTECS 2006 |

CAS = Chemical Abstracts Service; DOT/UN/NA/IMCO -Department of Transportation/United Nations/North America/ Intergovernmental Maritime Consultive Organization; EPA = Environmental Protection Agency; HSDB = Hazardous Substance Data Bank; NCI = National Cancer Institute; NIOSH = National Institute for Occupational Safety and Health; OHM/TADS = Oil and Hazardous Materials/Technical Assistance Data Base; RTECS = Registry of Toxic Effects of Chemical Substances

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Table 4-2. Physical and Chemical Properties of 1,1,2,2-Tetrachloroethane

| Property | Tetrachloroethane | References |
|---|---|---|
| Molecular weight | 167.85 | O'Neil et al. 2001 |
| Color | Colorless | Lewis 2001 |
| Physical state | Liquid | Lewis 2001 |
| Melting point | -42.4 °C | O'Neil et al. 2001 |
| Boiling point | 116.7 °C | Lide 2005 |
| Density (20 °C) | 1.6 | Lewis 2001 |
| Odor | Sweetish, suffocating, chloroform-like | O'Neil et al. 2001 |
| Odor threshold | | |
| Water | 0.5 ppm | Amoore and Hautala 1983 |
| Air | 1.5 ppm | Amoore and Hautala 1983 |
| Solubility | | |
| Water at 25 °C | 2.83x10 ³ mg/L | Horvath et al. 1999 |
| Organic solvents | Miscible with methanol, ethanol, benzene, ether, petroleum ether, carbon tetrachloride, chloroform, carbon disulfide, dimethylformamide, oils | O'Neil et al. 2001 |
| Partition coefficients | | |
| Log octanol/water | 2.39 | Hansch et al. 1995 |
| K _{oc} | 46, 83, 118, 173, 216, 240 | Borisover and Graber 1997; Chiou et al. 1979; Chu and Chan 2000; Valsaraj et al. 1999 |
| Vapor pressure | | |
| 25 °C | 4.62 mmHg | AICHE 1995 |
| Henry's law constant | | |
| atm/m ³ -molecule at 25 °C | 3.67x10 ⁻⁴ atm-m ³ /mol | Leighton and Calo 1981 |
| Conversion factors | | |
| ppm (v/v) to mg/m ³ in air (20 °C) | 1 ppm=6.98 mg/m ³ | Verschueren 2001 |
| mg/m ³ to ppm (v/v) in air (20 °C) | 1 mg/m ³ =0.14 ppm | Verschueren 2001 |
| Bioconcentration factor | | |
| BCF | 8, 2 | ASTER 1995; Barrows et al. 1980 |