PROPYLENE GLYCOL 67

3. CHEMICAL AND PHYSICAL INFORMATION

3.1 CHEMICAL IDENTITY

Information regarding the chemical identity of propylene glycol is located in Table 3-1.

3.2 PHYSICAL AND CHEMICAL PROPERTIES

Information regarding the physical and chemical properties of propylene glycol is located in Table 3-2.

Table 3-1. Chemical Identity of Propylene Glycol^a

Characteristic	Information
Chemical name	Propylene glycol
Synonyms and trade names	1,2-Dihydroxypropane; 1,2-propanediol; 1,2-propylene glycol; 2,3-propanediol; hydroxy-propanol; alpha-propylene glycol; methyl glycol; methylethyl glycol; monopropylene glycol; trimethyl glycol.
Registered trade name(s)	PG-12; Sirlene
Chemical formula	$C_3H_8O_2$
Chemical structure ^b	CH_3 CH_2 OH CH_2 CH_3 CH_2 CH_3 CH
Identification numbers:	
CAS registry	57-55-6
NIOSH RTECS	TY2000000
EPA hazardous waste	No data
OHM/TADS	7216877
DOT/UN/NA/IMDG shipping	No data
HSDB	174
NCI	No data

^aAll information obtained from HSDB 1994b, except where noted.

CAS = Chemical Abstracts Service; DOT/UN/NA/IMDG = Department of Transportation/United Nations/North America/International Maritime Dangerous Goods Code; EPA = Environmental Protection Agency; HSDB = Hazardous Substances Data Bank; NCI = National Cancer Institute; NIOSH = National Institute for Occupational Safety and Health; OHM/TADS = Oil and Hazardous Materials/Technical Assistance Data System; RTECS = Registry of Toxic Effects of Chemical Substances

^bEPA 1987a

Table 3-2. Physical and Chemical Properties of Propylene Glycol^a

Property	Propylene glycol
Molecular weight	76.11 ^b
Color	Colorless ^d
Physical state	Liquid ^b
Melting point	-60 °C ^{e,b} (forms glass)
Boiling point	187.6; 188.2 °C ^b
Density:	
at 20 °C (g/cm³)	1.0361°
Odor	Odorless
Odor threshold	No data
Solubility:	
Water at 20 °C	Miscible with water
Organic solvent(s)	Soluble in alcohol, ether, benzene, soluble in acetone, chloroform ^b .
Partition coefficients:	
Log K _{ow}	-0.92 ^{f,g}
Log K _{oc}	0.88 [†] , 0.76 ^g
Vapor pressure at 20 °C	0.07 mm Hg ^{also d}
Henry's law constant at 25 °C	1.2x10 ⁻⁸ atm-m ³ /mole; 1.7x10 ⁻⁸ atm-m ³ /mole ⁹
Autoignition temperature	421.26 °C ^h ; 371 °C ⁱ
Flashpoint	99.04 °C ^{h,i}
Flammability limits	2.6–12.5% ^{h,i}
Conversion factors	1 ppm = 3.11 mg/m ^{3j} 1 mg/L = 321.6 ppm ^j
Explosive limits	No data

^aAll information obtained from HSDB 1995b, except where noted. ^bMerck 1989 ^cWeast 1988 ^dLewis 1993

^eDaubert and Danner 1980

^fEPA 1987a

⁹ASTER 1995

^hDaubert and Danner 1989

NFPA 1994

^jRowe and Wolf 1982