

MICHIGAN DEPARTMENT OF ENVIRONMENTAL QUALITY

INTEROFFICE COMMUNICATION

October 26, 2001

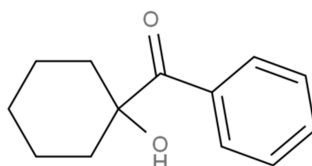
TO: File for 1-Hydroxycyclohexyl Phenyl Ketone (CAS# 947-19-3)
FROM: Michael Depa, Toxics Unit, Air Quality Division
SUBJECT: Screening Level Determination

The initial threshold screening level (ITSL) for 1-hydroxycyclohexyl phenyl ketone is 0.1 $\mu\text{g}/\text{m}^3$ based on an annual averaging time.

Other names found on the CAS registry printout for this CAS number include: α -hydroxy- α -cyclohexyl phenyl ketone and 1-benzoyl-1-hydroxycyclohexane.

The following references or databases were searched to identify data to determine the screening level: U.S. EPA Integrated Risk Information System (IRIS), Registry for Toxic Effects of Chemical Substances (RTECS), American Conference of Governmental and Industrial Hygienists (ACGIH) Threshold Limit Values (TLVs), National Institute for Occupational Safety and Health (NIOSH) Pocket Guide to Hazardous Chemicals, Environmental Protection Bureau Library, International Agency for Research on Cancer (IARC) Monographs, Chemical Abstract Service (CAS) Online (1967- October, 2001), National Library of Medicine (NLM), Health Effects Assessment Summary Tables (HEAST), and National Toxicology Program (NTP) Status Report. The EPA has not established a reference concentration (RfC) or reference dose (RfD) for 1-hydroxycyclohexyl phenyl ketone. The molecular weight of 1-hydroxycyclohexyl phenyl ketone is 206.03g. The molecular formula is $\text{C}_{13}\text{H}_{16}\text{O}_2$. There were no published records of a molecular structure for this compound; however, given the information from the CAS registry file the assumed molecular structure of 1-hydroxycyclohexyl phenyl ketone is shown in figure 1.

Figure 1



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