Chapter 16 Ethers, Epoxides, and Sulfides

16.1 Nomenclature of Ethers, Epoxides, and Sulfides

Substitutive IUPAC Names of Ethers

name as alkoxy derivatives of alkanes

CH₃OCH₂ CH₃ methoxyethane

CH₃CH₂OCH₂CH₂CH₂CH₂CI 1-chloro-3-ethoxypropane

CH₃CH₂OCH₂CH₃ ethoxyethane

Functional Class IUPAC Names of Ethers

name the groups attached to oxygen in alphabetical order as separate words; "ether" is last word

 $CH_3OCH_2CH_3$

ethyl methyl ether

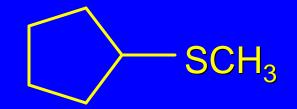
CH₃CH₂OCH₂CH₂CH₂CH₂CI 3-chloropropyl ethyl ether

CH₃CH₂OCH₂CH₃ diethyl ether Substitutive IUPAC Names of Ethers

name as alkylthio derivatives of alkanes

 $CH_3SCH_2CH_3$

methylthioethane

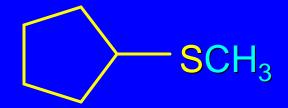


(methylthio)cyclopentane

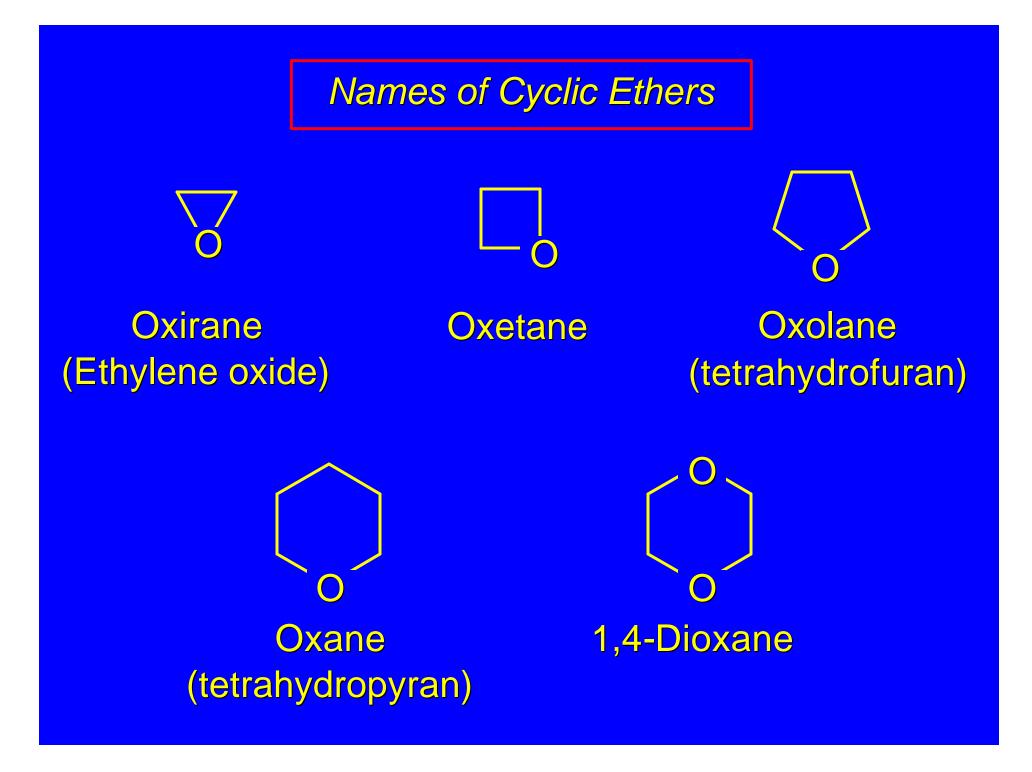
CH₃CH₂SCH₂CH₃ ethylthioethane Functional Class IUPAC Names of Sulfides

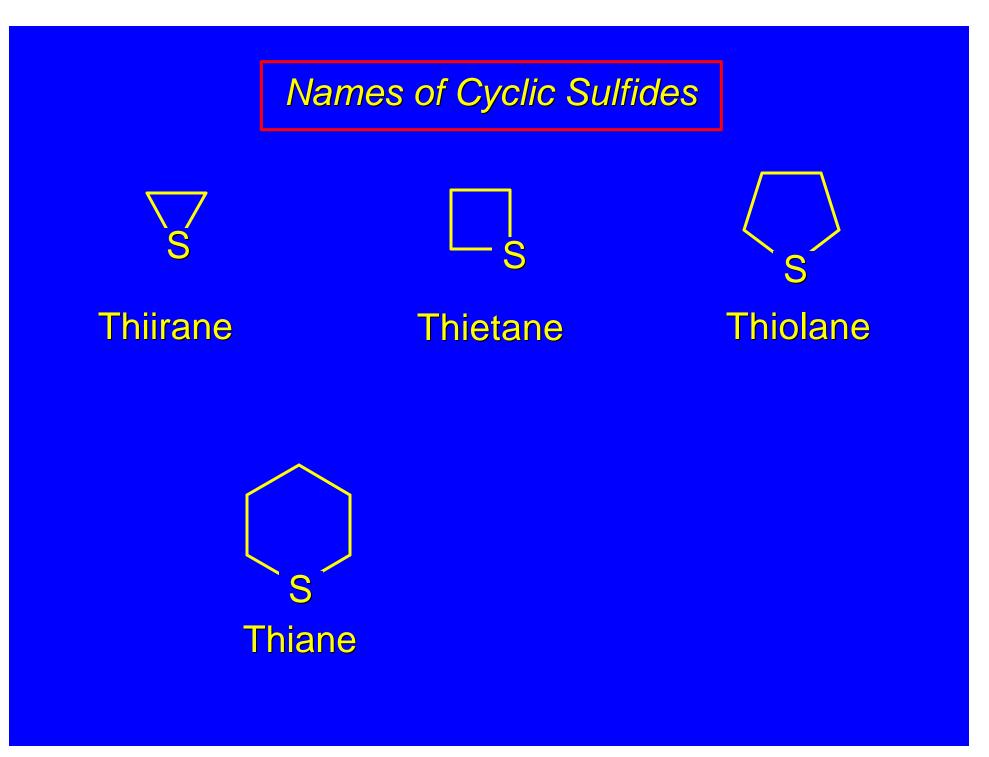
analogous to ethers, but replace "ether" as last word in the name by "sulfide."

CH₃SCH₂ CH₃ ethyl methyl sulfide



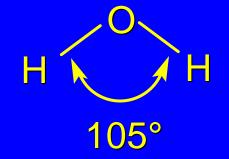
CH₃CH₂SCH₂CH₃ diethyl sulfide cyclopentyl methyl sulfide

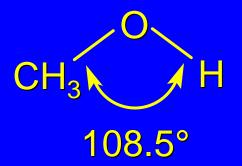


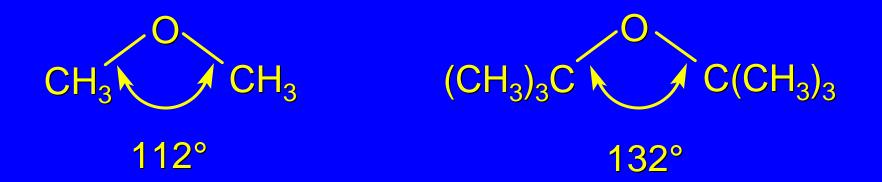


16.2 Structure and Bonding in Ethers and Epoxides

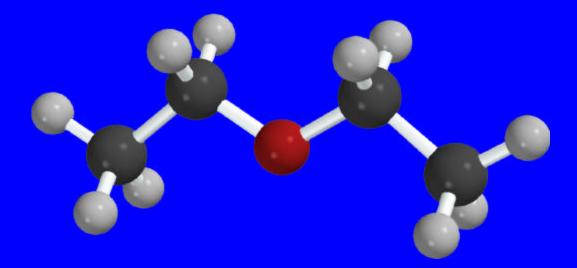
bent geometry at oxygen analogous to water and alcohols Bond angles at oxygen are sensitive to steric effects





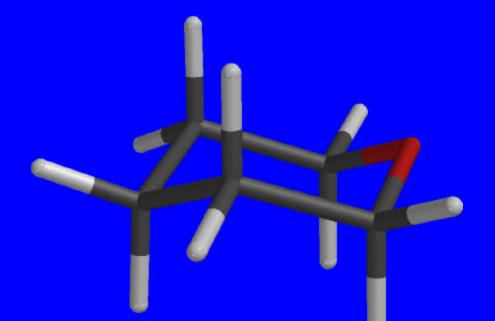


An oxygen atom affects geometry in much the same way as a CH₂ group



most stable conformation of diethyl ether resembles pentane

An oxygen atom affects geometry in much the same way as a CH₂ group

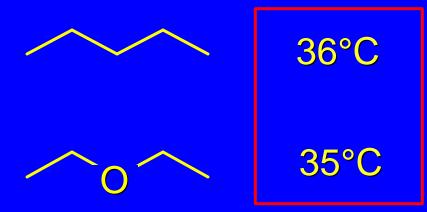


most stable conformation of tetrahydropyran resembles cyclohexane

16.3 Physical Properties of Ethers

Ethers resemble alkanes more than alcohols with respect to boiling point

boiling point



Intermolecular hydrogen bonding possible in alcohols; not possible in alkanes or ethers.



Ethers resemble alcohols more than alkanes with respect to solubility in water

solubility in water (g/100 mL)



very small



Hydrogen bonding to water possible for ethers and alcohols; not possible for alkanes.