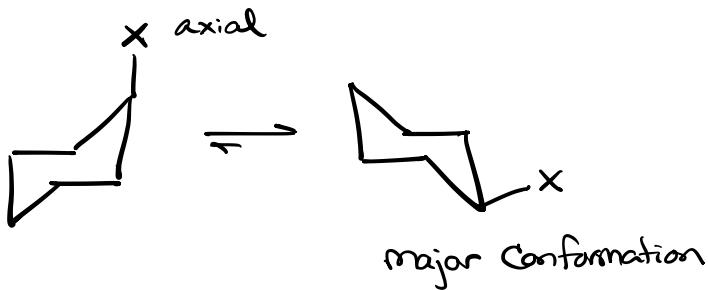
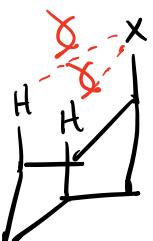


Six Member Ring Conformations

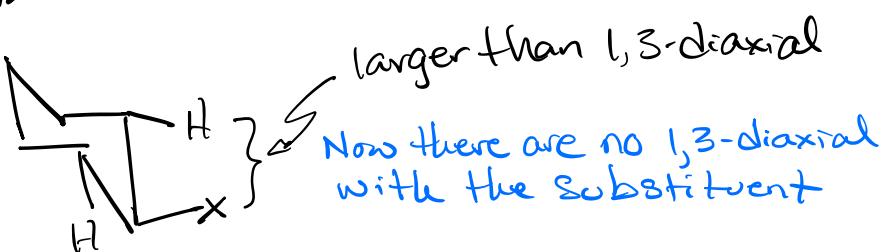


Two reasons

①

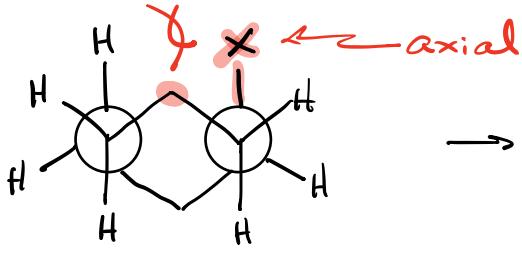


1,3-diaxial Steric interactions
Raise the energy of System

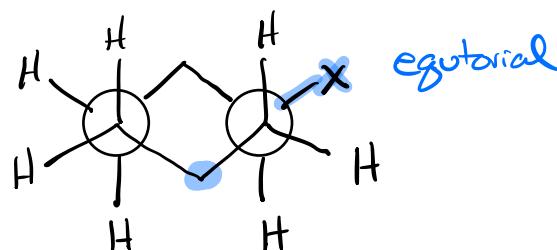


②

Gauche

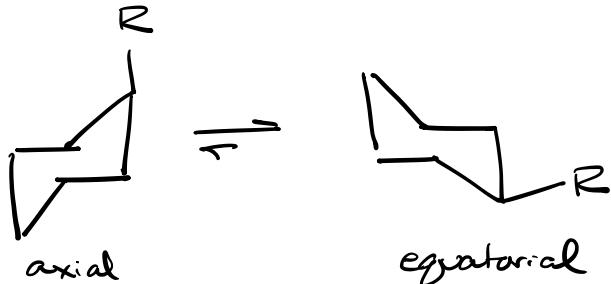


axial



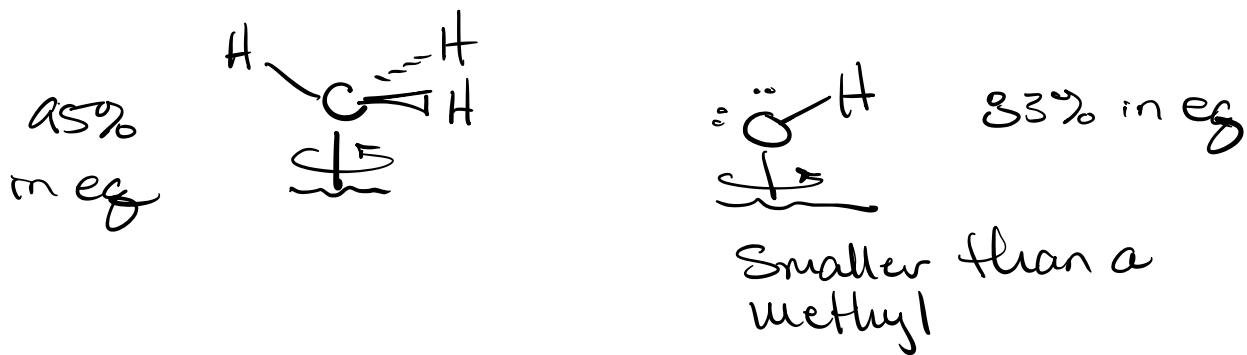
equatorial

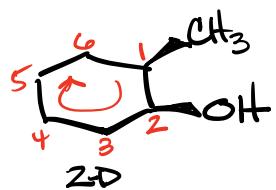
anti to ring
Carbon and is
lower energy



<u>R</u>	<u>% equatorial @ RT</u>
{-CH ₃	95%
-F	60%
t-butyl	99.99%
isopropyl	97%
-Cl	70%
-OH	83%
ethyl	96%

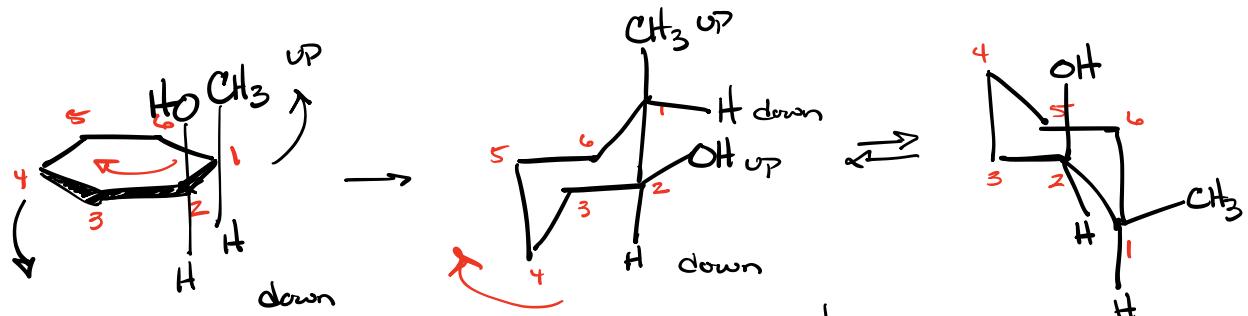
The larger the group the higher the % in equatorial position.



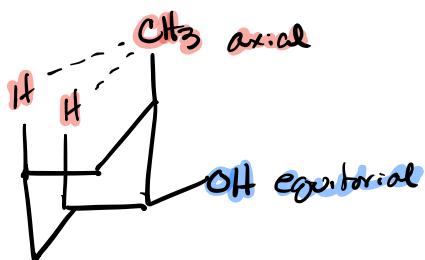


What is the lowest energy conformation?

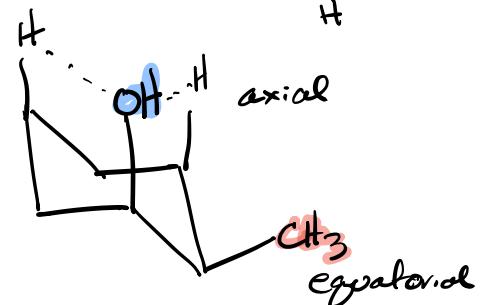
2-D \rightarrow 3-D \rightarrow Draw both conformations & analyse



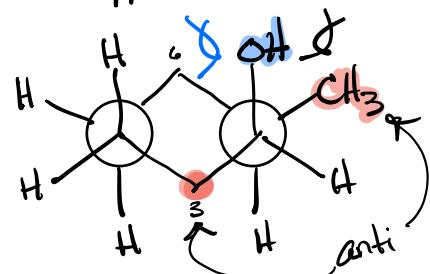
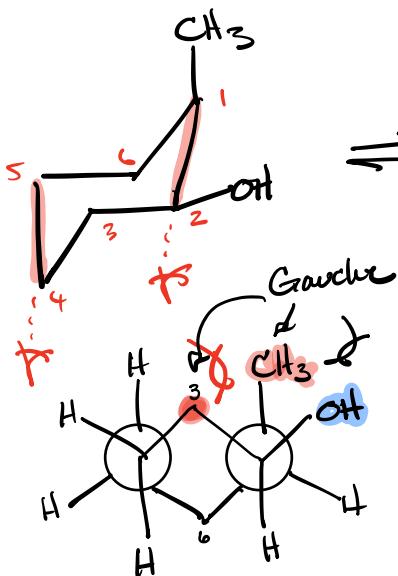
①



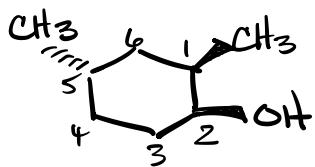
- strong 1,3-diaxial w/ large group (methyl)
- Small group equatorial



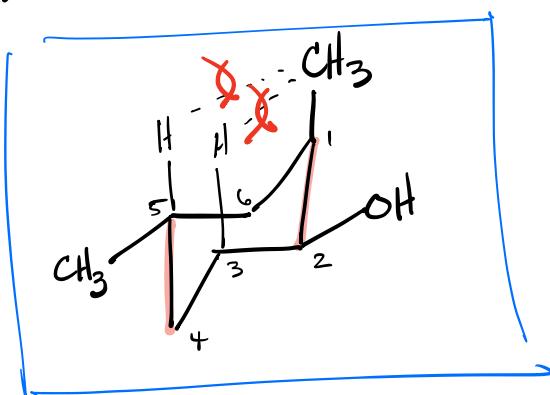
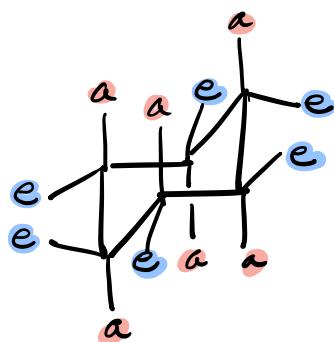
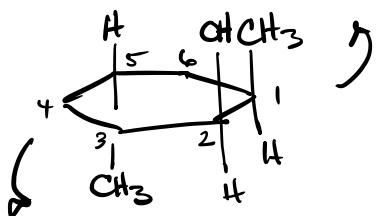
- lower 1,3-diaxials
- larger group in eq



3 Substituents



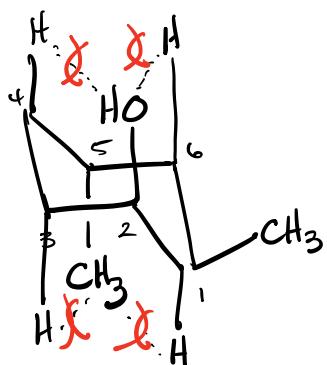
lowest energy conformation?



axial = 1 methyl

equatorial = 2 methyl, hydroxyl

Lower energy



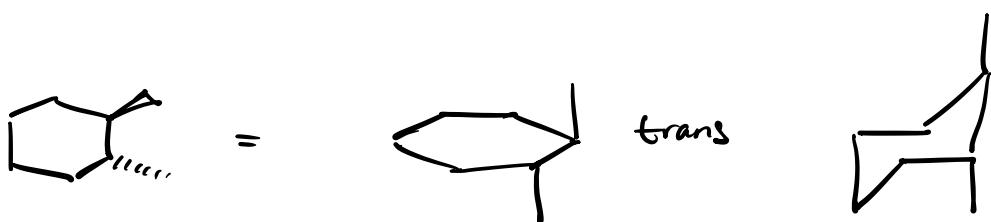
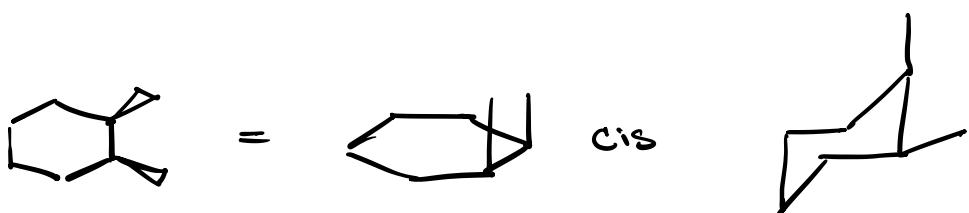
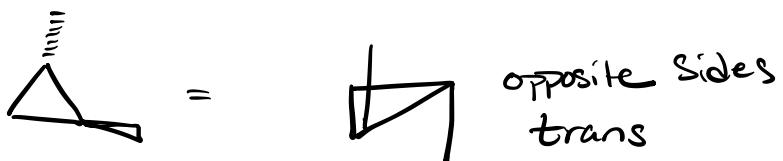
axial = 2 methyl, hydroxyl

equatorial = 1 methyl

Higher 1,3-diaxial

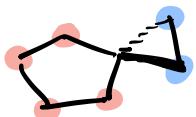
Higher energy

Stereochemistry



Multiple Rings - Joined Rings

Spiro Compounds



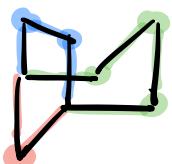
5 carbons total
pentane

Spiro [2,2] pentane

Spiro [2,4] Septane

#'s go in increasing order

Bicyclo

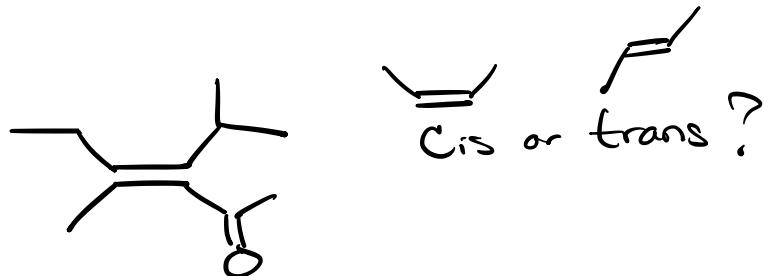


Bicyclo [3,2,1] Octane

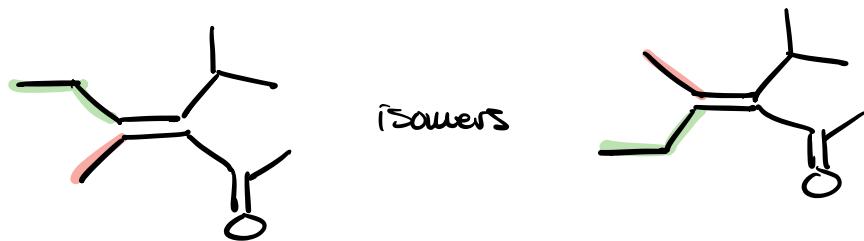
Bicyclo [4,3,0] Nonane

#'s go in decreasing order

Stereochemistry of alkenes



not well described by cis & trans



(E) vs (Z)

(E) Entgegen = "opposite" trans

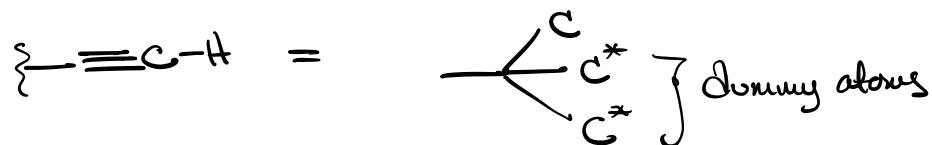
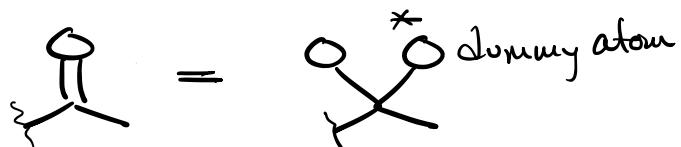
(Z) Zusammen = "Together" cis

Cahn-Ingold-Prelog Rules

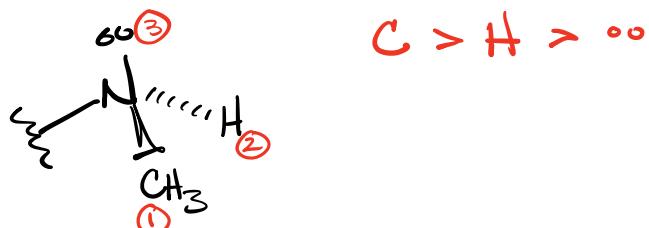
Assignment of priorities of Substituents

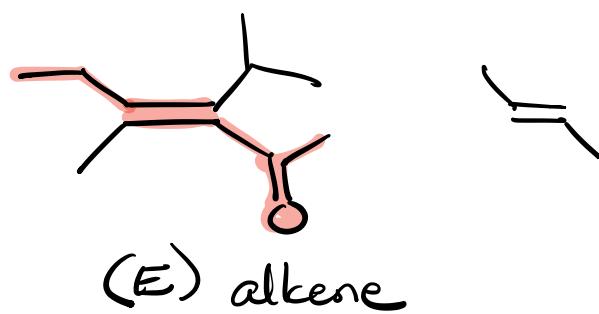
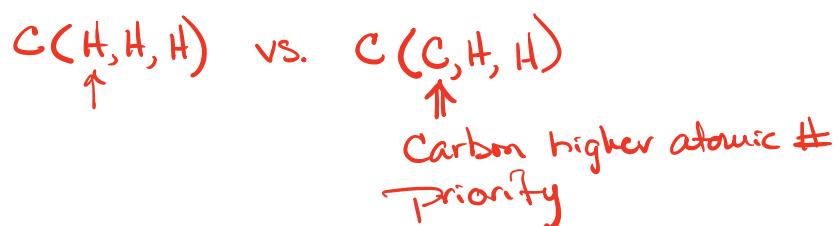
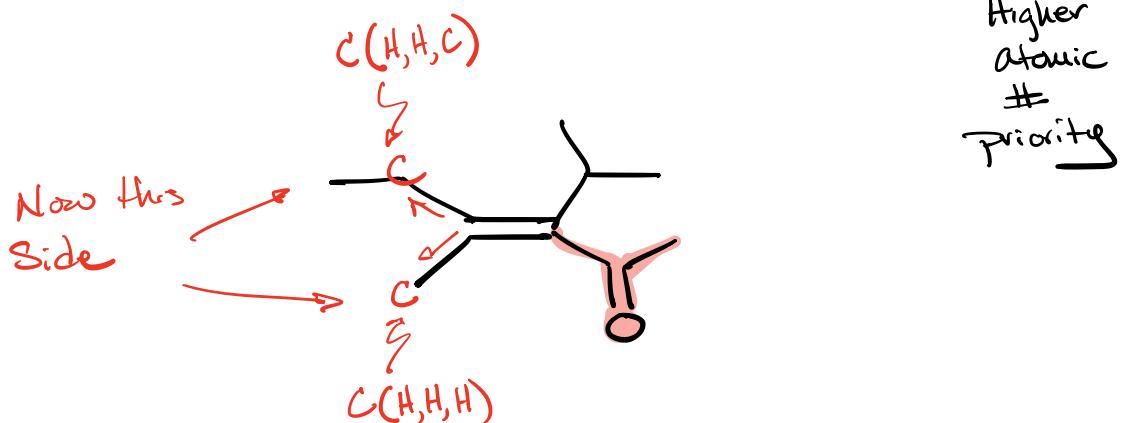
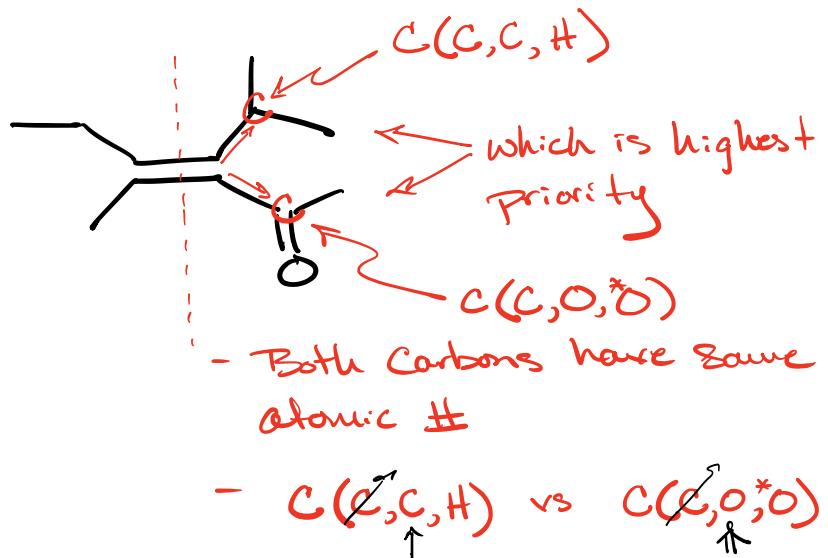
① The higher the atomic # at first point of difference, the higher the priority

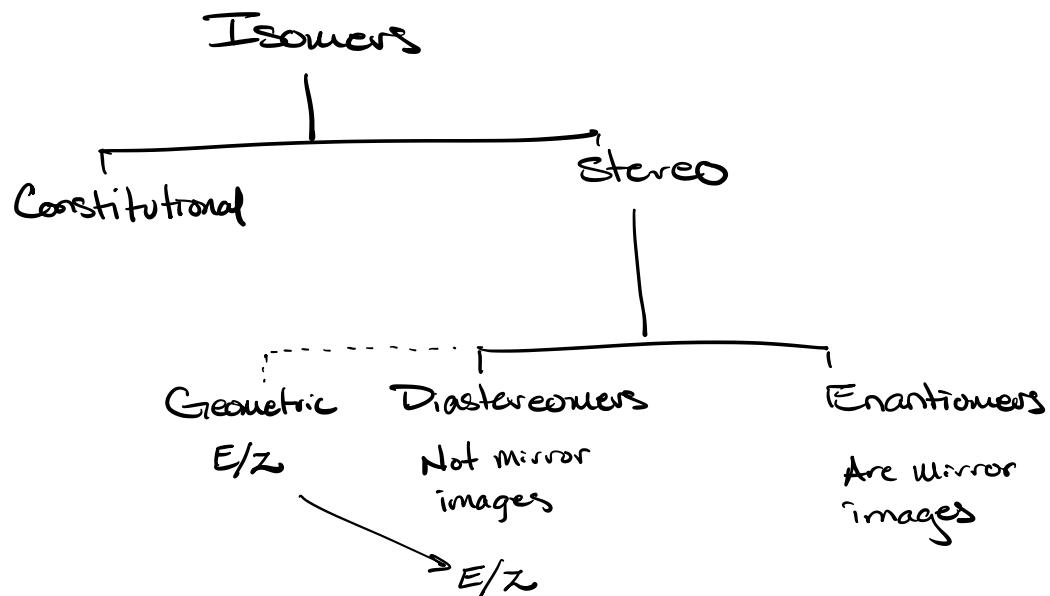
② Use dummy atoms for double bonds & triple bonds



③ Lone pairs have lowest priority as a group







If Rotation makes two molecules the same
 \Rightarrow Different Conformations of same molecule

If you need to break a bond and reconnect in new pattern to make two molecules the same
 \Rightarrow Isomers