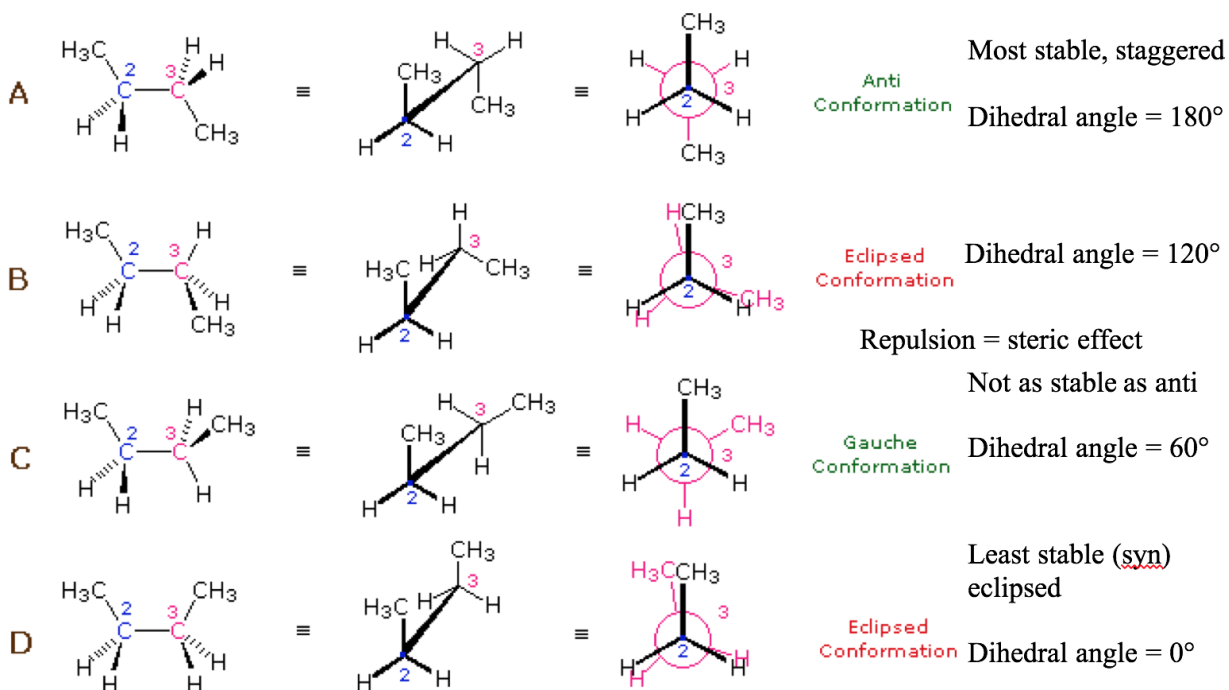


**Recall:** A molecular conformation is any spatial arrangement of atoms that can be interconverted by rotation about single bonds of the molecule.

Rotation around all bonds still very rapid.

Most stable (most populated conformation) is called anti and has groups as far away as possible.

**Note:** Newman Projection allows you to look at only one bond at a time. In the case of n-butane, we are looking through the central bond.

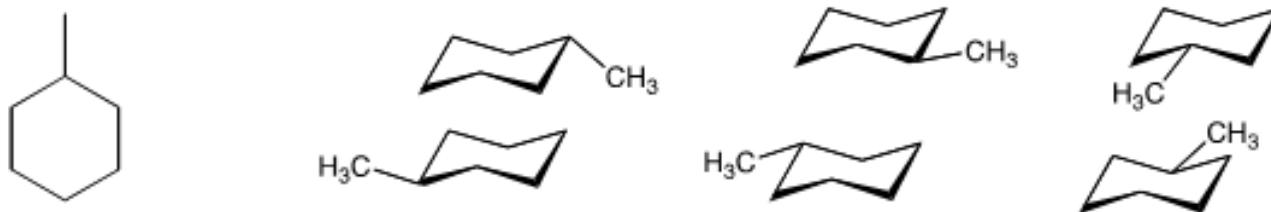


Cyclohexane molecules usually prefer to exist as chair conformations. Steric strain drives the lowest energy conformation.

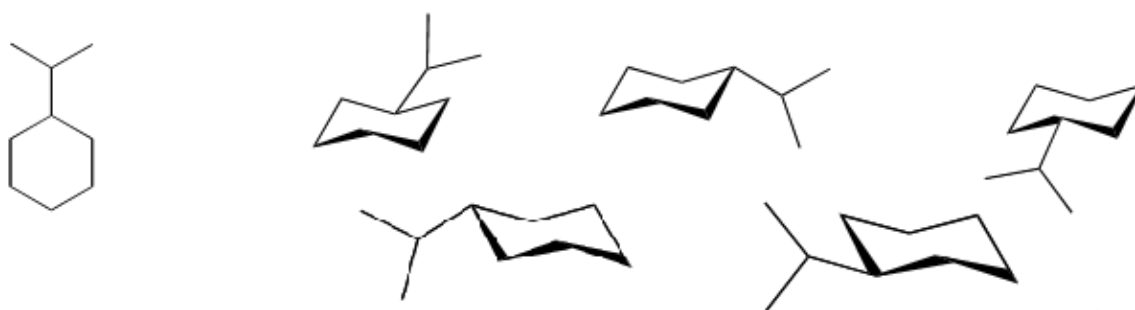
### Examples of Most Stable Conformation of Substituted Cyclohexanes :

#### 1. Methylcyclohexane

– 6 drawings of same molecule



## 2. Isopropylcyclohexane

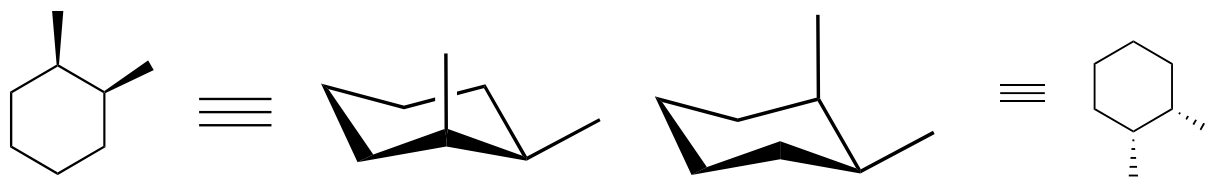


For most stable conformation, **largest group at equatorial position**

Examples of Most Stable Conformation of Polysubstituted cyclohexanes:

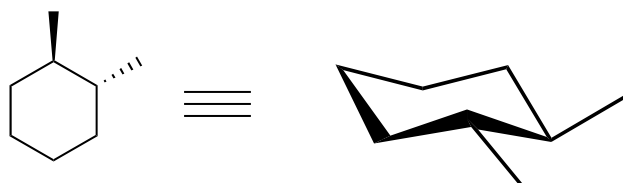
Examples:

1A) *cis*-1,2-dimethylcyclohexane



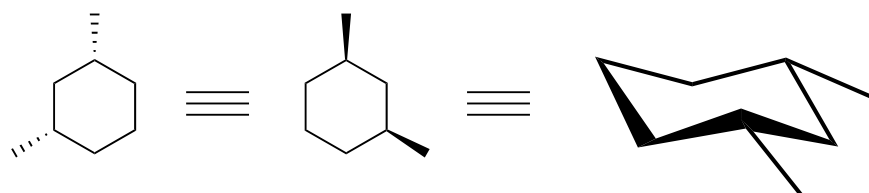
- One methyl group axial and one methyl group equatorial

1B) *trans*-1,2-dimethylcyclohexane (a stereoisomer (diastereomer) of above structures)



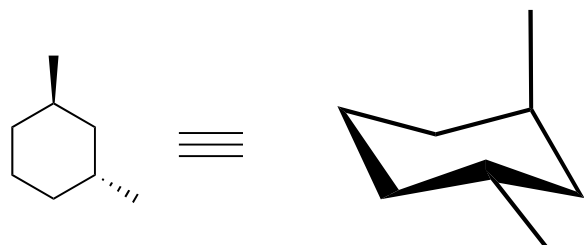
- Both methyl groups equatorial

2A) *cis*-1,3-dimethylcyclohexane (a structural isomer of above structures)



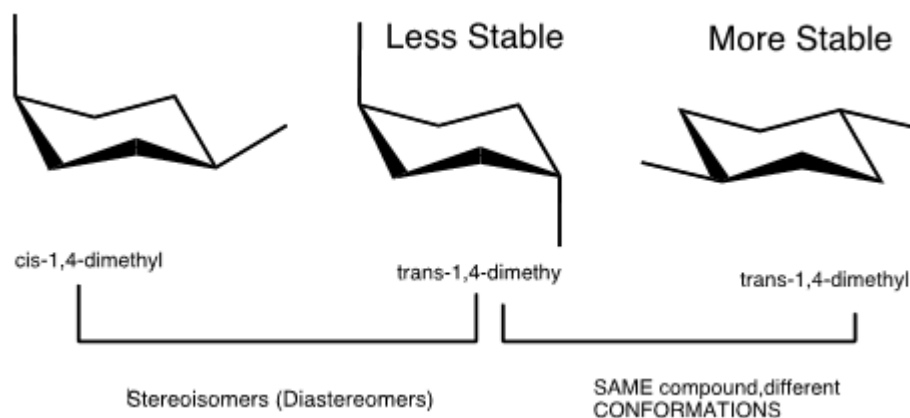
- Both methyl groups equatorial

2B) *trans*-1,3-dimethylcyclohexane: a stereoisomer of above *cis*-1,3-dimethylcyclohexane

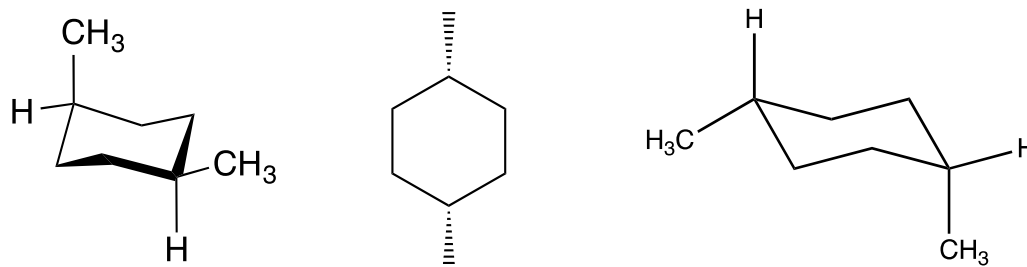


- One methyl group axial and one methyl group equatorial

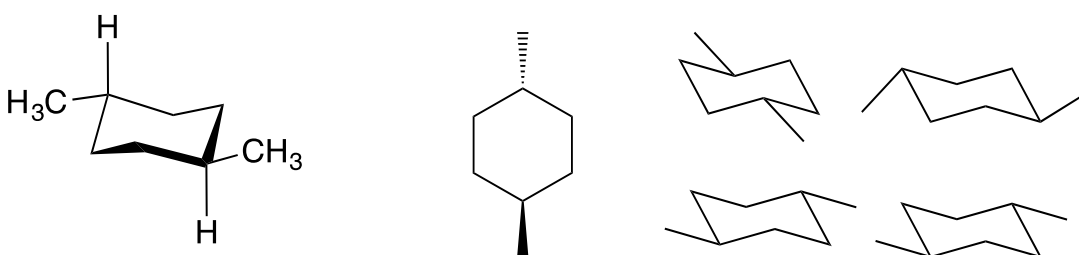
3A) *cis*-1,4-dimethylcyclohexane and *trans*-1,4-dimethylcyclohexane:



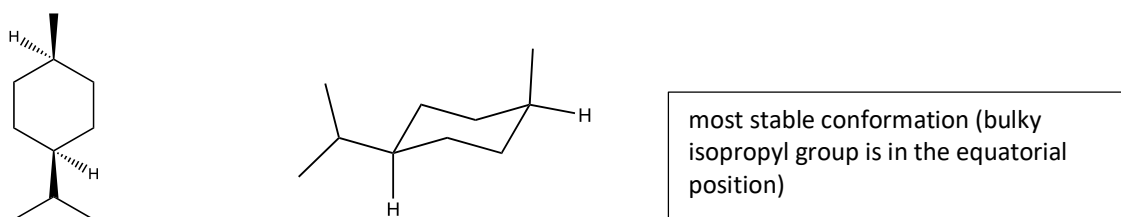
*cis*-1,4-dimethylcyclohexane:



3B) *trans*-1,4-dimethylcyclohexane:

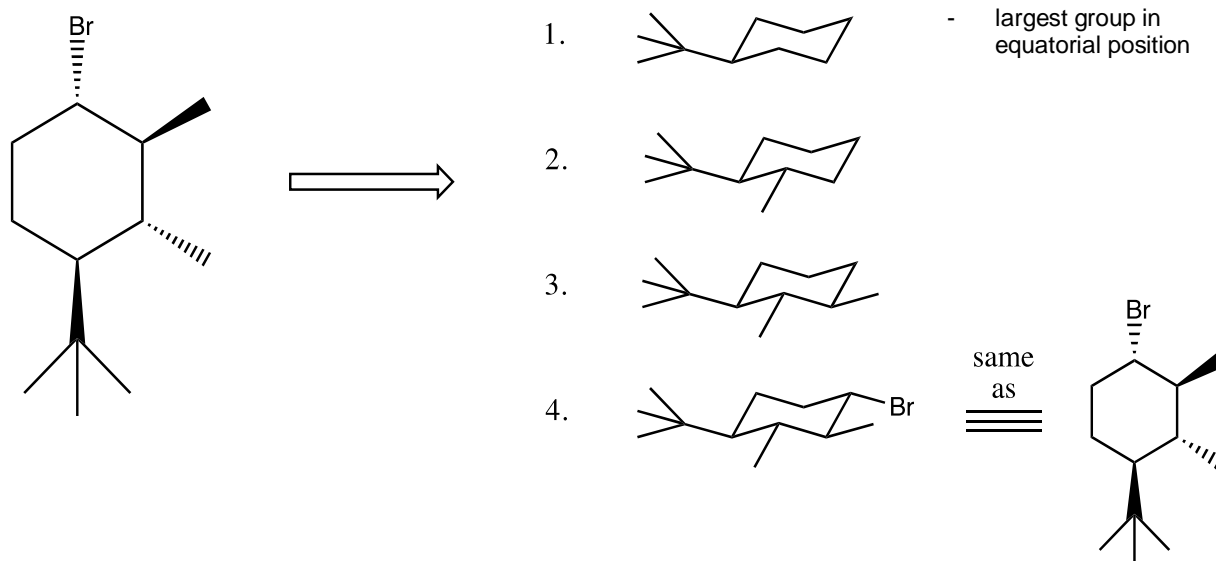


Application example: *cis*-1-isopropyl-4-methylcyclohexane

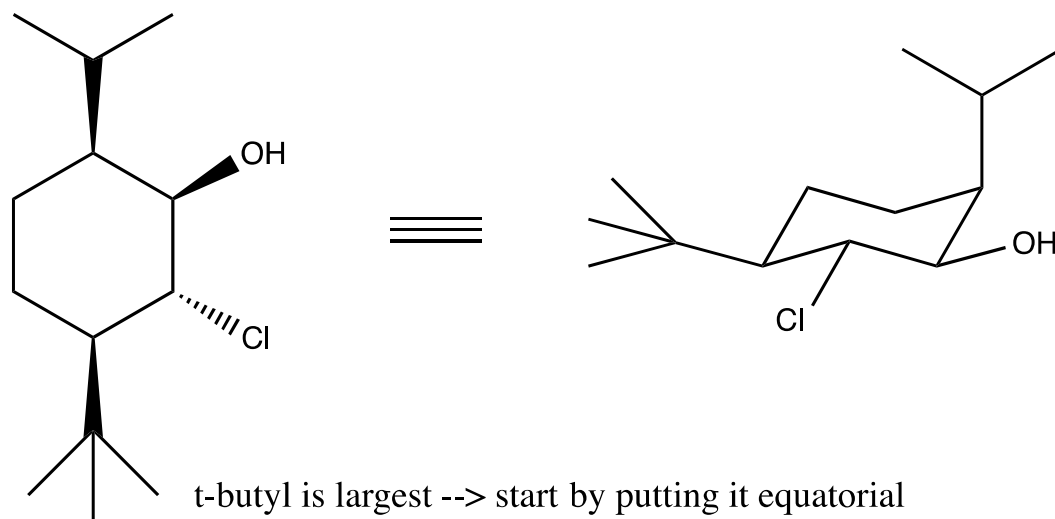
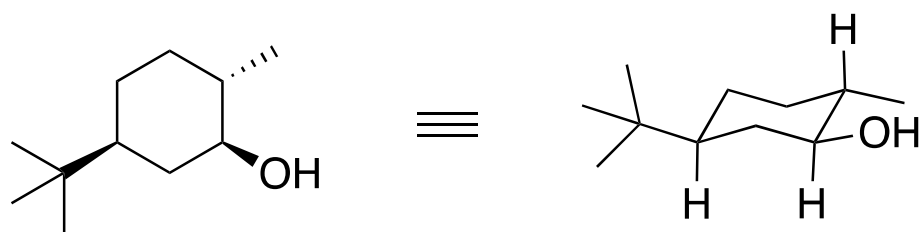


### How to draw the most stable conformation of substituted cyclohexanes:

1. Start by drawing the chair conformation of cyclohexane  
Put the largest group in an equatorial position
2. Draw the next group(s) on the correct atom(s) with respect to the largest group in correct geometry

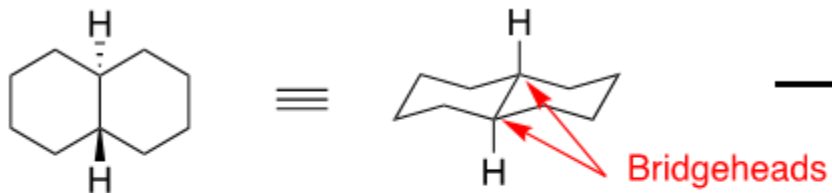


Note that the **largest substituent** (tertiary butyl) is placed in the **equatorial** position to **avoid** destabilizing **1,3-diaxial interactions**

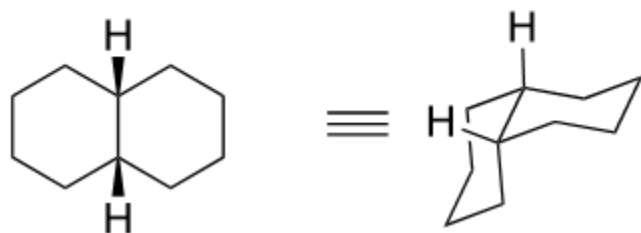
**Another example:****More example:**

**Bicyclic structures**

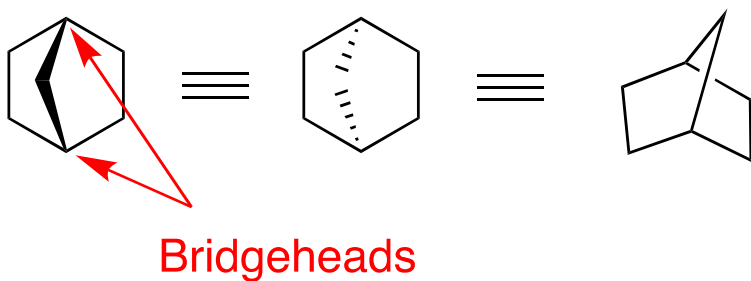
Trans-decalin (4.4.0 bicyclodecane)



Cis-decalin (4.4.0 bicyclodecane)

**diastereomers**  
(NOT conformational isomers)

Norbornane (2.2.1 bicycloheptane)

**Steroids:**