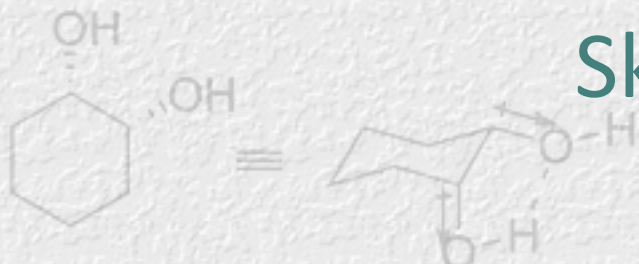
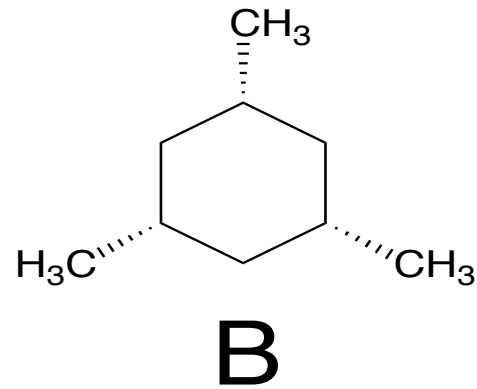
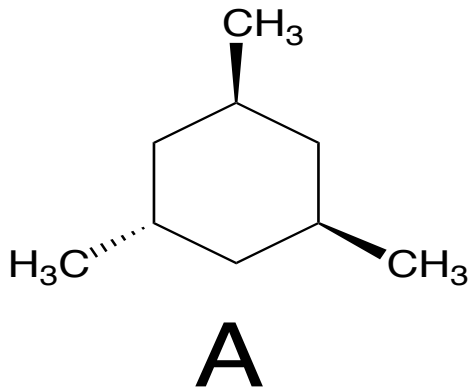


Skillbuilder 2



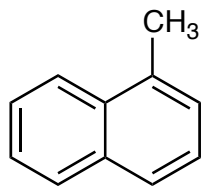
IR = 1714 cm⁻¹

First, draw the most stable chair conformation for each of the following configurational isomers. Second, circle the configurational isomer with the highest potential energy (least stable).

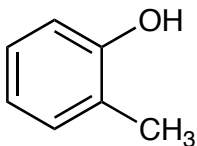


1

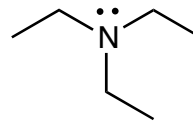
Circle the solutions that would be expected to dissolve each compound. More than one may apply. Circle all that apply.



CH_2Cl_2
3.0 M NaOH (aq)
3.0 M HCl (aq)
 H_2O



CH_2Cl_2
3.0 M NaOH (aq)
3.0 M HCl (aq)
 H_2O



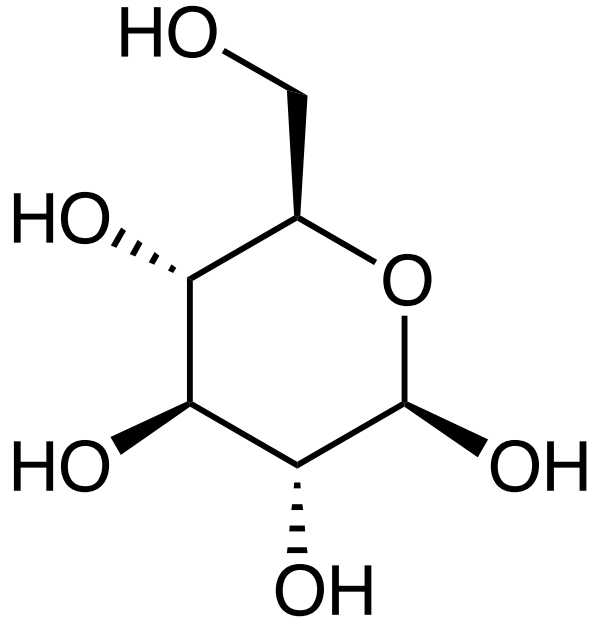
CH_2Cl_2
3.0 M NaOH (aq)
3.0 M HCl (aq)
 H_2O

3

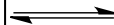
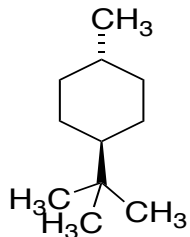
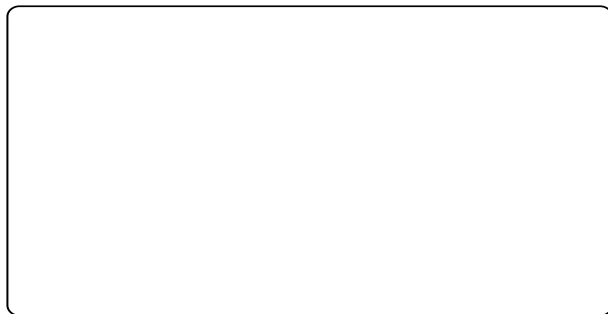
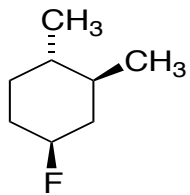
First, for the molecule below, draw a Newman projection for the most stable conformation that sights down C2-C3. **Second**, draw a Sawhorse projection of the least stable *staggered* conformation.

2-methylbutane

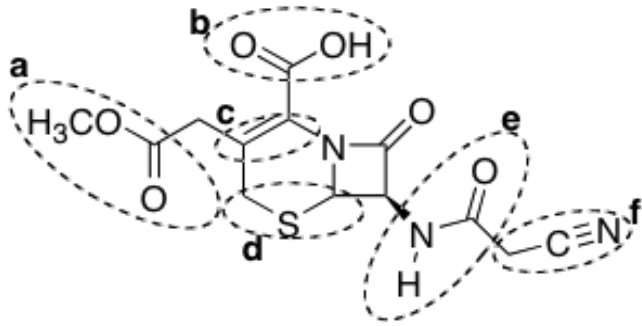
Below is the structure of β -D-glucose. In the most stable chair form, all groups are equatorial. Draw the chair form of glucose.



Draw the two chair conformations for each cycloalkane below. Then, circle the conformation for each configuration that is the most stable.



Below is a naturally occurring antibiotic, a derivative of penicillin. Label each circled functional group.



a _____

d _____

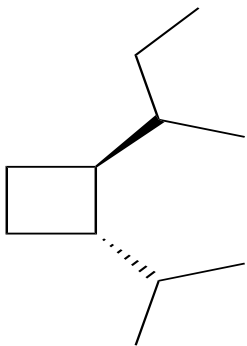
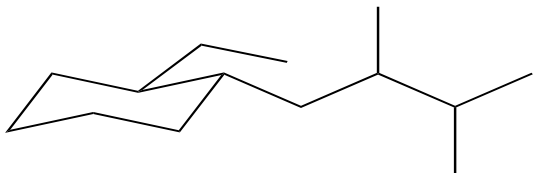
b _____

e _____

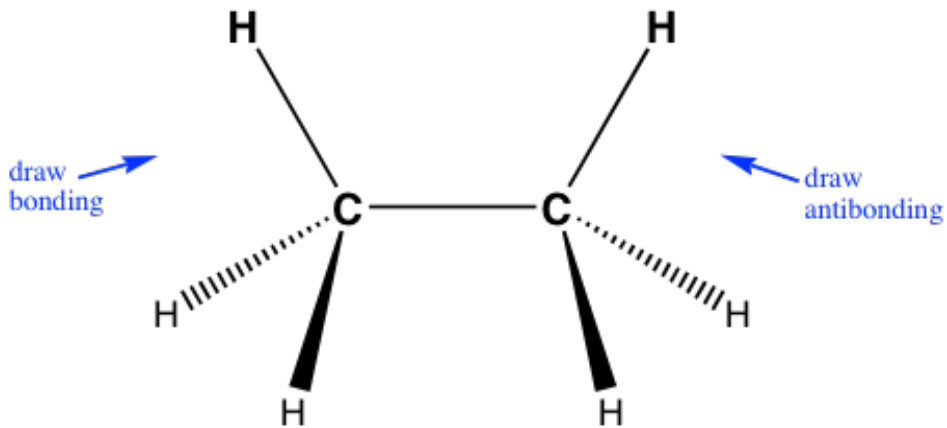
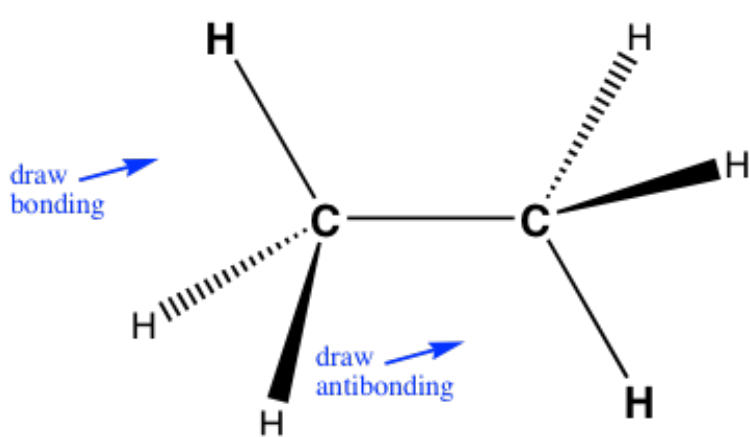
c _____

f _____

Provide an IUPAC name for each disubstituted cycloalkane. Your name should include *cis* or *trans* to describe the configurational relationship between the two groups on the ring.



First, draw the bonding and antibonding molecular orbitals for each of the C-H bonds indicated on top of the given structural formulas. **Second**, circle the structural formula that has the most stabilization by hyperconjugation.

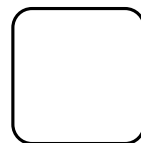
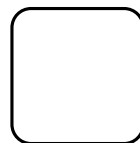
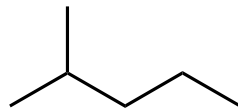
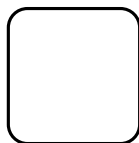
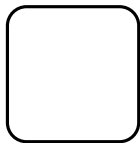
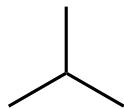
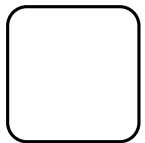
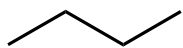


Using wedge-and-dash notation, draw the most polar and the least polar conformation of the following molecule.

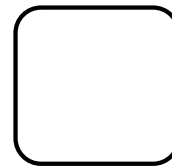
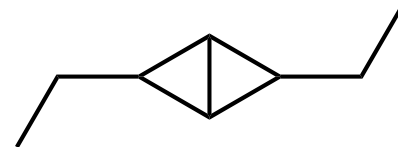
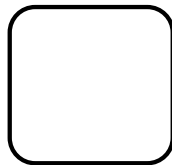
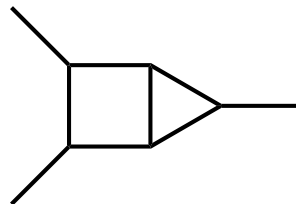
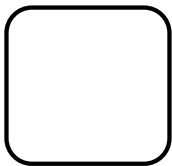
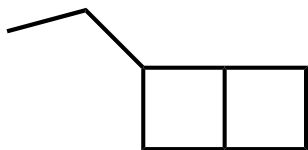
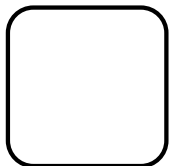
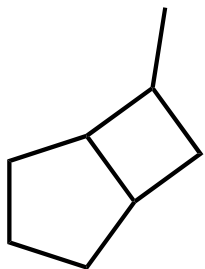
1,2-dibromoethane

10

Redraw the diagram below. Then, rank the molecules in order of increasing heat of combustion (the amount of energy released upon burning) (1 = lowest; 5 = highest).

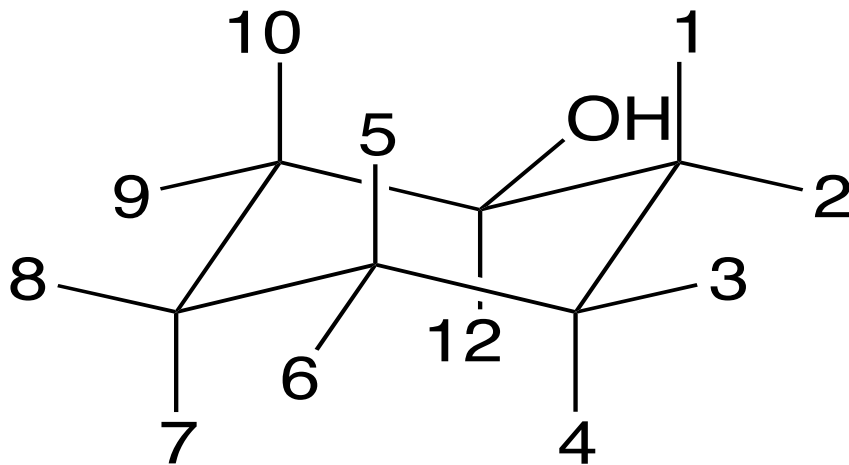


Redraw the figure below. Then, rank the following molecules in order of increasing heat of combustion (1 = lowest; 4 = highest).



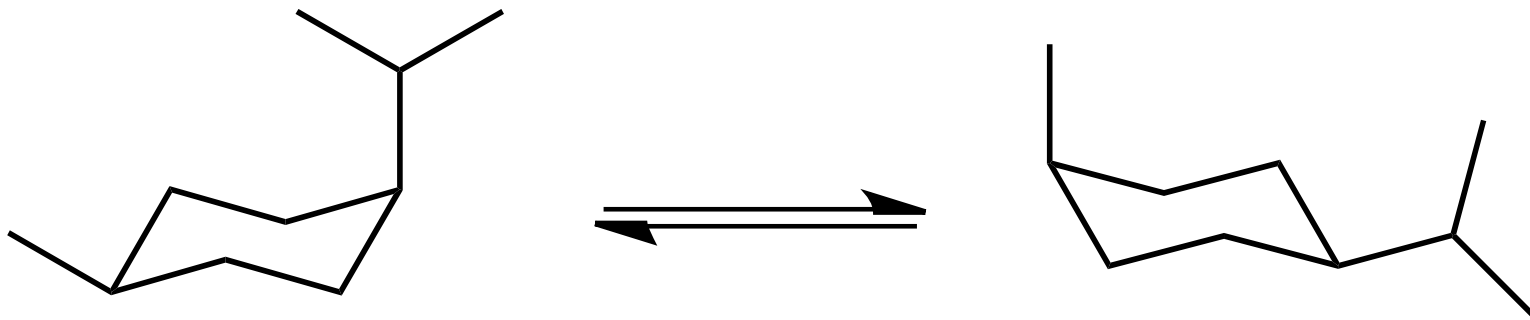
12

In ascending order, list the numbers that are *trans* to the hydroxyl (-OH) group in the cyclohexane molecule below.



13

Below is the ring-inversion equilibrium for cis-1-isopropyl-4-methylcyclohexane. **First**, replicate these structures. **Second**, draw all hydrogen atoms in both conformers that experience 1,3-diaxial interactions with either the methyl group or the isopropyl group. **Third**, circle the most stable conformer.



Draw conformers of butane that fit the descriptions below.

Newman that maximizes torsional strain and minimizes steric strain.



Bond-line (dash-wedge) that maximizes torsional and steric strain.

Draw a picture of an p, sp, sp², sp³ and s orbital. Label each. Make sure the relative size of the lobes indicates the %s-character in each orbital.