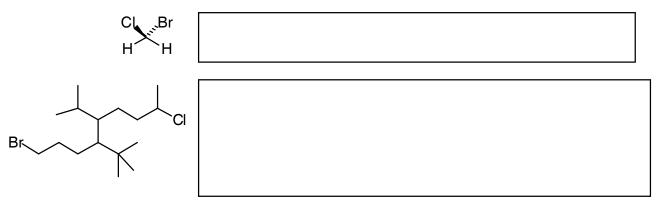
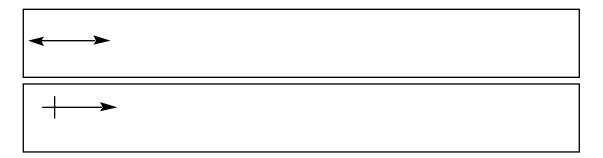
# **Organic Lingo**

1A. Provide a systematic name for the following compounds. Use common nomenclature

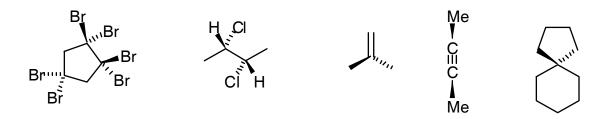


for any branched substituents. (30 pts)

- B. Draw a structure for the following names. For cycloalkanes use flat rings. For all others use bond-line notation.
- 2-cyclobutyl-5,5-dibutyl-2,3,7-trimethyldecane
- 1,1-dibromo-3-isobutylcyclohexane
- trans-1,3-difluorocycloheptane
- C. We have used many types of arrows to describe various phenomena. In the boxes below, identify each arrow using only a FEW words AND provide a REAL example of how it could be used.



D. One or more of the bond-line structures shown below contain improper use of wedgedash notation. Circle the incorrect one(s) AND explain why they are incorrect.



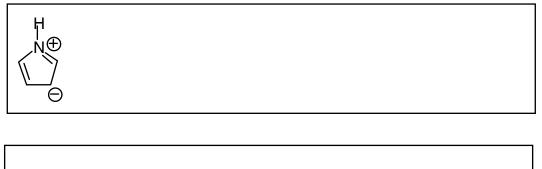
#### The Name is Bond. Molecular Bond.

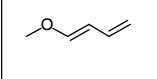
2. (20 pts)

A. Write the Lewis dot structures for each of the following molecules. Be sure to SHOW ALL FORMAL CHARGES AND INDICATE ALL BONDING AND NON-BONDING ELECTRONS AS DOTS (i.e do not use dashes).

		Ŧ
MeOH	H <sub>2</sub> CCHCH <sub>2</sub> Br	H <sub>3</sub> CNCCH <sub>3</sub>

B. Draw two additional resonance structures for each of the following molecules. Your structures must be reasonable.





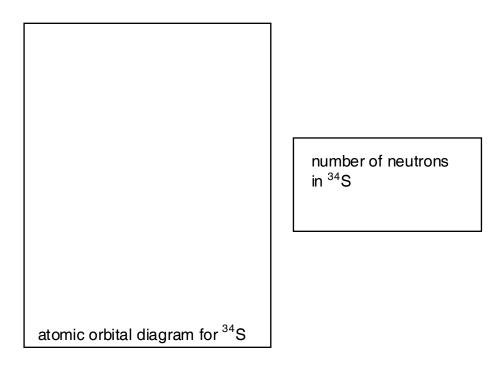
## AO/MO

3. (30 pts)

A. Which term best describes the  $2P_x$ ,  $2P_y$  and  $2P_z$  atomic orbitals (circle one):

valence nodeless Hund's rule

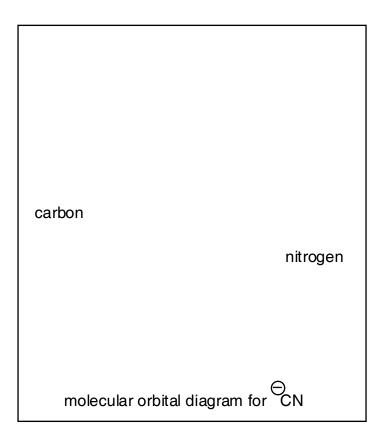
B. Draw the atomic orbital energy level diagram for 1S through 3P. On the diagram fill in the electrons for <sup>34</sup>S (atomic number 16). How many neutrons does <sup>34</sup>S have?



C. Construct the molecular orbital diagram for the cyanide ion, CN, using the following guidelines and labeling schemes. BE SURE TO INCLUDE EVERYTHING ASKED FOR BELOW.

- Assume that both the carbon and nitrogen are SP hybridized.
- The relative energies for carbon and nitrogen are provided below.
- Indicate the relative energy levels of any atomic and hybrid orbitals on the left and right hand sides of the diagram.
- Clearly indicate which orbitals are being combined to make molecular orbitals.
- Fill in all of the electrons.
- Label all the levels as  $\sigma$ ,  $\pi$ , etc.

• Label the HOMO and LUMO



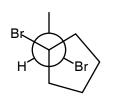
## **Freely Rotating**

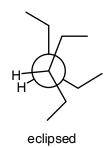
### 4. (30 pts)

A. Draw Newman projections for the bond line structures shown below. Make sure your projection is representative of the conformation shown.

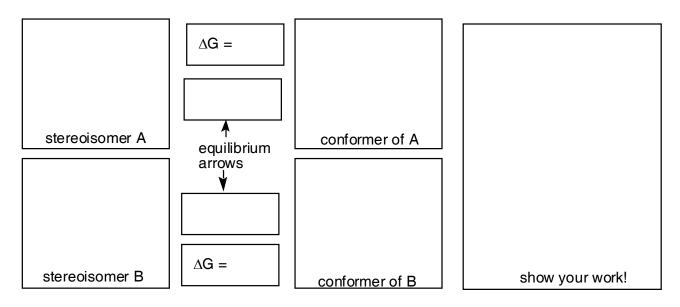


B. Draw bond-line structures, including wedges and dashes, that exactly represent the conformations depicted by the following Newman projections.





Questions 5a-d were on a subject we have not gotten to yet. 5E. In the boxes labeled stereoisomer A and B, draw the CHAIR structures for the two stereoisomers of 1-bromo-1,4-dimethylcyclohexane.

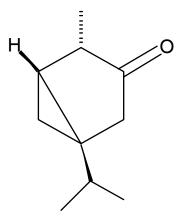


F. In the boxes labeled conformer A and B, draw the ring-flipped conformations of stereoisomers A and B.

G. Calculate  $\Delta G$  for each conformational equilibrium. Place your answer in the boxes labeled  $\Delta G$  =. In order to receive any credit you must show your work. H. Using appropriate arrows indicate which side the equilibrium would lie on for each case.

#### Absinthe

6. Most countries banned the green liquor known as absinthe at the turn of the century due to the presence of thujone, a hallucinogen that, if abused, can lead to serious medical complications. All of the following questions are related to this molecule. (40 pts)





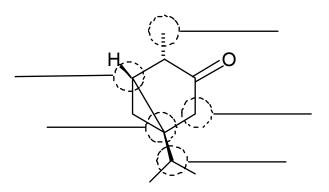
A. Would you describe the relationship between the methyl and isopropyl groups in thujone as (circle one):

cis	trans	resonance	equilibria
	conformers	rotamers	

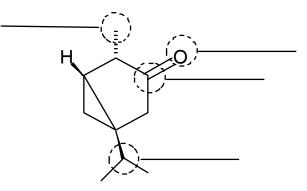
B. Provide a systematic name for the core of thujone.



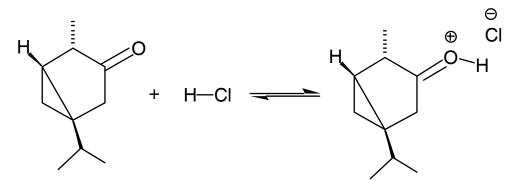
C. On the structure below label each of the circled carbons as primary, secondary, tertiary or quaternary.



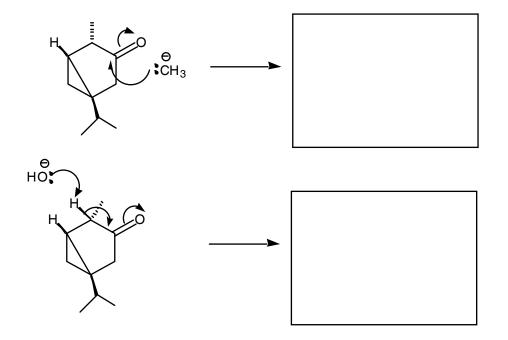
D. On the structure below describe the hybridization at each circled atom.



E. Draw an arrow-pushing mechanism that is consistent with the following reaction.



F. Show the product you would expect to get from the following arrow-pushing mechanism.



C. Below is a potential energy diagram that reflects the relative energies associated with rotation about the  $C_{\P}$ - $C_{\P}$  bond of 2,3,3-trimethylpentane. Match the Newman projection with each minimum and maximum on the diagram (put the letters in the boxes). Be sure to rotate in only one direction. **Hint:** Start by eliminating Newman projections that do not match the molecule in question.

