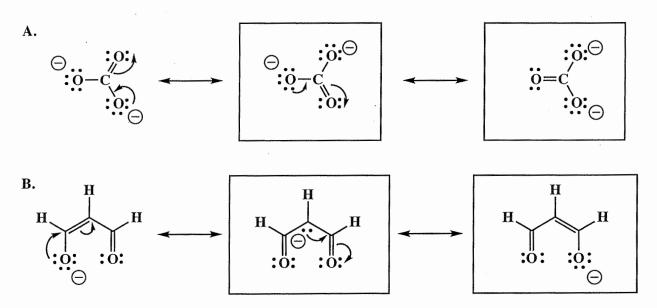
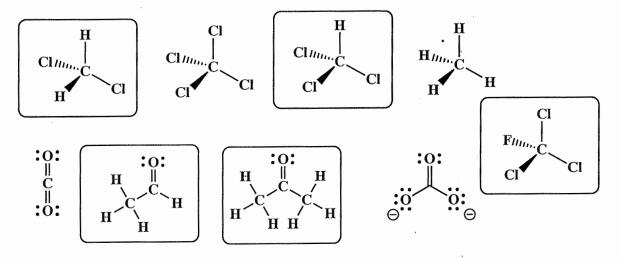
Signature	Pg 5	(29)
Dignatur C	- 0	()

7. (10 pts each) The following molecules are best represented as the hybrid of three contributing structures. Draw the second and third important contributing structures in the spaces provided, including all lone pairs and formal charges. For the two structures on the left in each problem, use arrows to indicate the movement of electrons to give the structures you drew. There is no need to draw any circles around any of these contributing structures. You might want to read these directions again to make sure you know what we want.



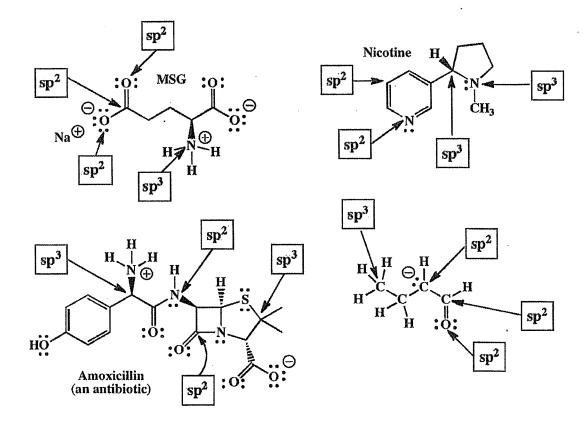
8. (1 pt each) Circle any molecule that has an overall molecular dipole moment.



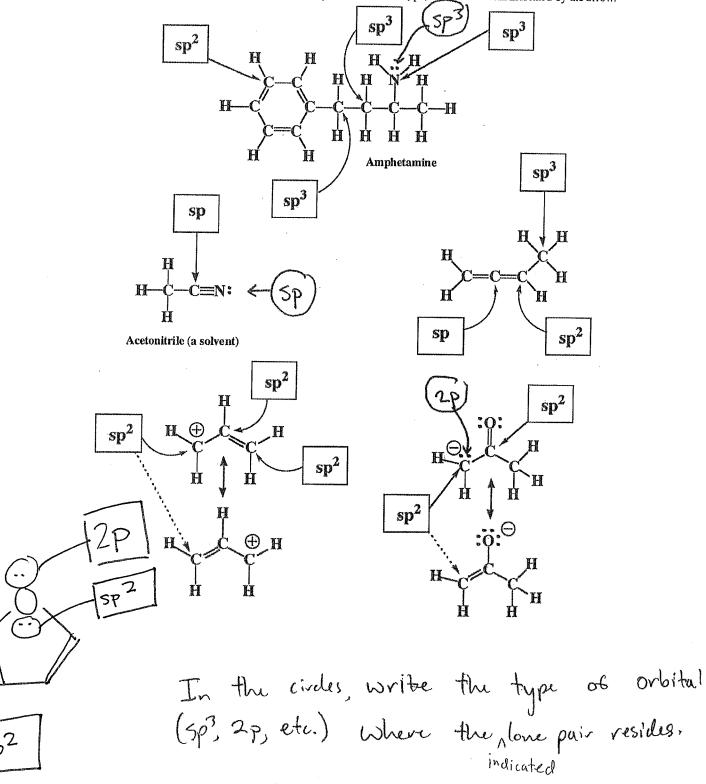
Signature			Pg 6	(30)
8. (2 pts each) Fill in each	blank with the word	or words that best co	ompletes the sentence	es.
For organic chemistry, it is	best to think of	electrons	as waves.	
According to the valence bo	nd approach, the a	atomic orbitals on ea	ch atom are	
combined first to create	hybridized	orbitals, th	at overlap to creat	te
sigma	bonds.			
Three (or more) atom "pi-v	ways" are the situa	tion resonance	contributing	structures are

usually trying to describe. For pi bonding and therefore pi delocalization to occur over more then two atoms (i.e. pi-ways), parallel and overlapping \_\_\_\_\_\_ orbitals are needed on ALL of the adjacent atoms involved. As a result, all of the atoms involved in pi-ways are usually \_\_\_\_\_\_ sp^2 \_\_\_\_\_ hybrized, and NEVER \_\_\_\_\_\_ sp^3 \_\_\_\_\_ hybridized.

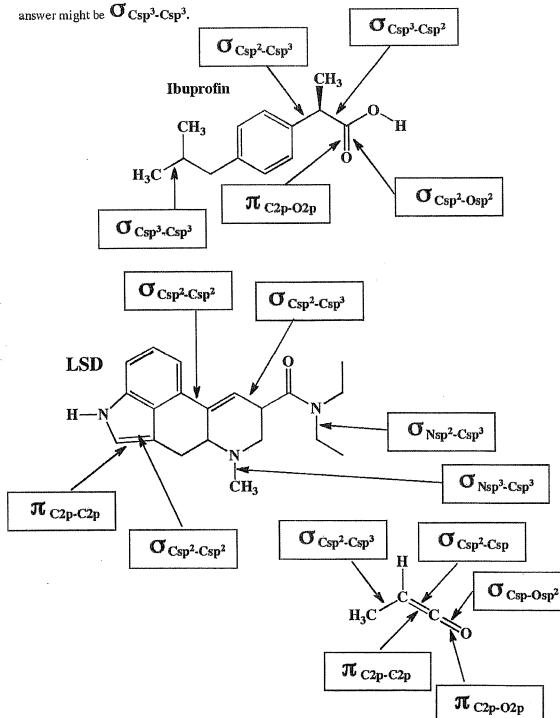
9. (1 pt each) For the following molecules, write the hybridization state of each atom indicated by the arrow.



Signature\_



8. (13 pts) In the box provided, write the hybridization state (sp<sup>3</sup>, etc.) of the atom indicated by the arrow.



**10.** (2 pts each) Describe each bond indicated with an arrow as the overlap of orbitals. For example, an answer might be  $\mathbf{O}$  Csp<sup>3</sup>-Csp<sup>3</sup>.

 **19.** (12 pts) The following three structures have delocalized pi systems that can best be described by the molecular orbital approach based on the overlap of three 2p orbitals. At the bottom of the page are three different molecular orbital diagrams, labeled as "A", "B", and "C". Underneath the structures in the boxes, write the letter ("A", "B", "C"or "D, none of the above") that most accurately describes that molecule's delocallized pi system.

