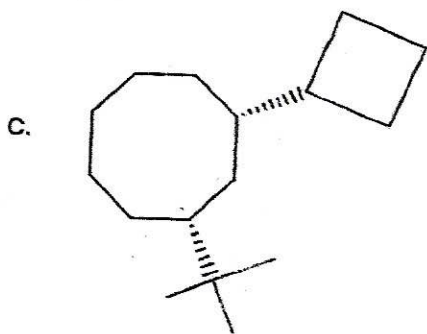
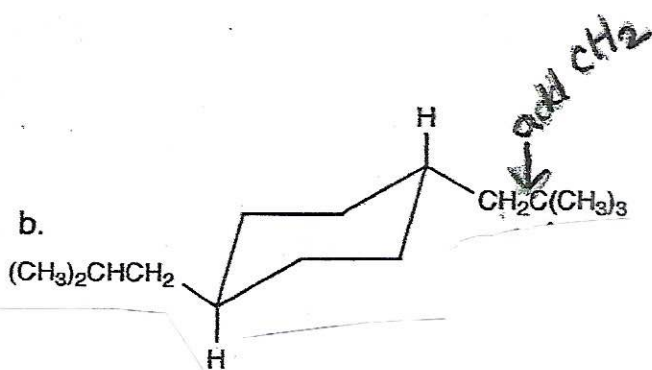
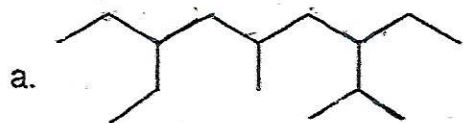


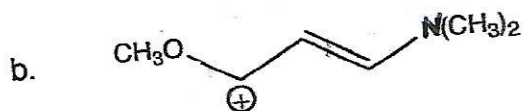
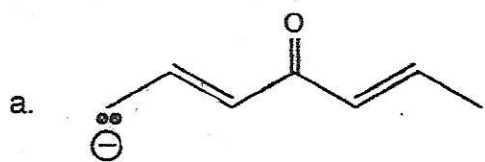
EXAM 1 / Fall 2002

1. NOMENCLATURE: (15 points, 5 pts. each)

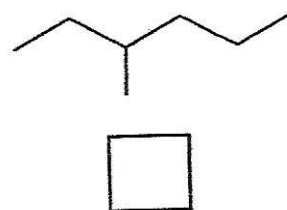
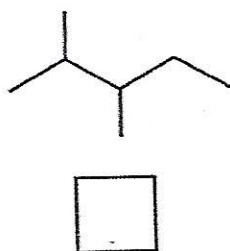
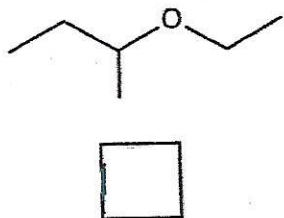
Give an acceptable IUPAC name for each of the following compounds. Be sure to indicate the stereochemistry where appropriate.



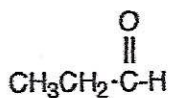
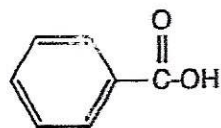
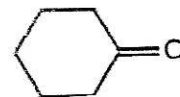
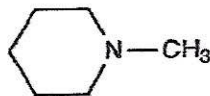
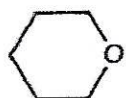
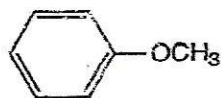
2. For each structure given, draw the important resonance contributors. Circle the major contributor (14 pts.) 2 pts each structure / 1 pt. each circle



3. Place the following compounds in order of increasing boiling point. (1=lowest, 3=highest) (6 pts.)

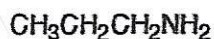


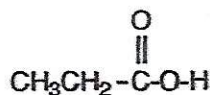
4. Circle the compounds that can form hydrogen bonds in a pure state. (4 pts.)

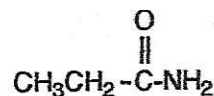


5. Place the following compounds in order of increasing acidity. (1=least acidic, 3=most acidic) (6 pts.)

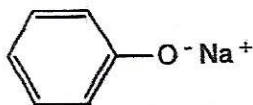
2 pts. each







6. Predict the products that would result from an acid/base reaction between the two compounds below and place the answers in the boxes provided. Predict the direction of the equilibrium, and indicate the direction by placing an arrow in the box. (6 pts.) *2 pts. each*

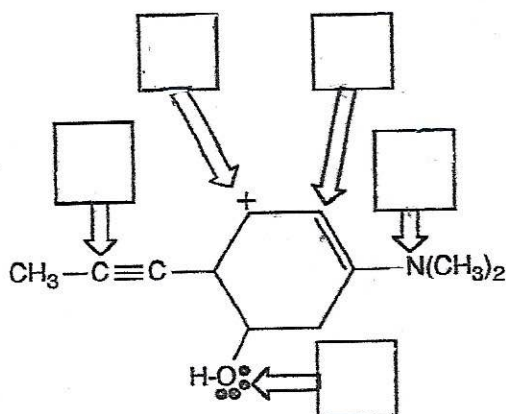


(arrow)

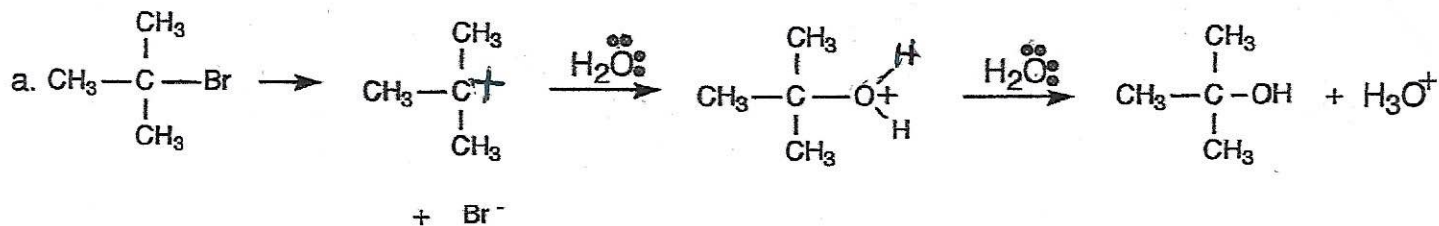
+

7. For the molecule below determine the hybridization of the indicated atoms and place the answers in the boxes provided. (10 pts.)

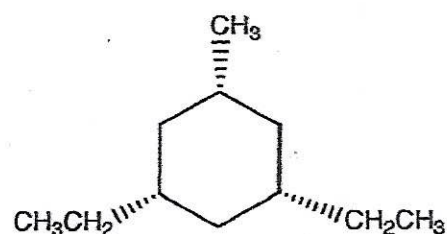
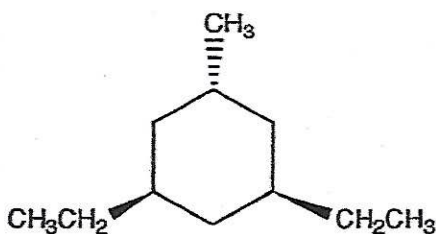
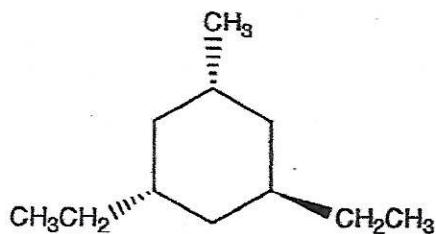
2 pts. each



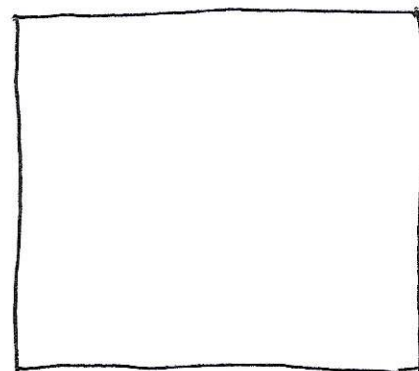
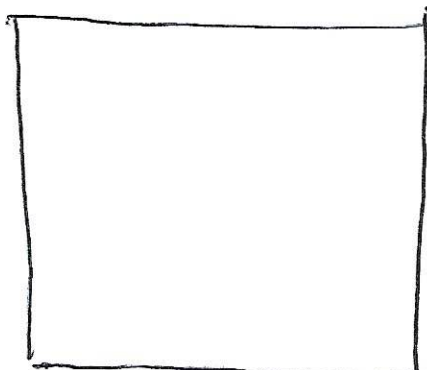
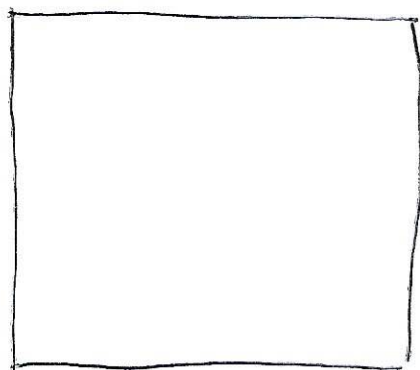
8. For the reactions below, provide curved arrows to indicate the movements of electron pairs associated with bond breaking and bond forming, i.e., provide the mechanisms. (10 pts.)



9. For each of the following substituted cyclohexanes, draw the more stable conformation in the large box. Place the cyclohexanes in order of increasing stability (1=least stable, 3=most stable). (12 pts.)

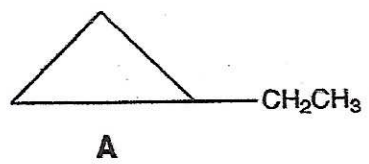


1 pt each

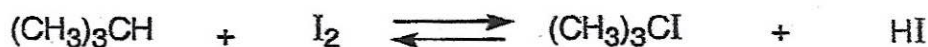


mono

10. Draw all structural isomers and geometric isomers (cis/trans) resulting from the ~~mono~~ bromination of ethylcyclopropane (structure A). Place geometric isomers next to each other and circle them. (11 pts.)



11. Consider the following reaction:



a. Given the bond dissociation energies (BDE) below, calculate the overall ΔH°

(4 pts.)

b. Are the products or reactants favored at equilibrium? (2 pts.)

<i>Bond</i>	<i>Bond-Dissociation Energy</i> <i>kcal/mol</i>	<i>Bond</i>	<i>Bond-Dissociation Energy</i> <i>kcal/mol</i>
<i>H—X bonds and X—X bonds</i>		<i>Bonds to secondary carbons</i>	
H—H	104	(CH ₃) ₂ CH—H	95
D—D	106	(CH ₃) ₂ CH—F	106
F—F	38	(CH ₃) ₂ CH—Cl	80
Cl—Cl	58	(CH ₃) ₂ CH—Br	68
Br—Br	46	(CH ₃) ₂ CH—I	53
I—I	36	(CH ₃) ₂ CH—OH	91
H—F	136	<i>Bonds to tertiary carbons</i>	
H—Cl	103	(CH ₃) ₃ C—H	91
H—Br	88	(CH ₃) ₃ C—F	106
H—I	71	(CH ₃) ₃ C—Cl	79
HO—H	119	(CH ₃) ₃ C—Br	65
HO—OH	51	(CH ₃) ₃ C—I	50
<i>Methyl bonds</i>		(CH ₃) ₃ C—OH	91
CH ₃ —H	104	<i>Other C—H bonds</i>	
CH ₃ —F	109	PhCH ₂ —H (benzylic)	85
CH ₃ —Cl	84	CH ₂ =CHCH ₂ —H (allylic)	87
CH ₃ —Br	70	CH ₂ =CH—H (vinyl)	108
CH ₃ —I	56	Ph—H (aromatic)	110
CH ₃ —OH	91	<i>C—C bonds</i>	
<i>Bonds to primary carbons</i>		CH ₃ —CH ₃	88
CH ₃ CH ₂ —H	98	CH ₃ CH ₂ —CH ₃	85
CH ₃ CH ₂ —F	107	CH ₃ CH ₂ —CH ₂ CH ₃	82
CH ₃ CH ₂ —Cl	81	(CH ₂) ₂ CH—CH ₃	84
CH ₃ CH ₂ —Br	68	(CH ₃) ₃ C—CH ₃	81
CH ₃ CH ₂ —I	53		
CH ₃ CH ₂ —OH	91		
CH ₃ CH ₂ CH ₂ —H	98		
CH ₃ CH ₂ CH ₂ —F	107		
CH ₃ CH ₂ CH ₂ —Cl	81		
CH ₃ CH ₂ CH ₂ —Br	68		
CH ₃ CH ₂ CH ₂ —I	53		
CH ₃ CH ₂ CH ₂ —OH	91		