

## Unit 2 Review for Midterm

- ① Substance A is an ionic solid. It has a relatively high melting point and conducts electricity in liquid state, not solid (Ionic compds conduct electricity when ions can move freely, like in liq.)  
Substance B is a metallic solid. It has high melting point and conducts electricity as solid  
C is molecular substance - low melting point, doesn't conduct electricity  
D is network covalent solid - very high melting point, does not conduct electricity

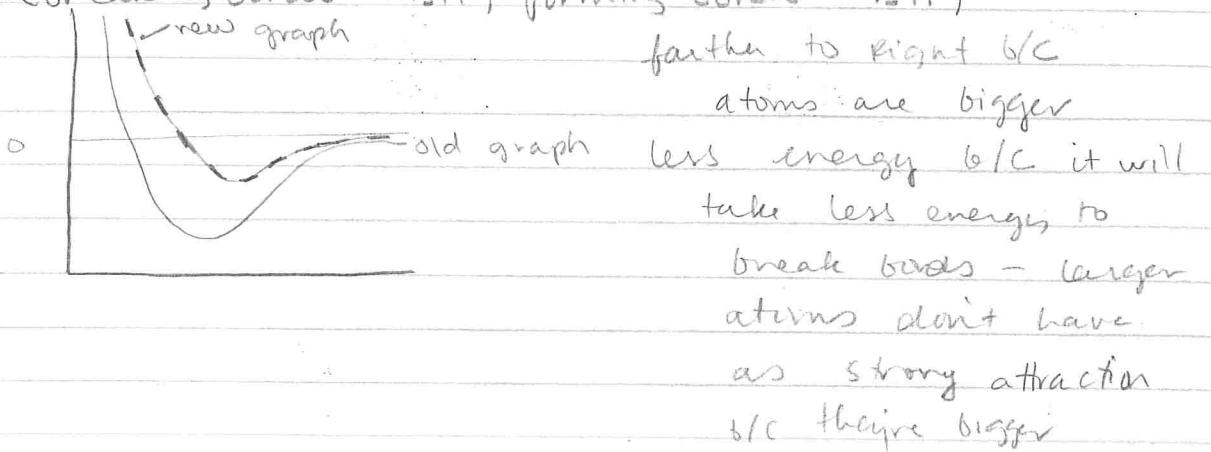
- ② a) A - atoms are too close together + are repelling (nuclei are repelling)  
B - less repulsion; balance between attraction + repulsion  
C - maximum attraction between atoms; this is where the bond will end up; indicates average bond length  
D - atoms are too far apart to have strong attractive force

b) C

c) same energy as point C but will be  $\oplus$

(breaking bonds +  $\Delta H$ ; forming bonds -  $\Delta H$ )

d)

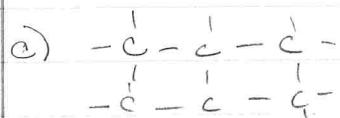




atoms are covalently bonded, molec. are separate



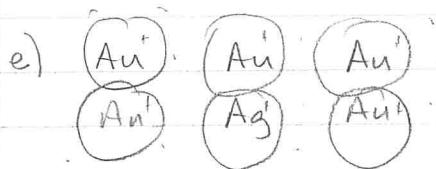
3D crystal structure of repeating  $\oplus + \ominus$  to maximize attraction + minimize repulsion.



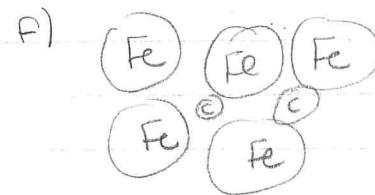
carbon atoms covalently bonded - network-covalent solid



copper ions surrounded by sea of  $e^-$  (metallic bonding)



similar size atoms,  
Ag substitutes for Au  
(and thine metals)



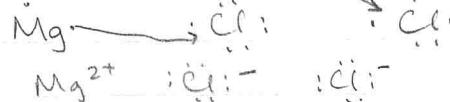
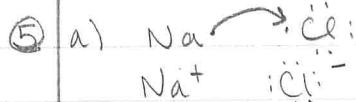
little atoms fit in between bigger Fe atoms



F is more electronegative      / both same



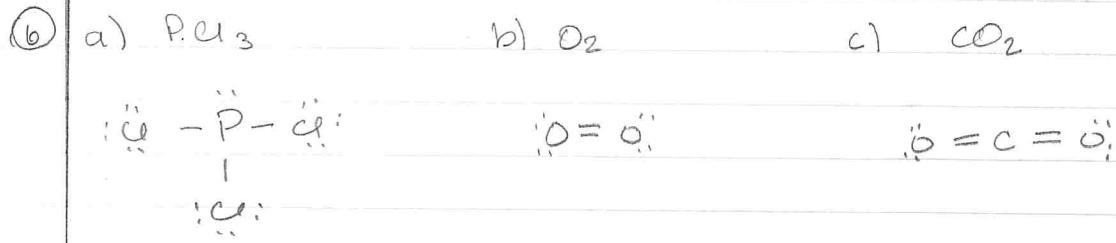
more ionic      more covalent



b) metals give up  $e^-$  - the outermost  $e^-$  are far away from  $\oplus$  nucleus + shielded by other core  $e^-$ , doesn't take much energy to remove, low IE

c) lattice energy - energy required to separate 1 mol solid crystal into gaseous ions.

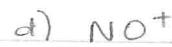
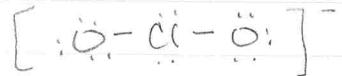
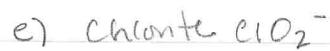
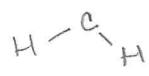
d)  $MgCl_2$  -  $Mg^{2+}$  has stronger attraction force b/c of +2 charge



EDG: tetrahedral  
MG: pyramidal  
angle < 109,5

EDG linear  
MG linear  
—

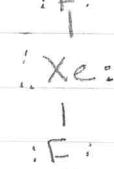
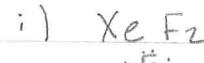
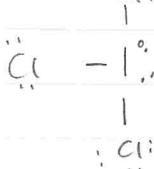
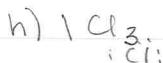
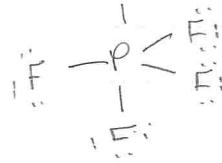
linear  
linear  
 $180^\circ$



EDG: trig planar  
MG: trig planar  
angle:  $120^\circ$

tetrahedral  
bent  
 $> 109,5$

linear

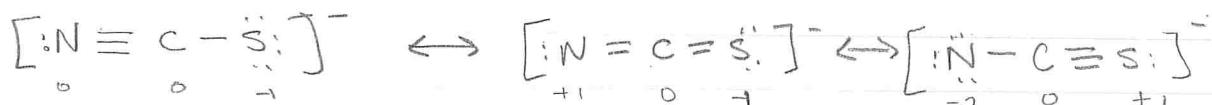


EDG: trig bipyramidal  
MG: trig bipyramidal  
angle:  $90^\circ, 120^\circ$

trig bipyramidal  
T-shaped  
 $90^\circ$

trig bipyramidal  
linear  
 $180^\circ$

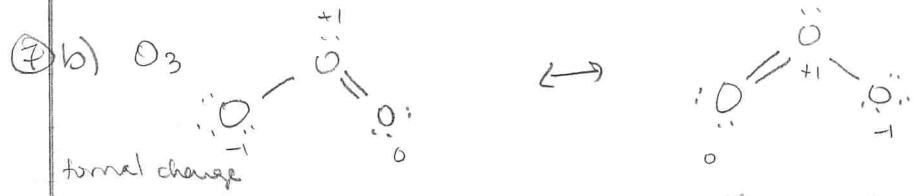
⑦ a)  $\text{NCS}^-$



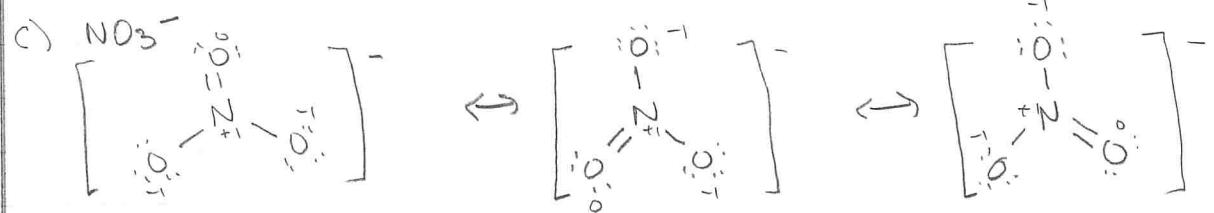
Formal charge

most dominant - least amount of formal charges,

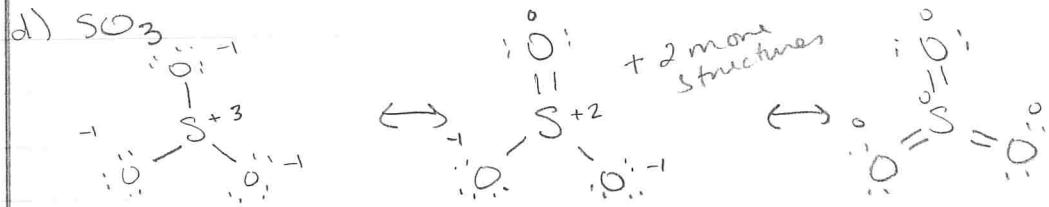
↪ formal charge on electronegative atom



formal charge  
neither is dominant over the other

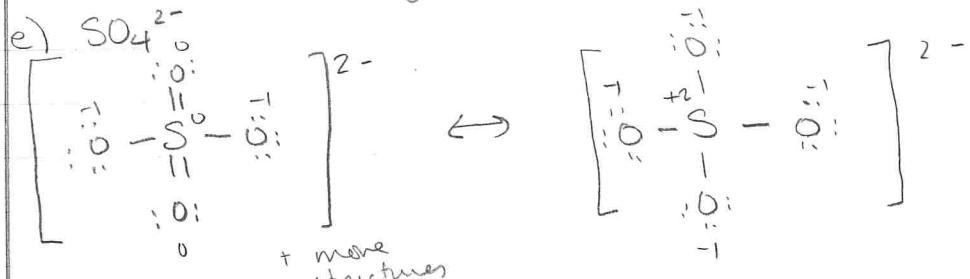


none of resonance structures are dominant

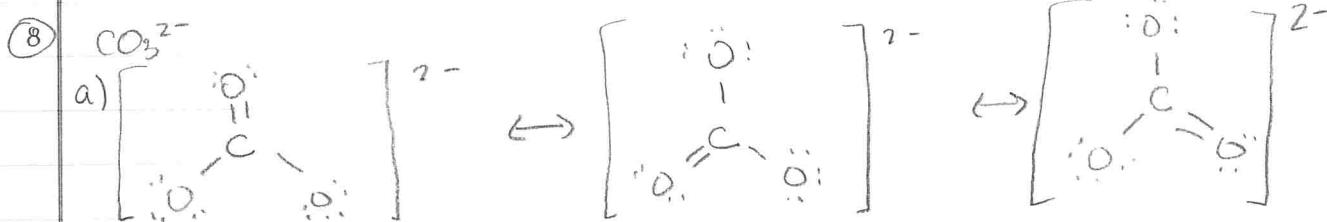


+ 2 more structures  
most dominant

least amt of formal charges

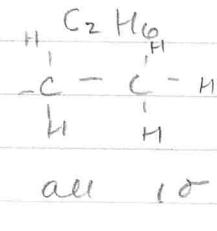
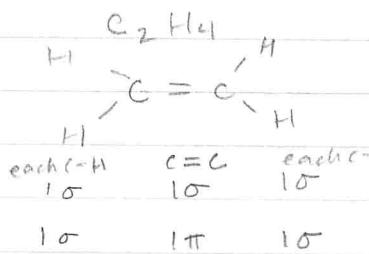
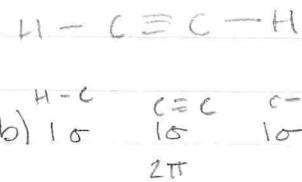


+ more structures  
more dominant, fewer formal charges

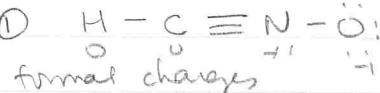


all bonds are same length + same strength  
all resonance structures contribute equally to overall structure, so each C-O bond is like a 1.33 bond.

(9)

c)  $C \equiv C$  shortest, strongest $C-C$  longest, weakestd)  $C \equiv C$   $sp$  $C=C$   $sp^2$  $C-C$   $sp^3$ 

AP Questions



a) the negative formal charge is on the more electronegative atom (O) in the left structure. in the right structure the negative formal charge is on C, which is not the more electronegative atom

b) breaking bonds

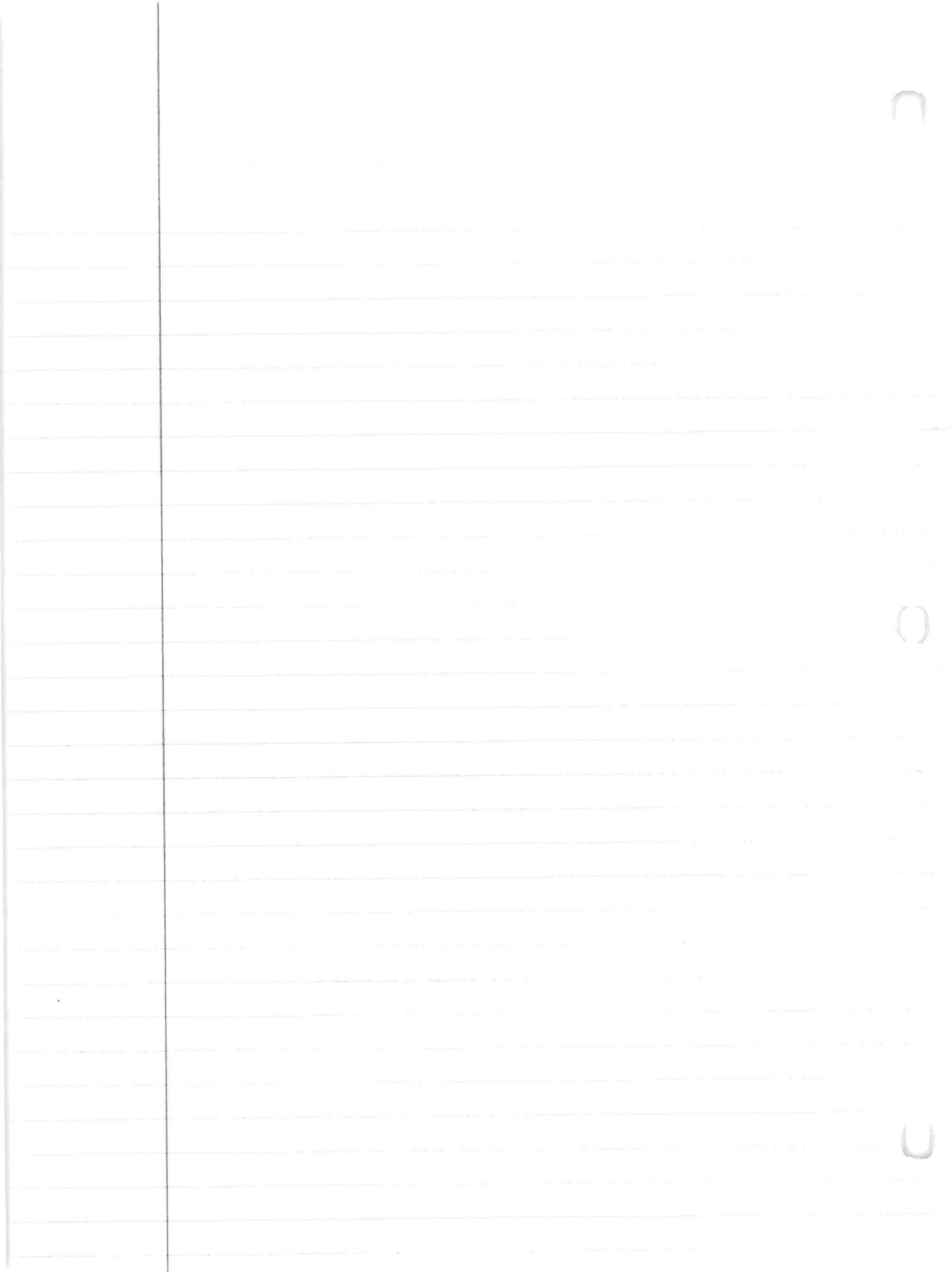
1505  $\xrightarrow{\text{endo}}$ 

forming bonds

1751  $\xrightarrow{\text{exo}}$ 

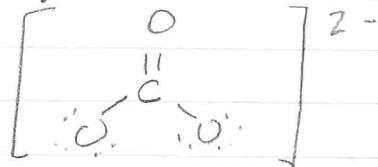
$$\Delta H = \text{bonds breaking} - \text{bonds forming}$$

$$= 1505 - 1751 = -246 \text{ kJ/mol}$$



② a) um... or (typo here?)

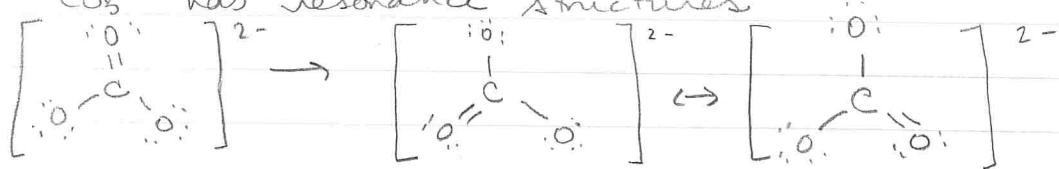
b) (should be i)



c) (should be ii)

each carbon-oxygen bond in  $\text{CO}_2$  is a double bond

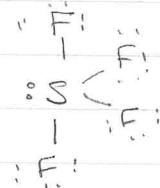
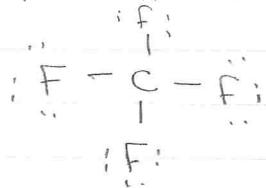
$\text{CO}_3^{2-}$  has resonance structures



the actual structure of  $\text{CO}_3^{2-}$  is a combination of all resonance structures so each carbon-oxygen bond is between a single + double

d) oops again

e) (should be d i)

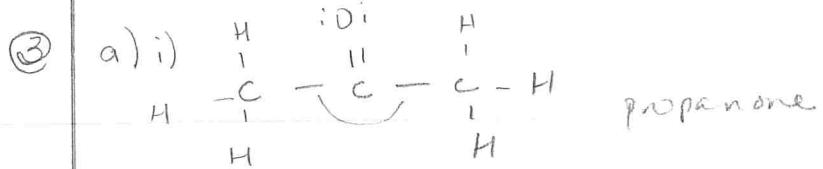


f) (should be d ii)

$\text{CF}_4$  is a tetrahedral molecule that is symmetrical  
the electron density is distributed evenly around the molecule.

$\text{SF}_4$  is seesaw shaped & not symmetrical.

$e^-$  can be pulled to one side of molecule,  
making it polar



ii)  $120^\circ$

b) i) propane vs propanone

19.0 kJ/mol      32.0 kJ/mol

Propanone is a polar molecule w/ dipole-dipole interactions between molecules

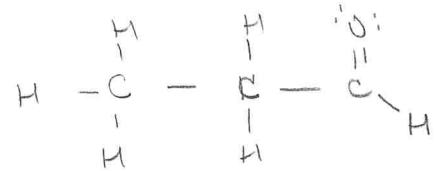
Propane is nonpolar w/ London Dispersion forces.  
Dipole-dipole interactions are stronger than LDF + take more energy to overcome + vaporize

ii) propanone + 1-propanol

32.0 kJ/mol      47.3 kJ/mol

Propanone has dipole-dipole interactions while 1-propanol has stronger hydrogen bonding between molecules, requiring more heat energy to vaporize

c) isomer - same formula, different structure



d) i) sp

ii) 6 total  $\sigma$  bonds (one for each C-H bond  
C-C, and one for  $\text{C}\equiv\text{C}$ )

2  $\pi$  bonds (2 for  $\text{C}\equiv\text{C}$ )

- (4) a)  $\text{C}_2\text{H}_2$        $\text{H}-\text{C}\equiv\text{C}-\text{H}$
- b) Ethyne - triple bonds are shorter than single bonds
- c) i) trigonal planar  
ii) tetrahedral
- d) False. Energy is used to overcome attractive forces between molecules. The covalent bonds within the molecules are unaffected when boiling.
- e)  $\text{CH}_3\text{CH}_3$  (or  $\text{C}_2\text{H}_6$ ) is nonpolar. The electron cloud is distributed evenly throughout molecule.
- f) Ethanol can form hydrogen bonds w/ water molecules. Ethanethiol cannot form H-bonds w/ water

