					page 1	pts
N A M E :						
Chem istry 231 Associate Professor: 1	Or. Gergens	TAKE-HOME EXA DUE 9/26	AM 1		FALL	2001
1. (12) Circle the m	ost (highest); un	derline the least (bwest)				
basiz	NH2	PI	4 <sub>2</sub>		SH <sup>-</sup>	
acidic	CH <sub>3</sub> CH <sub>2</sub> OH	CH3	3CO2H		CH <sub>3</sub> NH <sub>2</sub>	
preferred side of equilibrium	Ph <del>-</del> O-	+ CH <sub>4</sub>		Ph-O-H	+	CH3
highest boiling isom er	$\frown$	СН2ОН		CH <sub>2</sub> NH	I <sub>2</sub>	NH
stable chair structure for $C_7H_{13}C1$		H H H CH <sub>2</sub> CH <sub>2</sub> H H H H H CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>3</sub>	H CI	$ \begin{array}{c} \begin{array}{c} \begin{array}{c} \begin{array}{c} \begin{array}{c} \begin{array}{c} \end{array}\\ \end{array}\\ \end{array} \\ \begin{array}{c} \end{array} \\ \end{array} $		$\bigcup_{CH_2}^{CH_2} \bigoplus_{CH_3}^{H} \bigoplus_{H}^{H}$
electronegative atom	С	0	Ν		I	
2. (2) In the follow	wing, circle <u>ALL</u> c	om pounds that are covale	nt:			
(NH <sub>2</sub> ) <sub>2</sub> CO	CH <sub>3</sub> CO <sub>2</sub> Na	CO <sub>2</sub>	CH	HBr <sub>3</sub>	ammonium	bromide
3. (2) Draw in the positive, $\delta^{\dagger}$	direction of the b , and which is pa	ond dipole m om ent and incartial negative, $\delta^{-}$ :	licate whic	h end of the bo	nd is partia	1
		Br-Cl				
4. (2) Complete the	following for car	bon-14.				
Num ber of neutrons:	N	um ber of protons:	1	Numberofelect	crons:	

5. (4) A typical covalent bond is worth between \_\_\_\_\_ and \_\_\_\_\_ kcals/mole in bond strength. Nonbonding interactions are considerably weaker. A hydrogen bond is \_\_\_\_\_ kcals/mole, a dipole-dipole interaction is \_\_\_\_\_ kcals/m ole, and a dispersion force is \_\_\_\_\_ kcals/m ole.

### Structure and Bonding

6. (2) Mark with an "X" all form ulas that are <u>NOT</u> valid resonance structures of diazom ethane, CH<sub>2</sub>NN.



- 7. (2) For each valid structure in question 6, indicate the form alcharge on each atom.
- 8. (3) Between the valid structures of  $CH_2NN$  in question 6, which structure is the major contributor and why?

9. (3) Using atom is orbitals, draw a three-dimensional representation (using wedges and dashes lines) of the structure for the major contributor in question 6, showing how the sigma and pibonds are formed and where the electrons are located.

10. (4) Forthe compound bebw.	a. Draw the line-angle structure.
H <sub>2</sub> NCH <sub>2</sub> (CH <sub>2</sub> ) <sub>2</sub> CH <sub>2</sub> O $\underbrace{\bigcap_{C CHCHCH_2C}^{O}(CH_3)_3}_{1}$	
0 1 2 3 4 5 6	b. How many non bonding electron pairs are in this molecule?
0 1 2 3 4 5 6	c. What is the degree of unsaturation in this molecule
0 1 2 3 4 5 6	d. How many tertiary centers are in this molecule?
0 1 2 3 4 5 6	e. How m any atom s in this m olecule have sp <sup>2</sup> hybrization?
$\pi(p^{-}p)  \sigma(p^{2}-p^{2})  \sigma(p^{2}-p^{3})$	f. The bonding molecular orbital between the carbonyl and the CH, identified as number 1.

- 11. (3) Valence Bond Theory
- a. Based on the experimentally determined bond angle for water, the predicted H-S-H bond angle in H<sub>2</sub>S would be 104.5°. The actual experimental bond angle in the H-S-H bond angle in H<sub>2</sub>S has been experimentally determine to be 92°.
- b. Draw the valence bond picture for the experimentally observed bond angle for  $H_2S$ .



13. (2) What are the predicted bond angles for the num bered atom s in AZT: 1.\_\_\_\_; 2.\_\_\_\_; 3.\_\_\_\_.



(CH $_3$ CH $_2$ ) $_2$ C (CH $_3$ ) CH $_2$ CH (CH $_3$ ) $_2$	



#### 15. (6) <u>Conformation</u> Analysis

Draw the Newman projection for 2-m ethylbutane about the  $C_2-C_3$  bond in the eclipsed, staggered, and skewed conformations and kbelallgauche interactions. Draw a graph of the torsional strain observed in 2-m ethylbutane as it rotates about the  $C_2-C_3$  bond. Show the dihedralangle and calculated torsional strain for each conformation.



#### 16. (6) Chair Conform ations

## a. Draw the chair conform ation for compound in the box at the left.



- b. Circle the most stable conformer.
- c. Determ ine the energy difference between two chair conformers and calculate the equilibrium concentration of each conformerat room temperature.

17. (2) Label the mactants in these acid-base mactions as Lew is acids (electrophiles) or Lew is bases (nucleophiles).



18. (2) Use curved arrows to show the movement of electrons in the Lewis acid-base meactions of quetion 17.

19. (2) Complete the resonance structures in quetion 17.

20. (4) Complete the meaction and classify the meaction type as either substitution, addition, acid-base, or elimination. If the final product has resonance forms, draw them.

Reaction Type:



21. (2) Use curved arrows to show the movement of electrons in the Lewis acid-base reactions of questions 18.

22. (4) Using your knowledge of Lewis acid-base chemistry, draw a possible mechanism for the formation of carbonic acid. Show all steps, resonance structures, formal charge, and arrow pushing in your mechanism. Proceed step-wise. You should have a minimum of three individual steps in the mechanism.

NAME:\_\_\_\_

Chem istry 231

FALL 2001

Associate Professor: Dr. Gergens

1. (6) Name the following compounds:





2. (18) Mutiple Choice - Place your answer in the box at the right.

a.	A molecule of $NH_2NH_2$	has hybrization about	ut each n <b>ì</b> rogen atom .		
<b>A</b> .	sp <sup>3</sup>	B. sp <sup>2</sup>	C.sp	D. p <sup>3</sup>	

b.	A ion of sulfite, $SO_3^{2-}$	has hybrization abo	out the sulfur atom		
A	. sp <sup>3</sup>	B. sp <sup>2</sup>	C.sp	D. p <sup>3</sup>	<u> </u>

c.	Of the molecules listed	, the	one that would be line	ar is				
<b>A</b> .	BeF2	в.	SO <sub>2</sub>	c.	0 <sub>3</sub>	D.	H <sub>2</sub> S	

# d-g



# d. Indicate the hybridization about atom 1

•	Indicate th	e <u>geom et</u> :	ry about atom	1.			
Α.	linear B.	bent C.	trigonal plan	nar D. p	yramidal E	. tetraheo	dral

2'-deoxythymidine 5'-phosphate

g. The degree of unsaturation in this molecule is

A. 2	B. 3	C. 4	D. 5	E. 6	1979 1979 1979 1979 1979 1979 1979 1979

 h. The predicted bonding molecular orbital between the carbon atoms indicate by the curved arrow attached to number 2;

A. 
$$\pi(p^-p)$$
 B. $\sigma(p^-p^2)$  C. $\sigma(p^2-p^3)$  D. $\sigma(p^3-p^3)$ 

i. Draw in the direction of the bond dipole moment and indicate which end of the bond is partial positive,  $\delta^+$ , and which is partial negative,  $\delta^-$ :

## 3. (20) Circle the most (highest); underline the least (lowest)



4. (4) For the compound bebw.	a. Draw the line-angle structure. b. Show all non bonding electron pairs.
CH <sub>3</sub> COCO <sub>2</sub> H	
pyruvic acid	

- 5. (6) a. Name the compound in the box bebw :\_
- b. W hat is the difference in free-energy between the two form s?



- c. Which side of equilibrium is preferred?
- 6. (6) For each meaction, indicate the Lewis acid and base. Complete the meaction and use curved arrows to show the movement of electron pairs (i.e., arrow pushing of electrons). Show all form all charges and mesonance structures in the product.



NAME:					
Chem istry 231 Assistant Professor: D.D.Gergens, Ph.D.		<b>EXAM 1</b> 9/23		FALL 1997	
1. (18) Circle the m	ı ost (highest); underlin	e the least (bwest)			
polarbond	F-F	I- F	<del>I</del> Br		
electronegative atom	0	F	S Be		
basic	сн <sub>3</sub> соо <sup>-</sup>	CH3 <sup>-</sup>	NH	1 <sup>+</sup>	
highest boiling isom er	. СН <sub>3</sub> —О—СН <sub>3</sub>	CH <sub>3</sub> -N- H	CH <sub>2</sub> CH <sub>3</sub>	H-O-CH <sub>2</sub> CH <sub>2</sub> CH <sub>3</sub>	
preferred side of equilibrium	Ph <b>-</b> OH +	NH2	Ph-O	+ NH <sub>3</sub>	
stable conform er	CH <sub>3</sub>	CH <sub>3</sub>	CH <sub>3</sub>	CH <sub>3</sub> CH <sub>3</sub>	
stable conformer	CH <sub>3</sub>	СН <sub>3</sub>	CH <sub>3</sub>	CH <sub>3</sub> CH	

2. (4) In the following, circle <u>ALL</u> compounds that are ionic:

N	H <sub>4</sub> HCO <sub>3</sub>	CH <sub>2</sub> Cl <sub>2</sub>	CO <sub>2</sub>	LiCH <sub>3</sub>	sodium acetate
3. (8)	Nam e or draw	the following compou	inds		
	CH <sub>3</sub>   CH <sub>3</sub> CCH <sub>2</sub>   CH <sub>3</sub>	CH <sub>2</sub> CH (CH <sub>3</sub> ) <sub>2</sub>   CH <sub>2</sub> CCH <sub>2</sub> CH <sub>3</sub>   CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>3</sub>			
				trans-1-sec-buty-2-i	sopropykycbhexane