George Pappas CHM. 121.004 Inst. Mark Novak Lab Report – Experiment 10 11/16/17



### **Experiment Objective**

The objective of this experiment was to use the concept of Lewis dot structures to find the bond angle, hybridization, shape name, and overall polarity of a molecule.

### Conclusion

This was a very straightforward lab. We were able to implement what we learned about Lewis structures, bond angle, hybridization, shape name, and polarity. The part of the lab that took the longest was creating the dot structures. After creating the dot structures, the rest was simply plugging in the rest of the data using the given dot structure chart. The only resonance structure that I noticed was for the SO<sub>2</sub> molecule; the double bond is just switched to the other side. I then started to do the polarity of molecules worksheet. This was also pretty straight forward; the lab manual gives the nonpolar, polar, and ionic bond ranges. After finding the electronegativity difference, then it's simply finding which range the bond belongs in. Overall, the lab was fairly easy and straightforward.

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### CHM121 E10 STRUCTURE AND POLARITY OF MOLECULES

Example: In  $H_3PO_4$ , there is a P-O bond. Oxygen has an electronegativity of 3.5 and phosphorus has an electronegativity of 2.1

$$O = 3.5$$
 $- P = 2.1$ 

Electronegativity Difference = 1.4

Therefore, the P-O bond is a POLAR bond.

Here are the electronegativities of some elements:

bromine 2.8

fluorine 4.0

oxygen 3.5

carbon 2.5

hydrogen 2.1

sulfur 2.5

chlorine 3.0

nitrogen 3.0

Complete the following table, calculating the electronegativity difference and determining the bond polarity.

Bond	Electronegativity Difference	Bond Polarity	Bond	Electronegativity Difference	Bond Polarity
C – H	.4	nonfolar	0 - 0	0	Numpular
C – Cl	.5	Polar	N – N	0	nontolar
C – O	~	Polar	C – C	0	nonPolar
O – H	1.4	Polar	C – F	1.5	Polar
H – F	-1.9	Ionic	S - O	- 1	Polar
N - H	.9	Polar			

**B3)** After you have drawn the dot structure for each substance, look at the center atom in each one. Count how many atoms are directly bond to the center atom, and also how many spare electron **pairs** are on the center atom.

Add those two numbers together in order to get the "SUM" listed on the "DOT STRUCTURE CHART". From the "DOT STRUCTURE CHART" and the sum calculated, complete the "Second Chart On The Data Sheet", except for Overall Polarity.

Finally, follow the directions in the section "Predict From Lewis Dot Structure", to determine the overall polarity of each substance in the "Second Chart On The Data Sheet" and also for all substances in Part C.

### CHM121 E10 STRUCTURE AND POLARITY OF MOLECULES

### Part C Working with ChemSketch

In this part of the experiment you will create space-filling models of six compounds from the ChemSketch program. Then you will print out the color picture of each of the compounds on separate sheets. Finally, you will add, computer printed from a blank template you have to make, the molecular formula, overall polarity (POLAR or NONPOLAR), your name in the upper right-hand corner after being stapled (Of Course!) and the correct dot structure. If necessary, you may print the correct dot structure in ink, instead of computer printout. However, everything else must come from the computer.

The six compounds are:

1) cis (Z) -C<sub>2</sub>H<sub>2</sub>Br<sub>2</sub>

2) trans (E) -**C<sub>2</sub>H<sub>2</sub>Br<sub>2</sub>** 

3) no-name- $C_2H_2Br_2$ 

4) **CH<sub>4</sub>O** 

5) **SO**<sub>2</sub>

6) CH<sub>2</sub>O<sub>2</sub>

Even though two of the six compounds are included in the previous chart, sulfur dioxide and methanol, you still have to include their molecular formula, dot structure and overall polarity on the individual space-filling model picture sheet. Be aware that the three carbon-carbon double bond isomers **are NOT the same compounds** as those in the previous chart.

To start, first draw a correct Lewis dot structure for each compound. Since you can't draw a correct space-filling model without a correct dot structure, you may want to check your dot structures with your instructor. Once you have the correct dot structures, then you can determine if the bonds are single, double or triple, and the correct bond angles.

### **USING CHEMSKETCH**

Turn on the computer. You don't have to log on to the Chemistry laptops, but you may need to if you use another computer on campus. Double-click on ChemSketch icon on the desktop. Then also click on the OK in "Did you know?" tip box

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PART B
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<b>CHM121</b>

				NAME		
	OVERALL POLARITY (P or NP)	N	1)4	<u>a</u>	۵_	C
Ship except for Account	ISOELECTRONIC RESONANCE OR ISOMERS					
SKip	SYMMETRY (S or U)					
) CHM121 E10 PART B CHART	SHAPE NAME	Tetrobedia 1	Te fear poles i	Tetrahrdial	Bent (v-shafed)	Trigonal
121 E10	HYBRID- IZATION	2	Contraction of the Contraction o	25	200	202
СНМ	BOND ANGLE	50	109.5	109.5	109.5 503	06
	lews Streeture DOT STRUCTURE		:10-3-H	H-0-0-1	;) ( H	о; 11 — 0 — Н
×	FORMULA	CH <sub>4</sub>	CH <sub>2</sub> Cl <sub>2</sub>	CH40	H <sub>2</sub> O	CH <sub>2</sub> O <sub>2</sub>
	#1	1	2	m	4	rv ,

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# CHM121 E10 PART B CHART

			NAME		
OVERALL POLARITY (P or NP)	d	0	0_	2	NP
ISOELECTRONIC RESONANCE OR ISOMERS					
SYMMETRY (S or U)					
SHAPE NAME		Trigonal Pyramid	Bent (v-shuped)		Trigonal Plante
HYBRID- IZATION		E & S	25		52
BOND		109.5 513	2,803		70
DOT STRUCTURE	H - F.		: 'S' = 0:	:N=N:	H, C=C, H
FORMULA	HF (+)	NH <sub>3</sub>	H <sub>2</sub> O <sub>2</sub>	N <sub>2</sub>	C <sub>2</sub> H <sub>4</sub>
#1	9	7	<b>∞</b>	ത	10

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# CHM121 E10 PART B CHART

OVERALL POLARITY (P or NP)	<u>_</u>	X	<u> </u>	2	Commen
ISOELECTRONIC RESONANCE OR ISOMERS					:0:
SYMMETRY (S or U)					
SHAPE NAME	Trigonal	Trigonal	Jonal Denal	Linear	Best (U-Shaped)
HYBRID- IZATION	28	Sp	200	S	5,22
BOND	120	6	30	180	04
DOT STRUCTURE		T	T ) = )	H - 0 = 0 - H	
FORMULA	cis- C <sub>2</sub> H <sub>2</sub> F <sub>2</sub>	trans- C <sub>2</sub> H <sub>2</sub> F <sub>2</sub>	no- name- C <sub>2</sub> H <sub>2</sub> F <sub>2</sub>	C <sub>2</sub> H <sub>2</sub>	SO <sub>2</sub>
#1	11	12	13	14	15

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# CHM121 E10 PART B CHART

<u> </u>	FORMULA	DOT STRUCTURE	BOND	HYBRID- IZATION	SHAPE NAME	<u>SYMMETRY</u> (S or U)	ISOELECTRONIC RESONANCE OR ISOMERS	OVERALL POLARITY (P or NP)
SC	<sup>16</sup> SO <sub>4</sub> - <sup>2</sup>							
3	VE=32	:0: -5-0:	109.5	109.5 5p3	1 ftrantdral			2
17 CO <sub>2</sub>	CO <sub>2</sub>	:0:0::	0 \$	3	Lineal			5

NAME