



Name All (Review)

Period All

College Prep Chemistry

Assignment 3M – Unit 3 Review

30 Points

Complete the four forms of the dot structure for the following elements

group gives  
# val e<sup>-</sup> (dots)

| Element Group      |  |  |  |  |
|--------------------|--|--|--|--|
| SA                 |  |  |  |  |
| Val e <sup>-</sup> |  |  |  |  |
| 5                  |  |  |  |  |

1A = 1 4A = 4

5A = 5 6A = 6

7A = 7 8A = 8

Draw the molecular formulas including the lewis dot structures and covalent bonds

| Element Group               | 1A | 4A | 5A | 6A | 7A | 8A |
|-----------------------------|----|----|----|----|----|----|
| Number Bonds                | 1  | 4  | 3  | 2  | 1  | 0  |
| Number e <sup>-</sup> Pairs | 0  | 0  | 1  | 2  | 3  | 4  |

— = 2

oo = 2

Steps

1. Bond between atoms (—)

2. e<sup>-</sup> pairs

3. count e<sup>-</sup>

4. Add extra

H = 2 e<sup>-</sup>

4A - 8A = 8 e<sup>-</sup>

oo = 2

— = 2

Use chart

for # e<sup>-</sup> pairs

| O <sub>2</sub>                                                                                                     | HNBr <sub>2</sub>                                                                                                   | H <sub>2</sub> SiBr <sub>2</sub>                                                             |
|--------------------------------------------------------------------------------------------------------------------|---------------------------------------------------------------------------------------------------------------------|----------------------------------------------------------------------------------------------|
| <p>6 e<sup>-</sup> : 6 e<sup>-</sup> = double</p> <p>6 + 2 = 8      6 + 2 = 8</p> <p>O: 6A 2 pairs</p>             | <p>H: 1A N: 5A Br: 7A</p>                                                                                           | <p>H: 1A Si: 4A Br: 7A</p>                                                                   |
| HPSe                                                                                                               | SiPBr                                                                                                               | H <sub>2</sub> CO                                                                            |
| <p>6 e<sup>-</sup> : 6 e<sup>-</sup> = double</p> <p>2      6 + 2 = 8      6 + 2 = 8</p> <p>H: 1A P: 5A Se: 6A</p> | <p>4 e<sup>-</sup> : 4 e<sup>-</sup> = triple</p> <p>4 + 4 = 8      4 + 4 = 8      8</p> <p>Si: 4A P: 5A Br: 7A</p> | <p>H only needs 2 e<sup>-</sup></p> <p>6 + 2 = 8      6 + 2 = 8</p> <p>H: 1A C: 4A O: 6A</p> |

|        |      |    |     |       |       |      |       |      |      |      |
|--------|------|----|-----|-------|-------|------|-------|------|------|------|
| Number | 1    | 2  | 3   | 4     | 5     | 6    | 7     | 8    | 9    | 10   |
| Prefix | mono | di | tri | tetra | penta | hexa | hepta | octa | nona | deca |

Name or give the formula for the following molecules using the chart of prefixes above

|                                   |                        |                                  |                         |                        |
|-----------------------------------|------------------------|----------------------------------|-------------------------|------------------------|
| First atom<br>no mono             | Formula                | Name                             | Formula                 | Name                   |
|                                   | $\text{H}_3\text{N}_1$ | Trihydrogen Mononitride          | $\text{C}_2\text{Cl}_6$ | Dicarbon Hexachloride  |
| Second atom<br>-ide<br>Use chart! | Formula                | Name                             | Formula                 | Name                   |
|                                   | $\text{SiO}_2$         | no prefix = 1<br>Silicon dioxide | $\text{C}_2\text{F}_4$  | Dicarbon Tetrafluoride |

Calculate the Electronegativity for the following molecules, and polarity

|                  |                                                          |                      |
|------------------|----------------------------------------------------------|----------------------|
| Ionic – ED > 1.9 | Polar – ED <del>0.5</del> – 1.9<br><del>0.51</del> – 1.9 | Non-Polar – ED < 0.5 |
|------------------|----------------------------------------------------------|----------------------|

|                                                     |                              |          |                              |          |                              |           |
|-----------------------------------------------------|------------------------------|----------|------------------------------|----------|------------------------------|-----------|
| Use elektro.<br>chart<br>(given)                    | Ba - F                       |          | H - N                        |          | F - Cl                       |           |
|                                                     | Elect. Ba                    | Elect. F | Elect. H                     | Elect. N | Elect. F                     | Elect. Cl |
|                                                     | 0.9                          | 4.0      | 2.1                          | 3.0      | 4.0                          | 3.0       |
| ED always +<br>If -, make +                         | Electronegativity Difference |          | Electronegativity Difference |          | Electronegativity Difference |           |
|                                                     | $4.0 - 0.9 = 3.1$            |          | $3.0 - 2.1 = 0.9$            |          | $4.0 - 3.0 = 1.0$            |           |
| large-small<br>ED calc<br>Metal: Left<br>of zig-zag | Bond Type                    |          | Bond Type                    |          | Bond Type                    |           |
|                                                     | Ionic (metal)                |          | Polar Covalent (no metal)    |          | Polar Covalent (no metal)    |           |

Calculate the electronegativity, then use the electronegativity and molecule structure to determine molecule polarity.

|                                                                                                          |                              |                           |                                                                                                                                                                               |           |
|----------------------------------------------------------------------------------------------------------|------------------------------|---------------------------|-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|-----------|
| <u>Non Polar</u><br>Symm.<br><u>Polar</u><br>Polar Bonds<br>Not symm.<br><u>Don't forget</u><br>— and .. | H <sub>2</sub> O             |                           | Atomic Structure                                                                                                                                                              |           |
|                                                                                                          | Elect. H                     | Elect. O                  | <div>Symm = Non-Polar</div> <div><div>H</div><div>—</div><div><div>..</div><div>..</div>O</div><div>—</div><div>H</div></div> <div><div>2</div><div>8</div><div>2</div></div> |           |
|                                                                                                          | 2.1                          | 3.5                       |                                                                                                                                                                               |           |
|                                                                                                          | Electronegativity Difference |                           |                                                                                                                                                                               |           |
|                                                                                                          | 3.5 - 2.1<br>= 1.4           |                           | Bond Type                                                                                                                                                                     |           |
|                                                                                                          |                              |                           | Molecule Type                                                                                                                                                                 |           |
|                                                                                                          |                              | Polar Covalent (no metal) |                                                                                                                                                                               | Non-Polar |