| iA | $1 / \mathrm{A}$ | The Periodic Table |  |  |  |  |  |  |  |  |  |  |  |  |  |  | V/IMA |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $\begin{gathered} \mathrm{H} \\ 1 \end{gathered}$ |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  | He 2 |
| 1.01 |  |  |  |  |  |  |  |  |  |  |  | IIIA | IVA | VA | V/A | V/IIA | 4.00 |
| $\begin{gathered} \hline \mathrm{Li} \\ 3 \end{gathered}$ | Be 4 |  |  |  |  |  |  |  |  |  |  | B 5 | C | N 7 | 0 <br> 8 | F | Ne 10 |
| 6.94 | 9.01 |  |  |  |  |  |  |  |  |  |  | 10.81 | 12.01 | 14.01 | 16.00 | 19.00 | 20.18 |
| $\begin{gathered} \mathrm{Na} \\ 11 \end{gathered}$ | $\begin{gathered} \mathrm{Mg} \\ 12 \end{gathered}$ |  |  |  |  |  |  |  |  |  |  | AI 13 | Si 14 | P | S | $\mathrm{Cl}_{17}$ |  |
| 22.99 | 24.31 | Im | IVB | VB | V/B | V/IB | V/me | V/IMB | V/IIB | 18 | IIB | 26.98 | 28.09 | 30.97 | 32.07 | 35.45 | 39.95 |
| K | Ca | Sc | Ti | V | Cr | Mn | Fe | Co | Ni | Cu | Zn | Ga | Ge | As | Se | Br | Kr |
| 19 | 20 | 21 | 22 | 23 | 24 | 25 | 26 | 27 | 28 | 29 | 30 | 31 | 32 | 33 | 34 | 35 | 36 |
| 39.10 | 40.08 | 44.96 | 47.88 | 50.94 | 52.00 | 54.94 | 55.85 | 58.93 | 58.69 | 63.55 | 65.39 | 69.72 | 72.61 | 74.92 | 78.96 | 79.90 | 83.80 |
| Rb | Sr | Y | Zr | Nb | Mo | Tc | Ru | Rh | Pd | Ag | Cd | In | Sn | Sb | Te | I | Xe |
| 37 | 38 | 39 | 40 | 41 | 42 | 43 | 44 | 45 | 46 | 47 | 48 | 49 | 50 | 51 | 52 | 53 | 54 |
| 85.47 | 87.62 | 88.91 | 91.22 | 92.91 | 95.94 | (97.9) | 101.07 | 102.91 | 106.42 | 107.87 | 112.41 | 114.82 | 118.71 | 121.76 | 127.60 | 126.90 | 131.29 |
| Cs | Ba | La | Hf | Ta | W | Re | Os | Ir | Pt | Au | Hg | TI | Pb | Bi | Po | At | Rn |
| 55 | 56 | 57 | 72 | 73 | 74 | 75 | 76 | 77 | 78 | 79 | 80 | 81 | 82 | 83 | 84 | 85 | 86 |
| 132.91 | 137.33 | 138.91 | 178.49 | 180.95 | 183.85 | 186.21 | 190.2 | 192.22 | 195.08 | 197.97 | 200.59 | 204.38 | 207.2 | 208.98 | (209) | (210) | (222) |
| Fr | Ra | Ac | Rf | Db | Sg | Bh | Hs | Mt | Ds | Rg | Uub | Uut | Uuq | Uup |  |  |  |
| 87 | 88 | 89 | 104 | 105 | 106 | 107 | 108 | 109 | 110 | 111 | 112 | 113 | 114 | 115 |  |  |  |
| 223.02 | 226.03 | 227.03 | (261) | (262) | 263) | (262) | (265) | (266) | (271) | (272) | (285) | (284) | (289) | (288) |  |  |  |


| Ce <br> 58 <br> 140.12 | $\begin{array}{\|c\|} \hline \mathrm{Pr} \\ 59 \\ \mathbf{1 4 0 . 9 1} \\ \hline \end{array}$ | Nd 60 144.24 | $\begin{array}{\|c\|} \hline \text { Pm } \\ 61 \\ (145) \end{array}$ | $\begin{array}{\|c} \hline \text { Sm } \\ 62 \\ 150.36 \end{array}$ | Eu 63 152.97 | $\begin{array}{c\|} \hline \text { Gd } \\ 64 \\ 157.25 \end{array}$ | $\begin{array}{\|c\|} \hline \text { Tb } \\ 65 \\ 158.93 \\ \hline \end{array}$ | $\begin{gathered} \hline \text { Dy } \\ 66 \\ 162.50 \end{gathered}$ | $\mathrm{H} \circ$ 67 164.93 | $\begin{array}{\|c\|} \hline \mathrm{Er} \\ 68 \\ 167.26 \end{array}$ | $\begin{array}{\|c} \hline \mathrm{Tm} \\ 69 \\ 168.93 \end{array}$ | Yb <br> 70 <br> 173.04 | $\begin{array}{\|c\|} \hline \text { Lu } \\ 71 \\ 174.97 \end{array}$ |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| Th | $\begin{aligned} & \mathrm{Pa} \\ & 91 \end{aligned}$ | $\begin{aligned} & \mathrm{U} \\ & 92 \end{aligned}$ | $\begin{gathered} \mathrm{Np} \\ 93 \end{gathered}$ | $\begin{gathered} \mathrm{Pu} \\ 94 \end{gathered}$ | $\mathrm{Am}$ | $\begin{gathered} \mathrm{Cm} \\ 96 \end{gathered}$ | $\begin{gathered} \mathrm{Bk} \\ 97 \end{gathered}$ | $\begin{aligned} & \hline \text { Cf } \\ & 98 \end{aligned}$ | $\begin{gathered} \hline \text { Es } \\ 99 \end{gathered}$ | $\begin{aligned} & \hline \text { Fm } \\ & 100 \end{aligned}$ | $\begin{aligned} & \hline \text { Md } \\ & 101 \end{aligned}$ | $\begin{aligned} & \text { No } \\ & 102 \end{aligned}$ | Lr 103 |
| 232.04 | 231.04 | 238.03 | 237.05 | (240) | 243.06 | (247) | (248) | (251) | 252.08 | 257.10 | (257) | 259.10 | 262.11 |

Average Single Bond Lengths (Picometers)

|  | H | C | N | $\bigcirc$ | F | Si | P | S | CI | Br | 1 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| H | 74 | 110 | 98 | 94 | 92 | 145 | 138 | 132 | 127 | 142 | 161 |
| C |  | 154 | 147 | 143 | 141 | 194 | 187 | 181 | 176 | 191 | 210 |
| N |  |  | 140 | 136 | 134 | 187 | 180 | 174 | 169 | 184 | 203 |
| 0 |  |  |  | 132 | 130 | 183 | 176 | 170 | 165 | 180 | 199 |
| F |  |  |  |  | 128 | 181 | 174 | 168 | 163 | 178 | 197 |
| Si |  |  |  |  |  | 234 | 227 | 221 | 216 | 231 | 250 |
| P |  |  |  |  |  |  | 220 | 214 | 209 | 224 | 243 |
| S |  |  |  |  |  |  |  | 208 | 203 | 218 | 237 |
| Cl |  |  |  |  |  |  |  |  | 200 | 213 | 232 |
| Br |  |  |  |  |  |  |  |  |  | 228 | 247 |
| I |  |  |  |  |  |  |  |  |  |  | 266 |

Average Multiple Bond Lengths (Picometers)

| $\mathrm{C}=\mathrm{C}$ | 134 |
| :--- | :--- |
| $\mathrm{C}=\mathrm{N}$ | 127 |
| $\mathrm{C}=\mathrm{O}$ | 122 |
| $\mathrm{~N}=\mathrm{O}$ | 115 |$\quad$| $\mathrm{C} \equiv \mathrm{C}$ | 121 |
| :--- | :--- |
| $\mathrm{C} \equiv \mathrm{N}$ | 115 |
| $\mathrm{C} \equiv \mathrm{O}$ | 113 |
| $\mathrm{~N} \equiv \mathrm{O}$ | 108 |

$1 \mathrm{pm}=1 \times 10^{-12} \mathrm{~m}$

Average Single Bond Energies (kJ per mole)

|  | H | C | N | 0 | F | Si | P | S | Cl | Br | 1 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| H | 436 | 414 | 389 | 464 | 569 | 293 | 318 | 339 | 431 | 368 | 297 |
| C |  | 347 | 293 | 351 | 439 | 289 | 264 | 259 | 330 | 276 | 238 |
| N |  |  | 159 | 201 | 272 |  | 209 |  | 201 | 243 |  |
| 0 |  |  |  | 138 | 184 | 368 | 351 |  | 205 |  | 201 |
| F |  |  |  |  | 159 | 540 | 490 | 285 | 255 | 197 |  |
| Si |  |  |  |  |  | 176 | 213 | 226 | 360 | 289 |  |
| P |  |  |  |  |  |  | 213 | 230 | 331 | 272 | 213 |
| S |  |  |  |  |  |  |  | 213 | 251 | 213 |  |
| CI |  |  |  |  |  |  |  |  | 243 | 218 | 209 |
| Br |  |  |  |  |  |  |  |  |  | 192 | 180 |
| 1 |  |  |  |  |  |  |  |  |  |  | 151 |

Average Multiple Bond Energies (kJ per mole)

| $\mathrm{N}=\mathrm{N}$ | 418 |
| :--- | :--- |
| $\mathrm{~N} \equiv \mathrm{~N}$ | 946 |
| $\mathrm{~N}=\mathrm{O}$ | 590 |
| $\mathrm{C} \equiv \mathrm{N}$ | 891 |
| $\mathrm{O}=\mathrm{O}$ | 498 |$\quad$| $\mathrm{C}=\mathrm{C}$ | 611 |
| :--- | :--- | :--- |
| $\mathrm{C} \equiv \mathrm{C}$ | 837 |
| $\mathrm{C}=\mathrm{O}$ | 803 |
| $\mathrm{C}=\mathrm{O}$ | 745 |
| $\mathrm{C} \equiv \mathrm{O}$ | 1075 |

$\qquad$
$\qquad$

Question 1 Using noble gas notation, write the electron configuration for the following:

1. Co
2. Cu
3. $\mathrm{Fe}^{3+}$
4. $\mathrm{I}^{-}$ $\qquad$
5. Dy $\qquad$ (Dy = Element 66)

Question 2 Arrange the following elements in order of 5 Points increasing size, by ranking then from 1 (smallest) to 5 (largest)


Question 3 Arrange the following elements in order of 5 Points ionization energy, by ranking then from 1 (greatest) to 5 (smallest)


Question 4 Arrange the following elements in order of 5 Points metallic character, by ranking then from 1 (smallest) to 5 (greatest)

N
$\square$
Si


K
Al

Question 5 Draw the best Lewis Dot structure for the following 15 Points

| $\mathrm{N}_{2}$ | $\mathrm{ClO}_{3}{ }^{-}$ |
| :--- | :--- |
| $\mathrm{BeCl}_{2}$ | $\mathrm{XeF}_{4}$ |
| HCN |  |

Question 6 The following questions all relate to Ozone, $\mathrm{O}_{3}$

12 Points
6 Points

3 Points

3 Points

1. The molecule has two resonance structure. Draw them.
2. The bond $O-O-O$ bond angle is approximately:
3. The $O$ to $O$ bond energy in kJ per mole is:
(Circle the best choice)
a) $=498$
b) $>498$
c) $=138$
d) $>138$
e) $<138$

Question 7 The formal charge on the carbon and oxygen atoms in CO are:

6 Points

Question 8 Methane when combusted produces carbon dioxide and water according to:
6 Points
$C$ : $\qquad$ O: $\qquad$

Estimate the amount of energy produced upon the combustion of 1 mole of $\mathrm{CH}_{4}$ ?

Question 9 8 Points


What is the bond angle about the numbered atoms?

1. $\qquad$
2. $\qquad$
3. $\qquad$
4. $\qquad$

Question 10 28 Points

The following questions refer to the molecules depicted below.

| A |  | $\ddot{\mathrm{O}}=\ddot{\mathrm{s}}=\mathrm{O}$ | :C̈l—Be一C̈l: |
| :---: | :---: | :---: | :---: |
| E | : : | G | H $\text { : } \mathrm{F}-\dot{+} \mathrm{Xe}-\overrightarrow{\mathrm{F}}:$ |

1. List the structure(s) whose only bond angle is $\sim 180^{\circ}$
2. List the structures(s) whose epg is/are tetrahedral:
3. Give the electron pair geometry (epg) for:

A: $\qquad$ $C:$ $\qquad$
D: $\qquad$ F: $\qquad$
4. Give the molecular geometry for:

A: $\qquad$ E: $\qquad$
G: $\qquad$ H: $\qquad$
5. Two of the above molecules have an angular/bent molecular geometry. They are:
$\qquad$
$\qquad$ .. Which one has the largest bond angle? $\qquad$
6. Label the following molecules as either polar ( $P$ ) or non polar (NP)
A: $\qquad$
C: $\qquad$
D: $\qquad$
F: $\qquad$
H: $\qquad$

Do Not Write Below This Line

