

| $\begin{gathered} \hline \mathrm{Ce} \\ 58 \\ 140.12 \end{gathered}$ | $\begin{aligned} & \hline \mathrm{Pr} \\ & \hline 59 \end{aligned}$ | $\begin{array}{\|c\|} \hline \mathrm{Nd} \\ 60 \end{array}$ | $\begin{array}{\|c\|} \hline \text { Pm } \\ 61 \\ (145) \end{array}$ | $\underset{62}{\mathrm{Sm}}$ | $\begin{array}{\|c\|} \hline \mathrm{Eu} \\ 63 \\ \hline \end{array}$ | $\underset{64}{\mathrm{Gd}}$ | $\underset{65}{\mathrm{~Tb}}$ | $\begin{aligned} & \hline \text { Dy } \\ & 66 \end{aligned}$ | $\begin{aligned} & \mathrm{Ho} \\ & 67 \end{aligned}$ | $\begin{aligned} & \mathrm{Er} \\ & \hline 68 \end{aligned}$ | $\mathrm{Tm}_{69}$ | $\begin{aligned} & \mathrm{Yb} \\ & 70 \end{aligned}$ | $\mathrm{Lu}_{71}$ |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| Th | Pa | U | Np | Pu | Am | Cm | Bk | Cf | Es | m | Md | No | Lr |
| 90 | 91 | 92 | 93 | 94 | 95 | 96 |  | 98 | 99 | 100 | 101 | 102 | 103 |
| 2.04 | 11.0 | 88.0 | . 05 | (240) | 243.06 | (247) | (248) | (251) | 252.0 | 57.1 | (257) | 259 |  |



Some Approximate Single and Multiple Bond Lengths*

## Single Bond Lengths

|  | H | C | N | O | F | Si | P | S | Cl | Br | I |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 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 |
| O |  |  |  | 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 |
|  |  |  | Multiple Bond Lengths |  |  |  |  |  |  |  |  |
|  |  |  | $\mathrm{C}=\mathrm{C}$ |  |  | $\mathrm{C} \equiv \mathrm{C}$ | 121 |  |  |  |  |
|  |  |  | $\mathrm{C}=\mathrm{N}$ |  |  | $\mathrm{C} \equiv \mathrm{N}$ | 115 |  |  |  |  |
|  |  |  | $\mathrm{C}=\mathrm{O}$ |  |  | $\mathrm{C} \equiv \mathrm{O}$ | 113 |  |  |  |  |
|  |  |  | $\mathrm{N}=\mathrm{O}$ |  |  | $\mathrm{N} \equiv \mathrm{O}$ | 108 |  |  |  |  |

*In picometers (pm); $1 \mathrm{pm}=10^{-12} \mathrm{~m}$.


Last: First:


Circle the correct answers to the following questions, which relate to the orbital depicted on the left.

1. The orbital depicted is an $\mathbf{s}, \mathbf{p}, \mathbf{d}, \mathrm{f}$ or g orbital.
2. The principal quantum number for this orbital cannot be: 234
3. The likely specific designation for this orbital:
$2 s, 3 s, 2 p_{x}, 2 p_{y}, 2 p_{z}, 3 p_{x}, 3 p_{y}, 3 p_{z}, 2 d_{x y}, 2 d_{x z}$, $2 d_{y z}, 2 d_{z 2}, 2 d_{x 2-y 2}, 3 d_{x y}, 3 d_{x z}, 3 d_{y z}, 3 d_{z 2}, 3 d_{x 2-y 2}$

Question 2

Question 3 6 Points

Question 4 8 Points

Label the following atom/ions as either paramagnetic $(P)$ or diamagnetic ( $D$ ):

1. Be
2. 
3. 

$\qquad$

With respect to the elements, $\mathrm{Rb}, \mathrm{Cs}, \mathrm{K}$ and Na :
A. Which element would you expect to have the smallest atomic radius?
B. Which element would you expect to be most metallic?
C. Which element would you expect to have the largest ionization energy? $\qquad$
D. Which element would you expect to be least electronegative?


SID: $\square$ Last: First:

| Question 8 | Draw the three resonance structures for $\mathrm{CO}_{3}{ }^{2-}$. |
| :---: | :---: |
| $$ |  |
| Question 10 5 Points | The anticipated Carbon to Oxygen bond length in $\mathrm{CO}_{3}{ }^{2-}$ is: (check the correct answer) $\qquad$ 143pm $\qquad$ between 143pm and 122pm $\qquad$ 122 pm $\qquad$ between 122pm and 113pm $\qquad$ 113pm |
| Question 11 12 Points | Give the correct name for the following straight chain alkane, $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{3}$ $\qquad$ |
| $\begin{aligned} & \text { 흔 } \\ & \text { 흔 } \\ & \text { 옹 흔 } \\ & \hline \end{aligned}$ |   on the left represent the same molecule. Circle the structure that does not match the others |
| $\begin{aligned} & \text { 흔 } \\ & \text { 훌 } \\ & \text { 옹 흔 } \end{aligned}$ |  <br> Fill in the missing portions of the correct name given below for the molecule depicted on the left. <br> _-methyl $\qquad$ ane |



