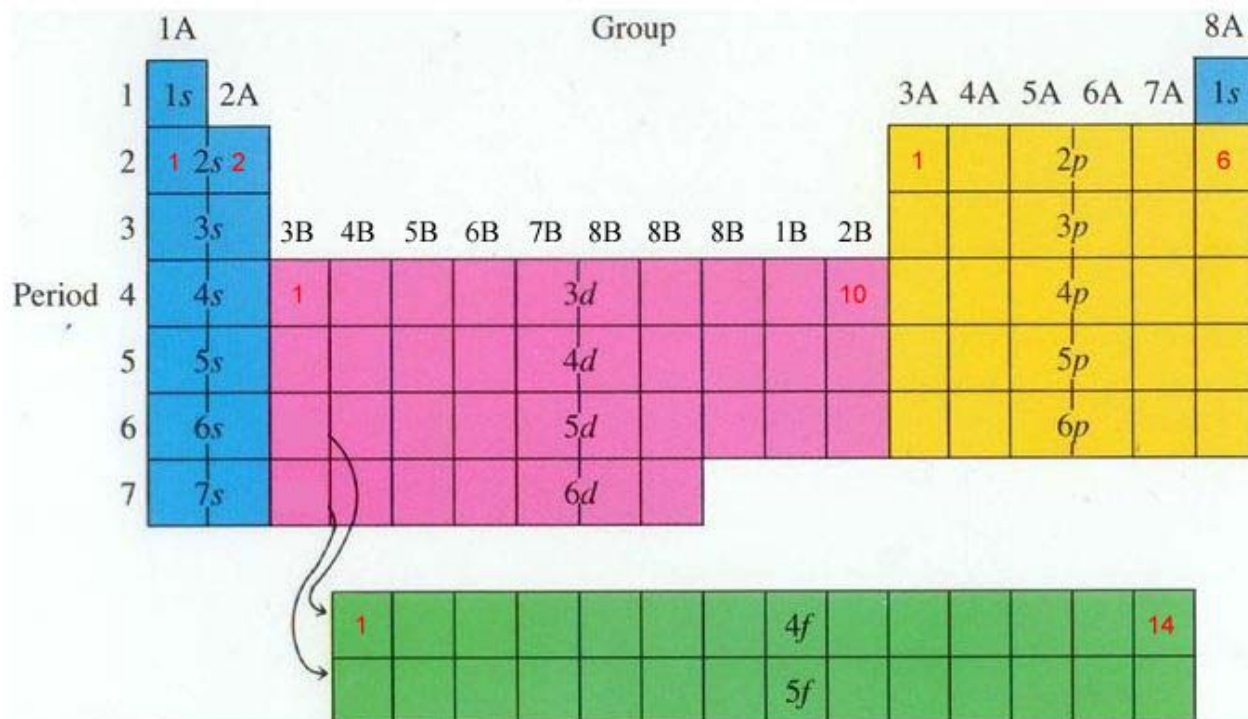


# The Periodic Table

|                           |                           |                           |                           |                           |                          |                           |                           |                           |                           |                           |                            |                            |                            |                            |                           |                          |                           |                          |
|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|--------------------------|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|----------------------------|----------------------------|----------------------------|----------------------------|---------------------------|--------------------------|---------------------------|--------------------------|
| <i>IA</i>                 |                           |                           |                           |                           |                          |                           |                           |                           |                           |                           |                            |                            |                            |                            |                           | <i>VIIIA</i>             |                           |                          |
| <b>H</b><br>1<br>1.01     |                           |                           |                           |                           |                          |                           |                           |                           |                           |                           | <b>He</b><br>2<br>4.00     |                            |                            |                            |                           |                          |                           |                          |
| <b>Li</b><br>3<br>6.94    |                           | <b>Be</b><br>4<br>9.01    |                           |                           |                          |                           |                           |                           |                           |                           |                            | <b>B</b><br>5<br>10.81     | <b>C</b><br>6<br>12.01     | <b>N</b><br>7<br>14.01     | <b>O</b><br>8<br>16.00    | <b>F</b><br>9<br>19.00   | <b>Ne</b><br>10<br>20.18  |                          |
| <b>Na</b><br>11<br>22.99  |                           | <b>Mg</b><br>12<br>24.31  |                           | <i>IIIB</i>               | <i>IVB</i>               | <i>VB</i>                 | <i>VIB</i>                | <i>VIIIB</i>              | <i>VIIIB</i>              | <i>VIIIB</i>              | <i>IB</i>                  | <i>IIB</i>                 | <b>Al</b><br>13<br>26.98   | <b>Si</b><br>14<br>28.09   | <b>P</b><br>15<br>30.97   | <b>S</b><br>16<br>32.07  | <b>Cl</b><br>17<br>35.45  | <b>Ar</b><br>18<br>39.95 |
| <b>K</b><br>19<br>39.10   | <b>Ca</b><br>20<br>40.08  | <b>Sc</b><br>21<br>44.96  | <b>Ti</b><br>22<br>47.88  | <b>V</b><br>23<br>50.94   | <b>Cr</b><br>24<br>52.00 | <b>Mn</b><br>25<br>54.94  | <b>Fe</b><br>26<br>55.85  | <b>Co</b><br>27<br>58.93  | <b>Ni</b><br>28<br>58.69  | <b>Cu</b><br>29<br>63.55  | <b>Zn</b><br>30<br>65.39   | <b>Ga</b><br>31<br>69.72   | <b>Ge</b><br>32<br>72.61   | <b>As</b><br>33<br>74.92   | <b>Se</b><br>34<br>78.96  | <b>Br</b><br>35<br>79.90 | <b>Kr</b><br>36<br>83.80  |                          |
| <b>Rb</b><br>37<br>85.47  | <b>Sr</b><br>38<br>87.62  | <b>Y</b><br>39<br>88.91   | <b>Zr</b><br>40<br>91.22  | <b>Nb</b><br>41<br>92.91  | <b>Mo</b><br>42<br>95.94 | <b>Tc</b><br>43<br>(97.9) | <b>Ru</b><br>44<br>101.07 | <b>Rh</b><br>45<br>102.91 | <b>Pd</b><br>46<br>106.42 | <b>Ag</b><br>47<br>107.87 | <b>Cd</b><br>48<br>112.41  | <b>In</b><br>49<br>114.82  | <b>Sn</b><br>50<br>118.71  | <b>Sb</b><br>51<br>121.76  | <b>Te</b><br>52<br>127.60 | <b>I</b><br>53<br>126.90 | <b>Xe</b><br>54<br>131.29 |                          |
| <b>Cs</b><br>55<br>132.91 | <b>Ba</b><br>56<br>137.33 | <b>La</b><br>57<br>138.91 | <b>Hf</b><br>72<br>178.49 | <b>Ta</b><br>73<br>180.95 | <b>W</b><br>74<br>183.85 | <b>Re</b><br>75<br>186.21 | <b>Os</b><br>76<br>190.2  | <b>Ir</b><br>77<br>192.22 | <b>Pt</b><br>78<br>195.08 | <b>Au</b><br>79<br>197.97 | <b>Hg</b><br>80<br>200.59  | <b>Tl</b><br>81<br>204.38  | <b>Pb</b><br>82<br>207.2   | <b>Bi</b><br>83<br>208.98  | <b>Po</b><br>84<br>(209)  | <b>At</b><br>85<br>(210) | <b>Rn</b><br>86<br>(222)  |                          |
| <b>Fr</b><br>87<br>223.02 | <b>Ra</b><br>88<br>226.03 | <b>Ac</b><br>89<br>227.03 | <b>Rf</b><br>104<br>(261) | <b>Db</b><br>105<br>(262) | <b>Sg</b><br>106<br>263  | <b>Bh</b><br>107<br>(262) | <b>Hs</b><br>108<br>(265) | <b>Mt</b><br>109<br>(266) | <b>Ds</b><br>110<br>(271) | <b>Rg</b><br>111<br>(272) | <b>Uub</b><br>112<br>(285) | <b>Uut</b><br>113<br>(284) | <b>Uuq</b><br>114<br>(289) | <b>Uup</b><br>115<br>(288) |                           |                          |                           |                          |

|                           |                           |                           |                           |                           |                           |                           |                           |                           |                           |                            |                           |                            |                            |
|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|----------------------------|---------------------------|----------------------------|----------------------------|
| <b>Ce</b><br>58<br>140.12 | <b>Pr</b><br>59<br>140.91 | <b>Nd</b><br>60<br>144.24 | <b>Pm</b><br>61<br>(145)  | <b>Sm</b><br>62<br>150.36 | <b>Eu</b><br>63<br>152.97 | <b>Gd</b><br>64<br>157.25 | <b>Tb</b><br>65<br>158.93 | <b>Dy</b><br>66<br>162.50 | <b>Ho</b><br>67<br>164.93 | <b>Er</b><br>68<br>167.26  | <b>Tm</b><br>69<br>168.93 | <b>Yb</b><br>70<br>173.04  | <b>Lu</b><br>71<br>174.97  |
| <b>Th</b><br>90<br>232.04 | <b>Pa</b><br>91<br>231.04 | <b>U</b><br>92<br>238.03  | <b>Np</b><br>93<br>237.05 | <b>Pu</b><br>94<br>(240)  | <b>Am</b><br>95<br>243.06 | <b>Cm</b><br>96<br>(247)  | <b>Bk</b><br>97<br>(248)  | <b>Cf</b><br>98<br>(251)  | <b>Es</b><br>99<br>252.08 | <b>Fm</b><br>100<br>257.10 | <b>Md</b><br>101<br>(257) | <b>No</b><br>102<br>259.10 | <b>Lr</b><br>103<br>262.11 |



### Average Single Bond Lengths (Picometers)

|    | 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 |

### Average Multiple Bond Lengths (Picometers)

|       |     |       |     |
|-------|-----|-------|-----|
| C = C | 134 | C ≡ C | 121 |
| C = N | 127 | C ≡ N | 115 |
| C = O | 122 | C ≡ O | 113 |
| N = O | 115 | N ≡ O | 108 |

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

=====

### Average Single Bond Energies (kJ per mole)

|    | H   | C   | N   | O   | F   | Si  | P   | S   | Cl  | Br  | I   |
|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| 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 |     |
| O  |     |     |     | 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 |     |
| Cl |     |     |     |     |     |     |     |     | 243 | 218 | 209 |
| Br |     |     |     |     |     |     |     |     |     | 192 | 180 |
| I  |     |     |     |     |     |     |     |     |     |     | 151 |






### Average Multiple Bond Energies (kJ per mole)

|       |     |       |      |                               |
|-------|-----|-------|------|-------------------------------|
| N = N | 418 | C = C | 611  |                               |
| N ≡ N | 946 | C ≡ C | 837  |                               |
| N = O | 590 | C = O | 803  | <i>In CO<sub>2</sub> Only</i> |
| C ≡ N | 891 | C = O | 745  |                               |
| O = O | 498 | C ≡ O | 1075 |                               |

SID

Last \_\_\_\_\_

First \_\_\_\_\_

|                                       |   |
|---------------------------------------|---|
| <p><b>Question 1</b><br/>3 Points</p> | <p>Each of the orbitals depicted has the <b>lowest</b> value of <b>n</b> possible for its type. Which one has the <b>highest n</b> value?</p> <div style="display: flex; justify-content: space-around; align-items: center;">    </div> <div style="display: flex; justify-content: space-around; margin-top: 10px;"> <span>a</span> <span>b</span> <span>c</span> </div> |
| <p><b>Question 2</b><br/>4 Points</p> | <div style="display: flex; align-items: center;">  <div style="flex: 1;"> <p>a) The orbital depicted on the left is what type of orbital? _____</p> <p>b) Based on its Radial Distribution depicted on the right you can label this orbital as? _____</p> </div>  </div>   |
| <p><b>Question 3</b><br/>4 Points</p> | <p>I am an <b>orbital</b> belong to a family whose <math>m_l</math> values are <b>-2, -1, 0, +1, +2</b> therefore I am a(n) _____ type orbital. I have a grand total of <b>5 nodes</b>, therefore my principal quantum number (<b>n</b>) is _____. Apart from me there are a <b>total of _____ other</b> orbital's that have this same principal quantum number. <b>Combined</b> we can accommodate a <b>total of _____ electrons</b>.</p>  |
| <p><b>Question 4</b><br/>6 Points</p> | <p>Give the <b>complete</b> electronic configuration for the following:</p> <p>a. <b>P</b> _____      b. <b>Al<sup>3+</sup></b> _____</p>   |
| <p><b>Question 5</b><br/>4 Points</p> | <p>How many <b>unpaired electrons</b> are in the <b>Fe</b> atom? _____</p> <p>Therefore <b>Fe</b> is <b>paramagnetic</b> or <b>diamagnetic</b>? _____</p>   |
| <p><b>Question 6</b><br/>6 Points</p> | <p>Give the <b>noble gas</b> configuration for the following</p> <p>a. <b>Kr</b> _____      c. <b>Cu</b> _____</p> <p>b. <b>Ni<sup>2+</sup></b> _____</p>   |
| <p><b>Question 7</b><br/>4 Points</p> | <p>Using only the periodic table <b>arrange</b> the following elements in order of <b>increasing atomic radius</b>:      <b>Na<sup>+</sup>, F<sup>-</sup>, O<sup>2-</sup>, Mg<sup>2+</sup></b></p> <p style="text-align: center;">             _____ <i>smallest</i>      _____      _____      _____ <i>largest</i> </p>   |

|                                  |  |                                |                        |               |                |
|----------------------------------|--|--------------------------------|------------------------|---------------|----------------|
| <p>Question 8<br/>6 Points</p>   | <p>a) Using only the periodic table <b>arrange</b> the following elements in order of <b>increasing atomic size</b>: S, Ca, F, Mg</p> <p>_____</p> <p><i>smallest</i> _____ _____ <i>largest</i></p> <p>b) Which one has the greatest Electron Affinity: _____</p> <p>c) Which one has the smallest first ionization energy: _____</p>   |                                |                        |               |                |
| <p>Question 9<br/>3 Points</p>   | <p>Using only the periodic table <b>arrange</b> the following elements in order of <b>decreasing ionization energy</b>: bromine, potassium, gallium</p> <p>_____</p> <p><i>largest</i> _____ <i>smallest</i></p>   |                                |                        |               |                |
| <p>Question 10<br/>12 Points</p> | <p>Draw the <b>best</b> Lewis Dot structure for the following</p> <table border="1" style="width: 100%; border-collapse: collapse;"> <tbody> <tr> <td style="width: 50%; padding: 10px; vertical-align: top;"> <math>\text{ClO}_2^-</math> (Cl=Chlorine)         </td> <td style="width: 50%; padding: 10px; vertical-align: top;"> <math>\text{HFCO}</math> </td> </tr> <tr> <td style="width: 50%; padding: 10px; vertical-align: top;"> <math>\text{BF}_3</math> </td> <td style="width: 50%; padding: 10px; vertical-align: top;"> <math>\text{XeF}_2</math> </td> </tr> </tbody> </table> | $\text{ClO}_2^-$ (Cl=Chlorine) | $\text{HFCO}$          | $\text{BF}_3$ | $\text{XeF}_2$ |
| $\text{ClO}_2^-$ (Cl=Chlorine)   | $\text{HFCO}$  |                                |                        |               |                |
| $\text{BF}_3$                    | $\text{XeF}_2$   |                                |                        |               |                |
| <p>Question 11<br/>4 Points</p>  | <p>Draw the <b>best</b> Lewis Dot structure for the following molecules on the rough work paper provided and then <b>classify each as either a free radical (yes) or not (no)</b></p> <p>a) <math>\text{NO}_2</math> _____                      c) <math>\text{BrO}_2</math> _____</p> <p>b) <math>\text{ClO}_2^-</math> _____ (Cl = chlorine)                      d) <math>\text{ClO}_2</math> _____</p>   |                                |                        |               |                |
| <p>Question 12<br/>6 Points</p>  | <p>Draw the <b>best</b> Lewis Dot structure for the following organic molecules</p> <table border="1" style="width: 100%; border-collapse: collapse;"> <tbody> <tr> <td style="width: 50%; padding: 10px; vertical-align: top;"> <math>\text{CH}_3\text{COCH}_3</math> </td> <td style="width: 50%; padding: 10px; vertical-align: top;"> <math>\text{C}_2\text{H}_2</math> </td> </tr> </tbody> </table>  | $\text{CH}_3\text{COCH}_3$     | $\text{C}_2\text{H}_2$ |               |                |
| $\text{CH}_3\text{COCH}_3$       | $\text{C}_2\text{H}_2$   |                                |                        |               |                |

Question 13  
8 Points  
(6 Points)

Draw all reasonable resonance structure for  $\text{NO}_3^-$ .

(2 Points)

Circle the best answer:

*Average bond length table is on the front page of this exam.*

The N to O bond length in pm is expected to be:

1. = 136

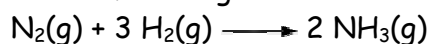
2. > 115

3. = 115

4. > 136

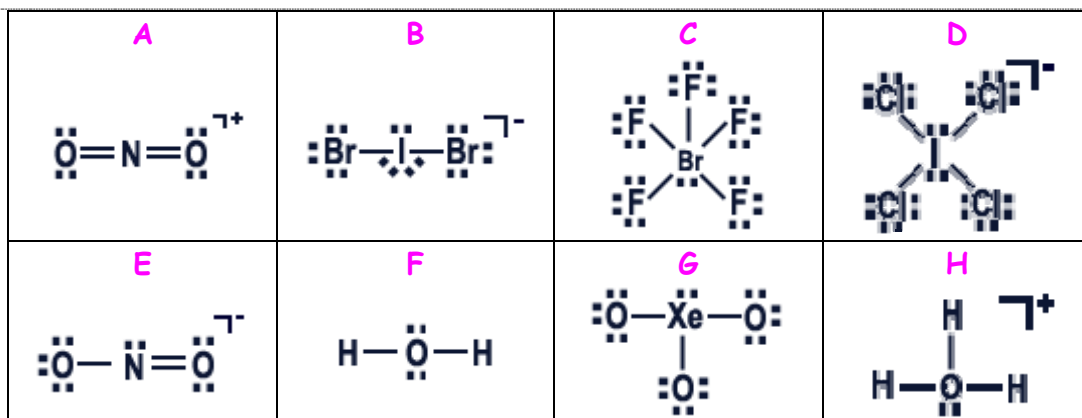
Question 14  
4 Points

Using the *Average bond energy table on the front page of this exam*, estimate the enthalpy change associated with the following reaction.



kJ

Question 15  
18 Points



1. List the **structure(s)** whose **only** bond angle is  $180^\circ$  \_\_\_\_\_

2. Give the **electron pair geometry (epg)** for:

**A:** \_\_\_\_\_

**C:** \_\_\_\_\_

**B:** \_\_\_\_\_

**E:** \_\_\_\_\_

3. Give the **molecular geometry** for:

**D:** \_\_\_\_\_

**F:** \_\_\_\_\_

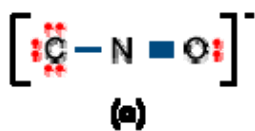
**G:** \_\_\_\_\_

**H:** \_\_\_\_\_

Question 16  
6 Points

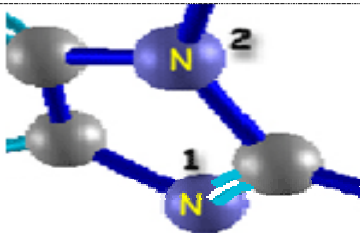
A resonance structure of  $\text{CNO}^-$  is given below:

Give the formal charge on:



C \_\_\_\_\_ N \_\_\_\_\_ O \_\_\_\_\_

Question 17  
6 Points



What is the **predicted bond angle** about the following atoms?

a) Nitrogen 1 \_\_\_\_\_

b) Nitrogen 2 \_\_\_\_\_

*Do Not Write Below This*

Exam II Score