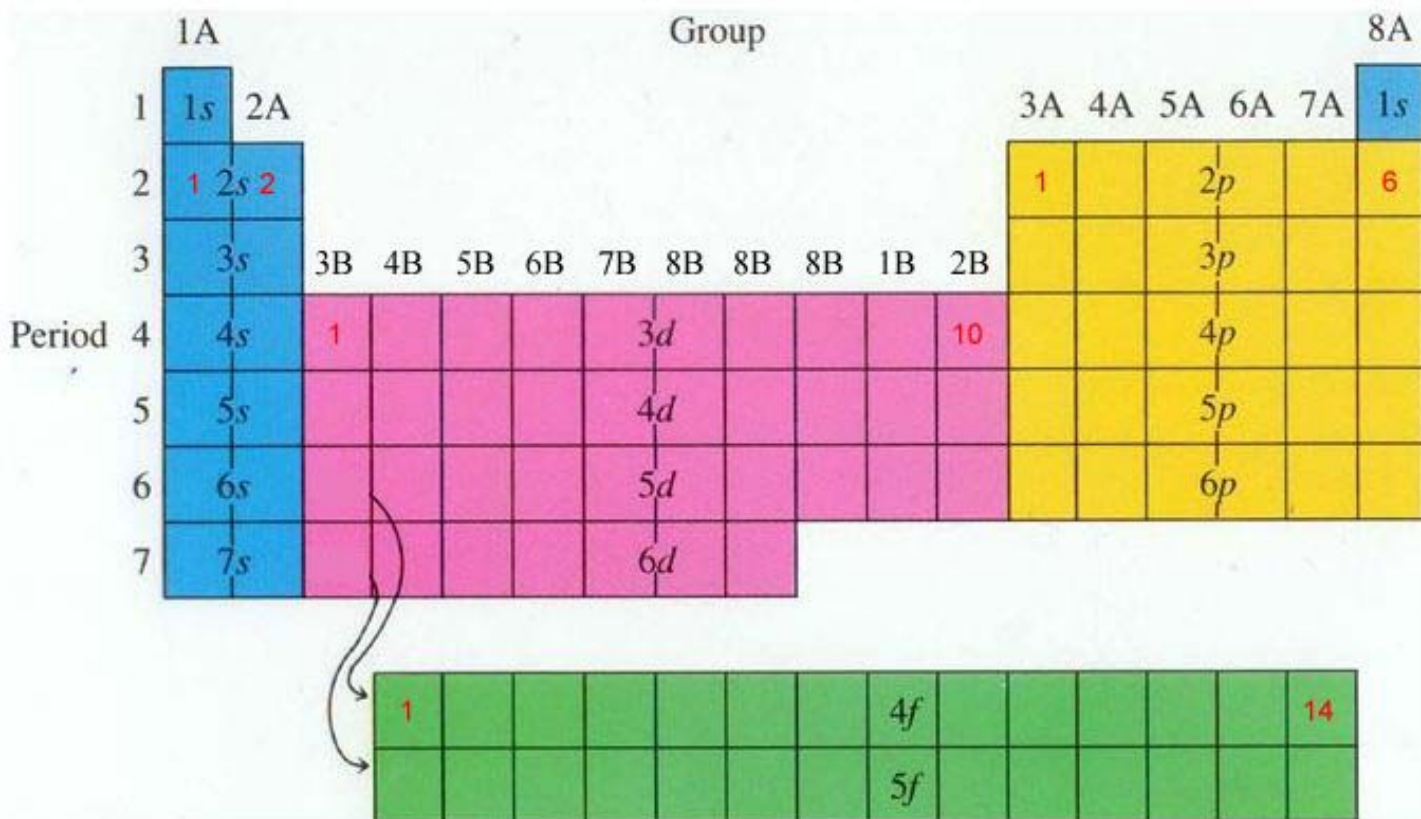


# The Periodic Table

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

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



### 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	

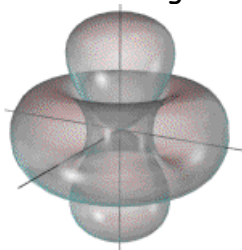
SID

Last \_\_\_\_\_

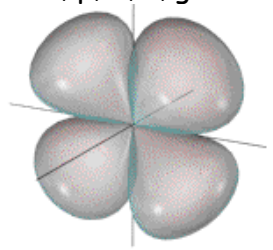
First \_\_\_\_\_

Question 1  
6 Points

Label the following orbital's as either: s, p, d, f, g?



\_\_\_\_\_



\_\_\_\_\_



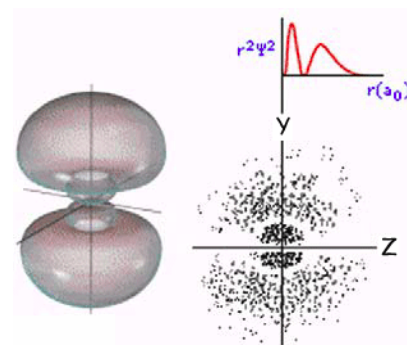
\_\_\_\_\_

Question 2  
6 Points

The orbital depicted on the left is:

- What **type** of orbital? \_\_\_\_\_
- Its **n** value is? \_\_\_\_\_
- Its **specific** designation is? \_\_\_\_\_

(x, y, z, xy, xz, yz, x<sup>2</sup>-y<sup>2</sup>, z<sup>2</sup>)



Question 3  
8 Points

Give the **complete** electronic configuration for the following:

- |                   |       |                     |       |
|-------------------|-------|---------------------|-------|
| a. S              | _____ | c. Mg <sup>2+</sup> | _____ |
| b. F <sup>-</sup> | _____ | d. Mn               | _____ |

Question 4  
8 Points

Give the **noble gas** configuration for the following

- |       |       |                     |       |
|-------|-------|---------------------|-------|
| a. Kr | _____ | c. Fe <sup>2+</sup> | _____ |
| b. Cu | _____ | d. Br <sup>-</sup>  | _____ |

Question 5  
2 Points

The element with the electronic configuration [Xe]6s<sup>2</sup>5d<sup>10</sup>4f<sup>14</sup>6p<sup>3</sup> has \_\_\_\_\_ valence electrons.

Question 6  
3 Points

Give the symbol for the **diamagnetic** element(s) in **period 4**? \_\_\_\_\_

Question 7  
5 Points

Using only the periodic table given with this exam rank the following elements **from 1 to 5** in order of **increasing ionization** (1 being the **smallest** ionization energy and 5 the **largest** ionization energy):

\_\_\_ Al      \_\_\_ Ca      \_\_\_ P      \_\_\_ Rb      \_\_\_ Ga

Question 8  
3 Points

Using only the periodic table given with this exam **arrange** the following elements in order of **increasing electron affinity**: **oxygen, fluorine, sulfur**

Smallest

Largest

Question 9  
12 Points

Draw the **best** Lewis Dot structure for the following molecules



Question 10  
8 Points

Draw the **best** Lewis Dot structure for the following organic molecules



Question 11  
9 Points  
(6 Points)

Draw all **reasonable** resonance structure for  $O_3$ .

Circle the best answer:

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

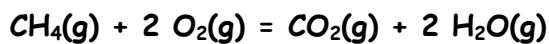
The O to O bond energy in  $\text{kJ}\cdot\text{mol}^{-1}$  is expected to be:

- = 498                       < 138                       = 138  
 > 498                       > 138

(3 Points)

Question 12  
6 Points

Using **average bond energies** (given on the front of this exam), **estimate** the **enthalpy change** for the following reaction:

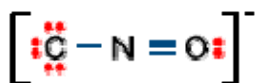


Show Work

$\text{kJ}\cdot\text{mol}^{-1}$

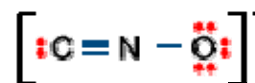
Question 13  
8 Points

Assign formal charges to the elements in each of the Lewis structures below.



A

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



B

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

The **best** Lewis structure for **CNO** - is: \_\_\_\_\_

Question 14  
6 Points

What is the **electron-pair geometry** for N in  $\text{NH}_3$ ? \_\_\_\_\_ ...

There are \_\_\_\_ **lone pair(s)** around the central atom, so the **molecular geometry** of  $\text{NH}_3$  is

\_\_\_\_\_.

Question 15  
6 Points

What is the **electron-pair geometry** for I in  $\text{IF}_3$  \_\_\_\_\_ ... There are

\_\_\_\_ **lone pair(s)** around the central atom, so the **molecular geometry** of  $\text{IF}_3$  is

\_\_\_\_\_.

Question 16  
4 Points

The planet **Whelanus** - far, far away in the galaxy - is somewhat like earth. Same order of elements, same order of filling orbitals, except that Hund's rule is different. Instead of 2 electrons per orbital there are **3**. This makes for a very different looking periodic table. On Whelanus give the symbols for the first 2 noble gases! \_\_\_\_\_

---

*Do Not Write Below This*

Exam II Score