

CHAPTER 2

2.8 Allowed values for the quantum numbers of electrons are as follows:

$$n = 1, 2, 3, \dots$$

$$l = 0, 1, 2, 3, \dots, n-1$$

$$m_l = 0, \pm 1, \pm 2, \pm 3, \dots, \pm l$$

$$m_s = \pm \frac{1}{2}$$

The relationships between n and the shell designations are noted in Table 2.1. Relative to the subshells,

$l = 0$ corresponds to an s subshell

$l = 1$ corresponds to a p subshell

$l = 2$ corresponds to a d subshell

$l = 3$ corresponds to an f subshell

For the K shell, the four quantum numbers for each of the two electrons in the $1s$ state, in the order of $nlm_l m_s$, are $100(\frac{1}{2})$ and $100(-\frac{1}{2})$. Write the four quantum numbers for all of the electrons in the L and M shells, and note which correspond to the s , p , and d subshells.

2.9 Give the electron configurations for the following ions: Fe^{2+} , Al^{3+} , Cu^+ , Ba^{2+} , Br^- , and O^{2-} .

2.18 The net potential energy between two adjacent ions, E_N , may be represented by the sum of Equations 2.9 and 2.11; that is,

$$E_N = -\frac{A}{r} + \frac{B}{r^n} \quad (2.17)$$

Calculate the bonding energy E_0 in terms of the parameters A , B , and n using the following procedure:

1. Differentiate E_N with respect to r , and then set the resulting expression equal to zero, since the curve of E_N versus r is a minimum at E_0 .
2. Solve for r in terms of A , B , and n , which yields r_0 , the equilibrium interionic spacing.
3. Determine the expression for E_0 by substitution of r_0 into Equation 2.17.

2.25 Compute the %IC of the interatomic bond for each of the following compounds: MgO , GaP , CsF , CdS , and FeO .