CHAPTER 12

12.13 Compute the atomic packing factor for the rock salt crystal structure in which $r_C/r_A = 0.414$.

12.16 Calculate the density of FeO, given that it has the rock salt crystal structure.

12.25 Compute the theoretical density of diamond, given that the C—C distance and bond angle are 0.154 nm and 109.5°, respectively. How does this value compare with the measured density?

12.27 Compute the atomic packing factor for the diamond cubic crystal structure (Figure 12.16). Assume that bonding atoms touch one another, that the angle between adjacent bonds is 109.5°, and that each atom internal to the unit cell is positioned a/4 of the distance away from the two nearest cell faces (a is the unit cell edge length).