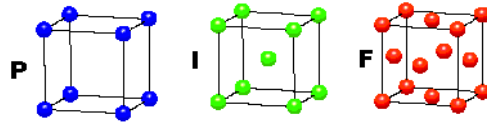


1.A.

CUBIC

$$a = b = c$$

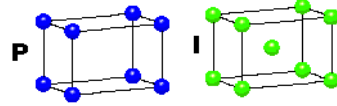
$$\alpha = \beta = \gamma = 90^\circ$$



TETRAGONAL

$$a = b \neq c$$

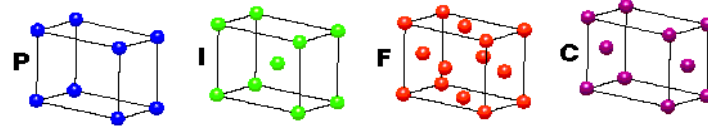
$$\alpha = \beta = \gamma = 90^\circ$$



ORTHORHOMBIC

$$a \neq b \neq c$$

$$\alpha = \beta = \gamma = 90^\circ$$

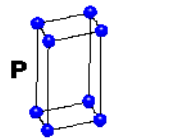


HEXAGONAL

$$a = b \neq c$$

$$\alpha = \beta = 90^\circ$$

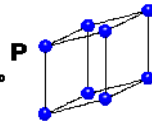
$$\gamma = 120^\circ$$



TRIGONAL

$$a = b = c$$

$$\alpha = \beta = \gamma \neq 90^\circ$$

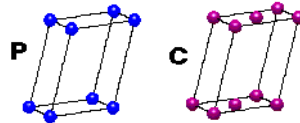


MONOCLINIC

$$a \neq b \neq c$$

$$\alpha = \gamma = 90^\circ$$

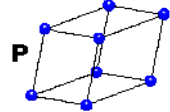
$$\beta \neq 120^\circ$$



TRICLINIC

$$a \neq b \neq c$$

$$\alpha \neq \beta \neq \gamma \neq 90^\circ$$



4 Types of Unit Cell

P = Primitive

I = Body-Centred

F = Face-Centred

C = Side-Centred

+

7 Crystal Classes

→ 14 Bravais Lattices

1.B.

$$p = \frac{h}{\lambda} = \frac{6.625 \times 10^{-34}}{85 \times 10^{-10}} = 7.79 \times 10^{-26}$$

$$E = \frac{p^2}{2m} = \frac{(7.79 \times 10^{-26})^2}{2 \times 9.11 \times 10^{-31}} = 3.33 \text{ J} = \underline{\underline{0.021 \text{ eV}}}$$

$$v = \frac{p}{m} = \frac{7.79 \times 10^{-26}}{9.11 \times 10^{-31}} = \underline{\underline{8.56 \times 10^4 \text{ m/s}}}$$

1.C.

$$\left(-\frac{\hbar^2}{2m} \frac{\partial^2}{\partial x^2} + V(x) \right) \psi(x,t) = j\hbar \frac{\partial}{\partial t} \psi(x,t)$$

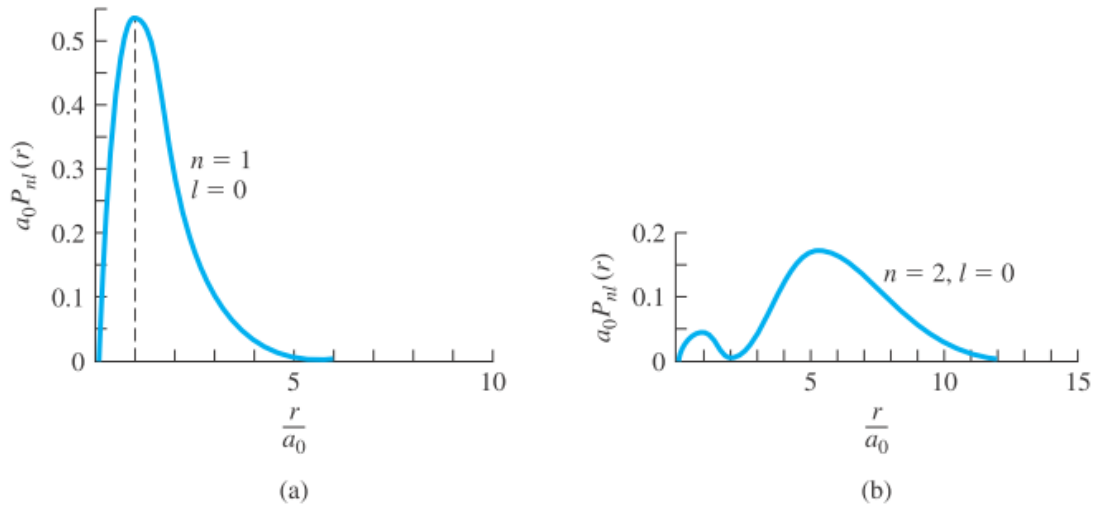


Figure 2.11 | The radial probability density function for the one-electron atom in the (a) lowest energy state and (b) next-higher energy state. (From Eisberg and Resnick [5].)

2.A.

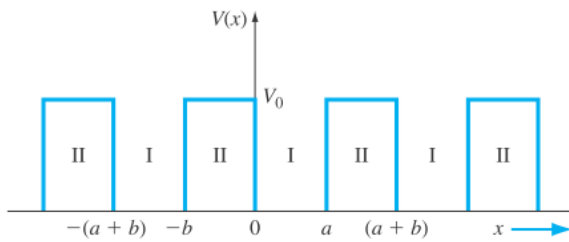


Figure 3.6 | The one-dimensional periodic potential function of the Kronig-Penney model.

1. In periodic potential, the wavefunction must be a Bloch function, where

$$\psi(x) = u(x)e^{jkx}$$

$$u(x) = u(x + a + b)$$

2. The solution of this wavefunction will satisfy

$$P \frac{\sin \alpha a}{\alpha a} + \cos \alpha a = \cos ka$$

$$P = \frac{mV_0 b a}{\hbar^2}$$

$$\alpha^2 = \frac{2mE}{\hbar^2}$$

Therefore, k is periodic in $\frac{2\pi}{a}$. The discontinuous allowed bands and forbidden bands by solving the equation numerical, as shown in the following figures.

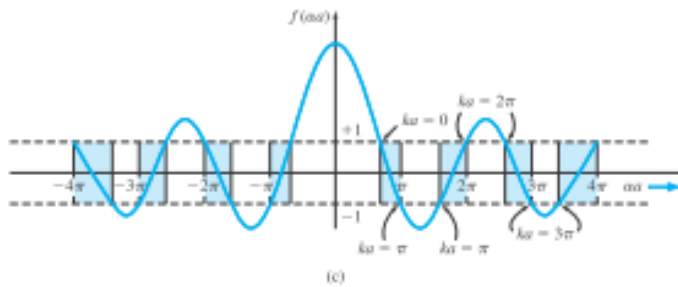


Figure 3.8 | A plot of (a) the first term in Equation (3.29), (b) the second term in Equation (3.29), and (c) the entire $f(ka)$ function. The shaded areas show the allowed values of (ka) corresponding to real values of k .

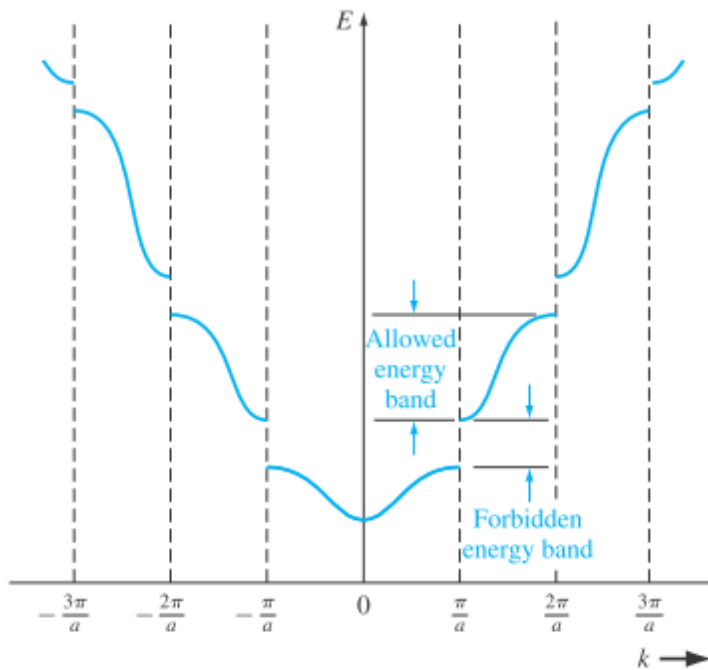


Figure 3.9 | The E versus k diagram generated from Figure 3.8. The allowed energy bands and forbidden energy bandgaps are indicated.

2.B.

$$f(E) = \frac{1}{1 + \exp\left(\frac{E - E_F}{kT}\right)} = \frac{1}{1 + \exp(5)} = 0.269$$

3.A.

$$E_{F_i} - E_{midgap} = \frac{3}{4}kT \ln\left(\frac{m_p^*}{m_n^*}\right) = \frac{3}{4} \times 0.0259 \times \ln(10) = 0.0447 \text{ eV}$$

3.B.

(i) This is p-type. Acceptors are added.

(ii)

$$E_F - E_{midgap} = -0.45$$

$$E_F - E_{F_i} = -0.45 - 0.0447 = -0.4947$$

$$p_0 = n_i \exp\left[-\frac{E_F - E_{F_i}}{kT}\right] = 10^5 \exp\left[\frac{0.4947}{0.0259}\right] = 1.97 \times 10^{13} \text{ cm}^{-3}$$