

# Problem Set 2

Due: Sep 27, 2024

## 1 Problem 1: Ashcroft & Mermin, Chapter 5, Problem 2

### 1.1 (a)

Using the primitive vectors given in Eq (4.9) and the construction (5.3) (or by any other method) show that the reciprocal of the simple hexagonal Bravais lattice is also simple hexagonal, with lattice constants  $2\pi/c$  and  $4\pi/\sqrt{3}a$ , rotated through  $30^\circ$  about the  $c$ -axis with respect to the direct lattice.

Relevant formula from the chapter: (4.9)

$$\begin{aligned}\mathbf{b}_1 &= 2\pi \frac{\mathbf{a}_2 \times \mathbf{a}_3}{\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)} \\ \mathbf{b}_2 &= 2\pi \frac{\mathbf{a}_3 \times \mathbf{a}_1}{\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)} \\ \mathbf{b}_3 &= 2\pi \frac{\mathbf{a}_1 \times \mathbf{a}_2}{\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)}\end{aligned}$$

and (5.3)

$$\mathbf{a}_1 = a\hat{x}, \quad \mathbf{a}_2 = \frac{a}{2}\hat{x} + \frac{\sqrt{3}a}{2}\hat{y}, \quad \mathbf{a}_3 = c\hat{z}.$$

#### 1.1.1 Solution

Our first task will be to evaluate the volume of the unit cell in real space, as given by  $\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)$ . First, we will evaluate the cross product of  $\mathbf{a}_2$  and  $\mathbf{a}_3$ :

$$\mathbf{a}_2 \times \mathbf{a}_3 = \left( \frac{a}{2}\hat{x} + \frac{\sqrt{3}a}{2}\hat{y} \right) \times c\hat{z} = -\left( \frac{a}{2}c \right)\hat{y} + \left( \frac{\sqrt{3}a}{2}c \right)\hat{x} \quad (1)$$

Next, we will evaluate the dot product of  $\mathbf{a}_1$  with this:

$$\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3) = a\hat{x} \cdot \left( -\left( \frac{a}{2}c \right)\hat{y} + \left( \frac{\sqrt{3}a}{2}c \right)\hat{x} \right) = \frac{\sqrt{3}a^2c}{2} \quad (2)$$

Now that we have the volume, we want to compute the numerators for the reciprocal lattice vectors:

$$\mathbf{a}_2 \times \mathbf{a}_3 = \left( \frac{a}{2} \hat{x} + \frac{\sqrt{3}a}{2} \hat{y} \right) \times c \hat{z} = - \left( \frac{a}{2} c \right) \hat{y} + \left( \frac{\sqrt{3}a}{2} c \right) \hat{x} \quad (3)$$

$$\mathbf{a}_3 \times \mathbf{a}_1 = c \hat{z} \times a \hat{x} = ca \hat{y} \quad (4)$$

$$\mathbf{a}_1 \times \mathbf{a}_2 = a \hat{x} \times \left( \frac{a}{2} \hat{x} + \frac{\sqrt{3}a}{2} \hat{y} \right) = \frac{\sqrt{3}a^2}{2} \hat{z} \quad (5)$$

Now we can compute the reciprocal lattice vectors by dividing these by the volume:

$$\mathbf{b}_1 = 2\pi \frac{\mathbf{a}_2 \times \mathbf{a}_3}{\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)} = 2\pi \frac{- \left( \frac{a}{2} c \right) \hat{y} + \left( \frac{\sqrt{3}a}{2} c \right) \hat{x}}{\frac{\sqrt{3}a^2 c}{2}} = \frac{2\pi}{a} \left( \hat{x} - \frac{1}{\sqrt{3}} \hat{y} \right) \quad (6)$$

$$\mathbf{b}_2 = 2\pi \frac{\mathbf{a}_3 \times \mathbf{a}_1}{\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)} = 2\pi \frac{ca \hat{y}}{\frac{\sqrt{3}a^2 c}{2}} = \frac{4\pi}{\sqrt{3}a} \hat{y} \quad (7)$$

$$\mathbf{b}_3 = 2\pi \frac{\mathbf{a}_1 \times \mathbf{a}_2}{\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)} = 2\pi \frac{\frac{\sqrt{3}a^2}{2} \hat{z}}{\frac{\sqrt{3}a^2 c}{2}} = \frac{2\pi}{c} \hat{z} \quad (8)$$

By taking the magnitude of the reciprocal lattice vectors, we can determine the new constants. For instance, the magnitude of  $\mathbf{b}_3$  is  $\frac{2\pi}{c}$ , while the magnitudes of  $\mathbf{b}_1$  and  $\mathbf{b}_2$  are  $\frac{4\pi}{\sqrt{3}a}$ . Now, we want to determine an angle of rotation. Since we know that the  $c$ -axis has the same definition in reciprocal space as defined now by  $\mathbf{b}_3$ , we now want to consider the angle between  $\mathbf{a}_1$  and  $\mathbf{b}_1$ , which is defined by the formula  $\cos(\theta) = \frac{\mathbf{a}_1 \cdot \mathbf{b}_1}{|\mathbf{a}_1| |\mathbf{b}_1|}$ . We can compute this as:

$$\mathbf{a}_1 \cdot \mathbf{b}_1 = a \hat{x} \cdot \left( \frac{2\pi}{a} \left( \hat{x} - \frac{1}{\sqrt{3}} \hat{y} \right) \right) = 2\pi \quad (9)$$

Now we already know the magnitudes, so we get:

$$\cos(\theta) = \frac{2\pi}{\left( a \times \frac{4\pi}{\sqrt{3}a} \right)} = \frac{\sqrt{3}}{2} \implies \theta = 30^\circ \quad (10)$$

So we show that the rotation between the direct and reciprocal lattices is  $30^\circ$ .

## 1.2 (b)

For what value of  $c/a$  does the ratio have the same value in both direct and reciprocal lattices? If  $c/a$  is ideal in the direct lattice, what is its value in the reciprocal lattice?

### 1.2.1 Solution

The ratio of interest is  $\frac{c}{a}$ . In the direct latest, this is simply  $\frac{c}{a}$ . In the reciprocal lattice, we can compute this as:

$$\frac{c}{a} = \frac{\frac{2\pi}{c}}{\frac{4\pi}{\sqrt{3}a}} = \frac{\sqrt{3}a}{2c} \quad (11)$$

Now, we want to find the value of  $c/a$  that makes these two ratios equal. We can set the two equal to each other and solve for  $c/a$ :

$$\frac{c}{a} = \frac{\sqrt{3}a}{2c} \implies c^2 = \sqrt{3}a^2 \implies \frac{c}{a} = \left(\frac{\sqrt{3}}{2}\right)^{1/2} \quad (12)$$

We know that in the direct lattice, the ideal ratio is  $\frac{c}{a} = \sqrt{\frac{8}{3}} \implies \frac{a}{c} = \sqrt{\frac{3}{8}}$ . Now we plug this in to the formula for the  $\frac{c}{a}$  ratio in the reciprocal lattice:

$$\frac{c}{a} = \frac{\sqrt{3}a}{2c} = \frac{\sqrt{3}}{2} \sqrt{\frac{3}{8}} = \frac{3}{4\sqrt{2}} \quad (13)$$

### 1.3 (c)

The Bravais lattice generated by three primitive vectors of equal length  $a$ , making equal angles  $\theta$  with one another, is known as the trigonal Bravais lattice (see Chapter 7). Show that the reciprocal of a trigonal Bravais lattice is also trigonal, with an angle  $\theta^*$  given by  $-\cos \theta^* = \cos \theta / [1 + \cos \theta]$ , and a primitive vector length  $a^*$ , given by  $a^* = (2\pi/a) (1 + 2 \cos \theta \cos \theta^*)^{-1/2}$ .

#### 1.3.1 Solution

Primitive vectors must have the same length and make equal angles with one another, which can be accomplished by permuting one nonzero quantity in the vector while keeping the others the same, i.e.,

$$\mathbf{a}_1 = \frac{a}{\sqrt{K^2 + 2}}(K, 1, 1) \quad (14)$$

$$\mathbf{a}_2 = \frac{a}{\sqrt{K^2 + 2}}(1, K, 1) \quad (15)$$

$$\mathbf{a}_3 = \frac{a}{\sqrt{K^2 + 2}}(1, 1, K) \quad (16)$$

where we have already enforced normalization. Now, we want to evaluate the reciprocal lattice vectors. First:

$$\mathbf{b}_1 = 2\pi \frac{\mathbf{a}_2 \times \mathbf{a}_3}{\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)} \quad (17)$$

Now, the first product is given by:

$$\mathbf{a}_2 \times \mathbf{a}_3 = \frac{a^2}{K^2 + 2} \begin{pmatrix} \hat{x} & \hat{y} & \hat{z} \\ 1 & K & 1 \\ 1 & 1 & K \end{pmatrix} = \frac{a^2}{K^2 + 2} (\hat{x}(K^2 - 1) - \hat{y}(K - 1) + \hat{z}(1 - K)) \quad (18)$$

Now, we can compute the dot product:

$$\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3) = \frac{a^3}{(K^2 + 2)^{\frac{3}{2}}} ((K - 1)(K + 1)K - (K - 1) - (K - 1)) = \frac{a^3(K + 2)(K - 1)^2}{(K^2 + 2)^{\frac{3}{2}}} \quad (19)$$

Now, we can compute the reciprocal lattice vector:

$$\mathbf{b}_1 = 2\pi \frac{\mathbf{a}_2 \times \mathbf{a}_3}{\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)} = 2\pi \frac{(K^2 + 2)^{\frac{1}{2}}}{a(K+2)(K-1)^2} (\hat{x}(K^2 - 1) - \hat{y}(K - 1) - \hat{z}(K - 1)) \quad (20)$$

$$= \frac{2\pi(K^2 + 2)^{1/2}}{a(K+2)(K-1)} (\hat{x}(K+1) - \hat{y} - \hat{z}) \quad (21)$$

Next, we can compute the other reciprocal lattice vectors. The relevant first product is

$$\mathbf{a}_3 \times \mathbf{a}_1 = \frac{a^2}{K^2 + 2} \begin{pmatrix} \hat{x} & \hat{y} & \hat{z} \\ 1 & 1 & K \\ K & 1 & 1 \end{pmatrix} = \frac{a^2}{K^2 + 2} (\hat{x}(1 - K) - \hat{y}(1 - K^2) + \hat{z}(1 - K)) \quad (22)$$

So we have:

$$\mathbf{b}_2 = 2\pi \frac{\mathbf{a}_3 \times \mathbf{a}_1}{\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)} = 2\pi \frac{(K^2 + 2)^{\frac{1}{2}}}{a(K+2)(K-1)^2} (\hat{x}(1 - K) - \hat{y}(1 - K^2) - \hat{z}(1 - K)) \quad (23)$$

$$= \frac{2\pi(K^2 + 2)^{1/2}}{a(K+2)(K-1)} (-\hat{x} + \hat{y}(K+1) - \hat{z}) \quad (24)$$

And then the final cross product is:

$$\mathbf{a}_1 \times \mathbf{a}_2 = \frac{a^2}{K^2 + 2} \begin{pmatrix} \hat{x} & \hat{y} & \hat{z} \\ K & 1 & 1 \\ 1 & K & 1 \end{pmatrix} = \frac{a^2}{K^2 + 2} (\hat{x}(1 - K) - \hat{y}(K - 1) + \hat{z}(K^2 - 1)) \quad (25)$$

So we have:

$$\mathbf{b}_3 = 2\pi \frac{\mathbf{a}_1 \times \mathbf{a}_2}{\mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)} = 2\pi \frac{(K^2 + 2)^{\frac{1}{2}}}{a(K+2)(K-1)^2} (\hat{x}(1 - K) - \hat{y}(K - 1) + \hat{z}(K^2 - 1)) \quad (26)$$

$$= \frac{2\pi(K^2 + 2)^{1/2}}{a(K+2)(K-1)^2} (\hat{x}(1 - K) - \hat{y}(K - 1) + \hat{z}(K^2 - 1)) \quad (27)$$

$$= \frac{2\pi(K^2 + 2)^{1/2}}{a(K+2)(K-1)} (-\hat{x} - \hat{y} + \hat{z}(K+1)) \quad (28)$$

So we see that the reciprocal lattice is also trigonal;  $(K+1)$  is permuted among entries.

## 2 Problem 2: Ashcroft & Mermin, Chapter 8, Problem 1

### Problem Description:

The band structure of the one-dimensional solid can be expressed quite simply in terms of the properties of an electron in the presence of a single-barrier potential  $t(x)$ . Consider therefore

an electron incident from the left on the potential barrier  $v(x)$  with energy  $\varepsilon = \hbar^2 K^2 / 2m$ . Since  $v(x) = 0$  when  $|x| \geq a/2$ , in these regions the wave function  $\psi_1(x)$  will have the form

$$\begin{aligned}\psi_1(x) &= e^{iKx} + re^{-iKx}, \quad x \leq -\frac{a}{2} \\ &= te^{iKx}, \quad x \geq \frac{a}{2}\end{aligned}$$

This is illustrated schematically in Figure 8.5a. The transmission and reflection coefficients  $t$  and  $r$  give the probability amplitude that the electron will tunnel through or be reflected from the barrier; they depend on the incident wave vector  $K$  in a manner determined by the detailed features of the barrier potential  $v$ . However, one can deduce many properties of the band structure of the periodic potential  $U$  by appealing only to very general properties of  $t$  and  $r$ . Because  $v$  is even,  $\psi_r(x) = \psi_r(-x)$  is also a solution to the Schrödinger equation with energy  $\varepsilon$ . From (8.65) it follows that  $\psi_r(x)$  has the form

$$\begin{aligned}\psi_r(x) &= te^{-iKx}, \quad x \leq -\frac{a}{2} \\ &= e^{-iKx} + re^{iKx}, \quad x \geq \frac{a}{2}\end{aligned}$$

Evidently this describes a particle incident on the barrier from the right, as depicted in Figure 8.5b. Since  $\psi_l$  and  $\psi_r$  are two independent solutions to the single-barrier Schrödinger equation with the same energy, any other solution with that energy will be a linear combination<sup>34</sup> of these two:  $\psi = A\psi_l + B\psi_r$ . In particular, since the crystal Hamiltonian is identical to that for a single ion in the region  $-a/2 \leq x \leq a/2$ , any solution to the crystal Schrödinger equation with energy  $\varepsilon$  must be a linear combination of  $\psi_l$  and  $\psi_r$  in that region:

$$\psi(x) = A\psi_l(x) + B\psi_r(x), \quad -\frac{a}{2} \leq x \leq \frac{a}{2}$$

Now Bloch's theorem asserts that  $\psi$  can be chosen to satisfy

$$\psi(x+a) = e^{ika}\psi(x)$$

for suitable  $k$ . Differentiating (8.68) we also find that  $\psi' = d\psi/dx$  satisfies

$$\psi'(x+a) = e^{ika}\psi'(x)$$

## 2.1

(a) By imposing the conditions (8.68) and (8.69) at  $x = -a/2$ , and using (8.65) to (8.67), show that the energy of the Bloch electron is related to its wave vector  $k$  by:

$$\cos ka = \frac{t^2 - r^2}{2t} e^{iKa} + \frac{1}{2t} e^{-iKa}, \quad \mathcal{E} = \frac{\hbar^2 K^2}{2m}$$

Verify that this gives the right answer in the free electron case ( $v \equiv 0$ ).

### 2.1.1 Solution

We begin by imposing the boundary conditions, which gives:

$$\psi\left(\frac{a}{2}\right) = e^{ika}\psi\left(-\frac{a}{2}\right) \quad (29)$$

$$\psi'\left(\frac{a}{2}\right) = e^{ika}\psi'\left(-\frac{a}{2}\right) \quad (30)$$

Now, we plug in for both  $\psi_l$  and  $\psi_r$ :

$$\psi_l\left(-\frac{a}{2}\right) = re^{iK\frac{a}{2}} + e^{-iK\frac{a}{2}} \quad (31)$$

$$\psi_r\left(-\frac{a}{2}\right) = te^{iK\frac{a}{2}} \quad (32)$$

and then for the derivatives

$$\psi'_l\left(-\frac{a}{2}\right) = iK\left(-re^{iK\frac{a}{2}} + e^{-iK\frac{a}{2}}\right) \quad (33)$$

$$\psi'_r\left(-\frac{a}{2}\right) = -iKte^{iK\frac{a}{2}}. \quad (34)$$

At the other boundary we have:

$$\psi_l\left(\frac{a}{2}\right) = te^{iK\frac{a}{2}} \quad (35)$$

$$\psi_r\left(\frac{a}{2}\right) = re^{iK\frac{a}{2}} + e^{-iK\frac{a}{2}} \quad (36)$$

and then for the derivatives

$$\psi'_l\left(\frac{a}{2}\right) = iKte^{iK\frac{a}{2}} \quad (37)$$

$$\psi'_r\left(\frac{a}{2}\right) = iK\left(re^{iK\frac{a}{2}} - e^{-iK\frac{a}{2}}\right). \quad (38)$$

Now, we plug in these results into the Bloch conditions:

$$A\psi_l\left(\frac{a}{2}\right) + B\psi_r\left(\frac{a}{2}\right) = e^{ika}\left(A\psi_l\left(-\frac{a}{2}\right) + B\psi_r\left(-\frac{a}{2}\right)\right). \quad (39)$$

Plugging and and redefining the reciprocal vector  $K \equiv G$  to avoid confusion, we have:

$$A\left(te^{iG\frac{a}{2}}\right) + B\left(re^{iG\frac{a}{2}} + e^{-iG\frac{a}{2}}\right) = e^{ika}\left(A\left(re^{iG\frac{a}{2}} + e^{-iG\frac{a}{2}}\right) + B\left(te^{iG\frac{a}{2}}\right)\right) \quad (40)$$

Grouping terms onto the left hand side and combining the exponentials, we have:

$$A\left(te^{iG\frac{a}{2}} - re^{ia\left(k+\frac{G}{2}\right)} - e^{ia\left(k-\frac{G}{2}\right)}\right) + B\left(re^{iG\frac{a}{2}} + e^{-iG\frac{a}{2}} - te^{ia\left(k+\frac{G}{2}\right)}\right) = 0 \quad (41)$$

For the derivative we have the condition:

$$A\psi'_l\left(\frac{a}{2}\right) + B\psi'_r\left(\frac{a}{2}\right) = e^{ika}\left(A\psi'_l\left(-\frac{a}{2}\right) + B\psi'_r\left(-\frac{a}{2}\right)\right). \quad (42)$$

Plugging in and redefining the reciprocal vector  $K \equiv G$  to avoid confusion, we have:

$$A \left( iG t e^{iG \frac{a}{2}} \right) + B \left( iG \left( r e^{iG \frac{a}{2}} - e^{-iG \frac{a}{2}} \right) \right) = e^{ika} \left( A \left( iG \left( -r e^{iG \frac{a}{2}} + e^{-iG \frac{a}{2}} \right) \right) + B \left( -iG t e^{iG \frac{a}{2}} \right) \right) \quad (43)$$

Grouping terms onto the left hand side and combining the exponentials and dividing by  $iG$ , we have:

$$A \left( t e^{iG \frac{a}{2}} + r e^{ia(k+\frac{G}{2})} - e^{ia(k-\frac{G}{2})} \right) + B \left( r e^{iG \frac{a}{2}} - e^{-iG \frac{a}{2}} + t e^{ia(k+\frac{G}{2})} \right) = 0 \quad (44)$$

Next, we can define the matrix equation:

$$\begin{pmatrix} T_{11} & T_{12} \\ T_{21} & T_{22} \end{pmatrix} \begin{pmatrix} A \\ B \end{pmatrix} = 0 \quad (45)$$

where  $T_{11} = t e^{iG \frac{a}{2}} - r e^{ia(k+\frac{G}{2})} - e^{ia(k-\frac{G}{2})}$ ,  $T_{12} = r e^{iG \frac{a}{2}} + e^{-iG \frac{a}{2}} - t e^{ia(k+\frac{G}{2})}$ ,  $T_{21} = t e^{iG \frac{a}{2}} + r e^{ia(k+\frac{G}{2})} - e^{ia(k-\frac{G}{2})}$ , and  $T_{22} = r e^{iG \frac{a}{2}} - e^{-iG \frac{a}{2}} + t e^{ia(k+\frac{G}{2})}$ . The determinant of this matrix must be zero, so we have:

$$T_{11}T_{22} - T_{12}T_{21} = 0 \quad (46)$$

Plugging in the expressions for  $T_{11}$ ,  $T_{12}$ ,  $T_{21}$ , and  $T_{22}$ , we have:

$$\left( t e^{iG \frac{a}{2}} - r e^{ia(k+\frac{G}{2})} - e^{ia(k-\frac{G}{2})} \right) \left( r e^{iG \frac{a}{2}} - e^{-iG \frac{a}{2}} + t e^{ia(k+\frac{G}{2})} \right) \quad (47)$$

$$- \left( r e^{iG \frac{a}{2}} + e^{-iG \frac{a}{2}} - t e^{ia(k+\frac{G}{2})} \right) \left( t e^{iG \frac{a}{2}} + r e^{ia(k+\frac{G}{2})} - e^{ia(k-\frac{G}{2})} \right) = 0 \quad (48)$$

Sympy simplifies this to:

$$2 \left( -r^2 e^{2iGa} + t^2 e^{2iGa} - t e^{ia(G-k)} - t e^{ia(G+k)} + 1 \right) e^{ia(-G+k)} = 0 \quad (49)$$

Dividing both sides by the factor of  $2e^{ia(-G+k)}$ , we have:

$$-r^2 e^{2iGa} + t^2 e^{2iGa} - t e^{ia(G-k)} - t e^{ia(G+k)} + 1 = 0 \quad (50)$$

Now, let us group the terms:

$$(t^2 - r^2) e^{2iGa} - t (e^{ia(G-k)} + e^{ia(G+k)}) + 1 = 0 \quad (51)$$

Now, notice the identity of:

$$e^{ia(G-k)} + e^{ia(G+k)} = 2e^{iaG} \cos(ka) \quad (52)$$

Substitution with further manipulation gives the desired result:

$$\cos(ka) = \frac{t^2 - r^2}{2t} e^{iGa} + \frac{1}{2t} e^{-iGa} \quad (53)$$

Now, we can verify this in the free electron case by setting  $v \equiv 0$ . In this case, we have  $t = 1$  and  $r = 0$ , so the equation becomes:

$$\cos(ka) = \frac{1}{2} e^{iGa} + \frac{1}{2} e^{-iGa} = \cos(Ga) \quad (54)$$

## 2.2

Equation (8.70) is more informative when one supplies a little more information about the transmission and reflection coefficients. We write the complex number  $t$  in terms of its magnitude and phase:

$$t = |t|e^{i\delta}$$

The real number  $\delta$  is known as the phase shift, since it specifies the change in phase of the transmitted wave relative to the incident one. Electron conservation requires that the probability of transmission plus the probability of reflection be unity:

$$1 = |t|^2 + |r|^2$$

This, and some other useful information, can be proved as follows. Let  $\phi_1$  and  $\phi_2$  be any two solutions to the one-barrier Schrodinger equation with the same energy:

$$-\frac{\hbar^2}{2m}\phi_i'' + r\phi_i = \frac{\hbar^2 K^2}{2m}\phi_i, \quad i = 1, 2$$

Define  $w(\phi_1, \phi_2)$  (the "Wronskian") by

$$w(\phi_1, \phi_2) = \phi_1'(x)\phi_2(x) - \phi_1(x)\phi_2'(x)$$

(b) Prove that  $w$  is independent of  $x$  by deducing from (8.73) that its derivative vanishes. We want to start by differentiating the Wronskian with respect to  $x$ :

### 2.2.1 Solution

$$\frac{d}{dx}w(\phi_1, \phi_2) = \frac{d}{dx}(\phi_1'\phi_2 - \phi_1\phi_2') = \phi_1''\phi_2 + \phi_1'\phi_2' - \phi_1'\phi_2' - \phi_1\phi_2'' = \phi_1''\phi_2 - \phi_1\phi_2'' \quad (55)$$

Now, we rearrange the Schrodinger equation for  $\phi_i''$

$$\phi_i'' = -\frac{2m}{\hbar^2} \left( \frac{\hbar^2 K^2}{2m} - r \right) \phi_i \quad (56)$$

Now we substitute in the forms for  $\phi_1''$  and  $\phi_2''$  into the derivative of the Wronskian:

$$\phi_1''\phi_2 - \phi_1\phi_2'' = -\frac{2m}{\hbar^2} \left( \frac{\hbar^2 K^2}{2m} - r \right) \phi_1\phi_2 + \frac{2m}{\hbar^2} \left( \frac{\hbar^2 K^2}{2m} - r \right) \phi_1\phi_2 = 0 \quad (57)$$

So we have shown that the derivative of the Wronskian vanishes, which implies that the Wronskian is independent of  $x$ .

## 2.3

(c) Prove (8.72) by evaluating  $w(\psi_l, \psi_l^*)$  for  $x \leq -a/2$  and  $x \geq a/2$ , noting that because  $v(x)$  is real  $\psi_l^*$  will be a solution to the same Schrödinger equation as  $\psi_l$ .



### 2.3.1 Solution

With the wave incoming from the left, recall that when  $x \leq -\frac{a}{2}$ , the wave function is given by:

$$\psi_l(x) = e^{iKx} + |r|e^{i\delta}e^{-iKx} \quad (58)$$

with derivative  $\psi'_l(x) = iK(e^{iKx} - |r|e^{i\delta}e^{-iKx})$ . The complex conjugate of the wave function is:

$$\psi_l^*(x) = e^{-iKx} + |r|e^{-i\delta}e^{iKx} \quad (59)$$

with derivative  $\psi_l^*(x) = -iK(e^{-iKx} - |r|e^{-i\delta}e^{iKx})$ . Now, we can evaluate the Wronskian for  $x \leq -\frac{a}{2}$ :

$$w(\psi_l, \psi_l^*) = \psi'_l \psi_l^* - \psi_l \psi_l^{*'} = iK(e^{iKx} - |r|e^{i\delta}e^{-iKx})(e^{-iKx} + |r|e^{-i\delta}e^{iKx}) \quad (60)$$

$$- (e^{iKx} + |r|e^{i\delta}e^{-iKx})(-iK(e^{-iKx} - |r|e^{-i\delta}e^{iKx})) \quad (61)$$

$$= iK(1 + |r|e^{-i\delta}e^{2iKx} - |r|e^{i\delta}e^{-2iKx} - |r|^2) \quad (62)$$

$$+ iK(1 - |r|e^{-i\delta}e^{2iKx} + |r|e^{i\delta}e^{-2iKx} - |r|^2) \quad (63)$$

$$= 2iK(1 - |r|^2) \quad (64)$$

With the wave coming from the left, recall that when  $x \geq \frac{a}{2}$ , the wave function is given by:

$$\psi_l(x) = |t|e^{i(\delta+Kx)} \quad (65)$$

with derivative  $\psi'_l(x) = iK|t|e^{i(\delta+Kx)}$ . The complex conjugate of the wave function is:

$$\psi_l^*(x) = |t|e^{-i(\delta+Kx)} \quad (66)$$

with derivative  $\psi_l^*(x) = -iK|t|e^{-i(\delta+Kx)}$ . Now, we can evaluate the Wronskian for  $x \geq \frac{a}{2}$ :

$$w(\psi_l, \psi_l^*) = \psi'_l \psi_l^* - \psi_l \psi_l^{*'} = iK|t|e^{i(\delta+Kx)}|t|e^{-i(\delta+Kx)} - |t|e^{i(\delta+Kx)}(-iK|t|e^{-i(\delta+Kx)}) \quad (67)$$

$$= 2iK|t|^2 \quad (68)$$

Setting that two results equal to each other, we find:

$$2iK(-|r|^2) = 2iK|t|^2 \implies 1 = |t|^2 + |r|^2 \quad (69)$$

## 2.4

(d) By evaluating  $w(\psi_l, \psi_r^*)$  prove that  $rt^*$  is pure imaginary, so  $r$  must have the form

$$r = \pm i|r|e^{i\delta}$$

where  $\delta$  is the same as in ( 8.71 )

### 2.4.1 Solution

We start by just considering the left boundary of  $x = -\frac{a}{2}$ , where the left-moving wave function is given by:

$$\psi_l(x) = e^{iKx} + re^{-iKx} \quad (70)$$

with derivative  $\psi'_l(x) = iK(e^{iKx} - re^{-iKx})$ . The complex conjugate of the right-moving wave function at the left boundary is:

$$\psi_r^*(x) = t^*e^{iKx} \quad (71)$$

with derivative  $\psi_r^{*'}(x) = iKt^*e^{iKx}$ . Now, we can evaluate the Wronskian for  $x \leq -\frac{a}{2}$ :

$$w(\psi_l, \psi_r^*) = \psi'_l \psi_r^* - \psi_l \psi_r^{*'} = iK(e^{iKx} - re^{-iKx})t^*e^{iKx} - (e^{iKx} + re^{-iKx})(iKt^*e^{iKx}) \quad (72)$$

$$= iKt^*(e^{2iKx} - r - e^{2iKx} - r) \quad (73)$$

$$= -2iKt^*r \quad (74)$$

Next, we consider the right boundary of  $x = \frac{a}{2}$ , where the left-moving wave function is given by:

$$\psi_l(x) = te^{iKx} \quad (75)$$

with derivative  $\psi'_l(x) = iKte^{iKx}$ . The complex conjugate of the right-moving wave function at the right boundary is:

$$\psi_r^*(x) = e^{iKx} + r^*e^{-iKx} \quad (76)$$

with derivative  $\psi_r^{*'}(x) = iK(e^{iKx} - r^*e^{-iKx})$ . Now, we can evaluate the Wronskian for  $x \geq \frac{a}{2}$ :

$$w(\psi_l, \psi_r^*) = \psi'_l \psi_r^* - \psi_l \psi_r^{*'} = iKte^{iKx}(e^{iKx} + r^*e^{-iKx}) - te^{iKx}(iK(e^{iKx} - r^*e^{-iKx})) \quad (77)$$

$$= iKt(e^{2iKx} + r^* - e^{2iKx} + r^*) \quad (78)$$

$$= 2iKtr^* \quad (79)$$

Setting the two results equal to each other, we find:

$$-2iKt^*r = 2iKtr^* = (-2iKt^*r)^* \quad (80)$$

For this to be true, we must have that  $rt^*$  is pure imaginary, so if we know  $t = |t|e^{i\delta} \implies t^* = |t|e^{-i\delta}$ , then  $r$  must have the form:  $r = \pm i|r|e^{i\delta}$  for the product  $rt^*$  to be pure imaginary.

## 2.5

(e) Show as a consequence of (8.70), (8.72) and (8.75) that the energy and wave vector of the Bloch electron are related by

$$\frac{\cos(Ka + \delta)}{|t|} = \cos ka, \quad \varepsilon = \frac{\hbar^2 K^2}{2m}$$

Since  $|t|$  is always less than one, but approaches unity for large  $K$  (the barrier becomes increasingly less effective as the incident energy grows). the left side of (8.76) plotted against

$K$  has the structure depicted in Figure 8.6. For a given  $k$ , the allowed values of  $K$  (and hence the allowed energies  $\varepsilon(k) = \hbar^2 K^2 / 2m$ ) are given by the intersection of the curve in Figure 8.6 with the horizontal line of height  $\cos(ka)$ . Note that values of  $K$  in the neighborhood of those satisfying

$$Ka + \delta = n\pi$$

give  $|\cos(Ka + \delta)| > 1$ , and are therefore not allowed for any  $k$ . The corresponding regions of energy are the energy gaps. If  $\delta$  is a bounded function of  $K$  (as is generally the case), there will be infinitely many regions of forbidden energy, and also infinitely many regions of allowed energies for each value of  $k$ .

### 2.5.1 Solution

We started with the recently derived definition of the transmission and reflection coefficients:

$$t = |t|e^{i\delta} \quad (81)$$

$$r = \pm i|r|e^{i\delta} \quad (82)$$

And then squaring these:

$$t^2 = |t|^2 e^{2i\delta} \quad (83)$$

$$r^2 = -|r|^2 e^{2i\delta} \quad (84)$$

We can now plug these into the expression for the cosine of the wave vector:

$$\cos(ka) = \frac{t^2 - r^2}{2t} e^{iGa} + \frac{1}{2t} e^{-iGa} \quad (85)$$

$$= \frac{|t|^2 e^{2i\delta} + |r|^2 e^{2i\delta}}{2|t|e^{i\delta}} e^{iGa} + \frac{1}{2|t|e^{i\delta}} e^{-iGa} \quad (86)$$

$$(87)$$

Now use the identity that the sum of the square magnitudes is unity in the first term and then combining all the exponentials

$$\cos(ka) = \frac{e^{2i\delta}}{2|t|e^{i\delta}} e^{iGa} + \frac{1}{2|t|e^{i\delta}} e^{-iGa} = \frac{e^{i(\delta+Ga)} + e^{-i(\delta+Ga)}}{2|t|} \quad (88)$$

Now, we make use of the exponential sum identity  $e^{i\theta} + e^{-i\theta} = 2\cos(\theta)$  to simplify the expression:

$$\cos(ka) = \frac{2\cos(\delta + Ga)}{2|t|} = \frac{\cos(\delta + Ga)}{|t|} \quad (89)$$

## 2.6

(f) Suppose the barrier is very weak (so that  $|t| \approx 1$ ,  $|r| \approx 0$ ,  $\delta \approx 0$ ). Show that the energy gaps are then very narrow, the width of the gap containing  $K = n\pi/a$  being

$$\varepsilon_{\text{gap}} \approx 2\pi n \frac{\hbar^2}{ma^2} |r|$$

### 2.6.1 Solution

We are interested in finding the width of the anergy gap  $\Delta\epsilon$  near a given value of  $K = n\pi/a$ . Since we are interested in small deviations from  $K$ , we want to consider  $K = n\pi/a + \Delta K$ . Then, we have:

$$Ka + \delta = n\pi + \Delta Ka + \delta = n\pi + \Delta\theta \quad (90)$$

but since we know that the  $\delta$  term is small, we can define a  $\Delta\theta$  where  $n\pi + \Delta\theta = n\pi + \Delta Ka$  and then we have  $\cos(Ka + \delta) = \cos(n\pi + \Delta\theta) = (-1)^n \cos(\Delta\theta)$ . Now, since we know that the transmission coefficient will be nearly 1, we are really interested in its inverse that appears in the equation, so if we define  $t = 1 - \epsilon$ , then we have  $\frac{1}{t} = \frac{1}{1-\epsilon} = 1 + \epsilon + \dots$ . Plugging in everything we just found gives:

$$\cos(ka) = \frac{\cos(Ka + \delta)}{|t|} = \frac{(-1)^n \cos(\Delta\theta)}{1 - \epsilon} \approx (-1)^n \cos(\Delta\theta)(1 + \epsilon) \quad (91)$$

Now we know that the magnitude of the cosine function is close to 1 near the energy gap, implying:

$$|\cos(\Delta\theta)(1 + \epsilon)| \approx 1 \implies |\cos(\Delta\theta)| \approx \frac{1}{1 + \epsilon} \approx 1 - \epsilon \implies \cos(\Delta\theta) \approx \pm(1 - \epsilon) \quad (92)$$

Now, we know that the Taylor expansion of the cosine function for a small argument is  $\cos(\Delta\theta) \approx 1 - \frac{1}{2}\Delta\theta^2$ , so we equate the two to find:

$$1 - \epsilon \approx 1 - \frac{1}{2}\Delta\theta^2 \implies \epsilon \approx \frac{1}{2}\Delta\theta^2 \implies \Delta\theta \approx \sqrt{2\epsilon} \quad (93)$$

Now, we know that we originally defined the  $\Delta\theta$  as  $\Delta\theta = n\pi + \frac{\pi}{a}\Delta K$ , so we can solve for  $\Delta K$ :

$$\Delta K = \frac{a}{\pi} \left( \sqrt{2\epsilon} - n\pi \right) = \frac{a}{\pi} \sqrt{2\epsilon} - n \quad (94)$$

And the energy is defined as  $\Delta\epsilon = \epsilon(K + \Delta K) - \epsilon(K) = \frac{\hbar^2}{2m} (K + \Delta K)^2 - \frac{\hbar^2}{2m} K^2 \approx \frac{\hbar^2}{m} K \Delta K$ , where we have neglected the second order term  $\Delta K^2$ . Since we know the relationship  $\Delta K = \Delta\theta/a = \sqrt{2\epsilon}/a$ , which we plug in along with  $K = n\pi/a$  to find:

$$\Delta\epsilon = \frac{\hbar^2 n\pi}{ma^2} \sqrt{2\epsilon} = 2\pi n \frac{\hbar^2}{ma^2} |r| \quad (95)$$

where in the final expression we make use of the fact that we defined initially for  $\epsilon$  as measuring the difference from full transmission, so we have  $\epsilon = \frac{1}{2}|r|^2$ .

## 2.7

(g) Suppose the barrier is very strong, so that  $|t| \approx 0, |r| \approx 1$ . Show that allowed bands of energies are then very narrow, with widths

$$\epsilon_{\max} - \epsilon_{\min} = O(|t|)$$

### 2.7.1 Solution

Again, we start with the relation

$$\cos(ka) = \frac{\cos(Ka + \delta)}{|t|} \quad (96)$$

but we know that the left hand side always is grounded by one, even though we have that the denominator on the right hand side is going to 0, which implies that the numerator too must banish:  $\cos(Ka + \delta) \approx 0$ . Now, we know that the cosine function is zero at integer multiples of  $\frac{\pi}{2}$ , so we have  $Ka + \delta = (n + \frac{1}{2})\pi$ , which implies that  $\delta = (n + \frac{1}{2})\pi - Ka$ . Now, we rearrange to have  $K_n = \frac{(n + \frac{1}{2})\pi - \delta}{a}$ . Now, this was always an approximation, so in reality there is a small deviation away from 0 or  $K = K_n + \Delta K$ , so we have:

$$K = \frac{(n + \frac{1}{2})\pi - \delta}{a} + \Delta K \implies Ka + \delta = \left(n + \frac{1}{2}\right)\pi + a\Delta K \quad (97)$$

We can use the trigonometric identity of  $\cos(\theta + \frac{\pi}{2}) = -\sin(\theta)$  to find:

$$\cos(Ka + \delta) = \cos\left(\left(n + \frac{1}{2}\right)\pi + a\Delta K\right) = -\sin(a\Delta K) \quad (98)$$

Now, our original relation is

$$\cos(ka) = \frac{\cos(Ka + \delta)}{|t|} = \frac{-\sin(a\Delta K)}{|t|} \quad (99)$$

and recall that the Corson function on the left has to be bounded by 1, so we must have

$$\left|\frac{-\sin(a\Delta K)}{|t|}\right| \leq 1 \implies |\sin(a\Delta K)| \leq |t| \quad (100)$$

We know that in this situation the transmission coefficient is supposed to be small, so we can utilize the tailor expansion of the sign function at small in goes to get:

$$|\sin(a\Delta K)| \approx |a\Delta K| \leq |t| \implies \Delta K \leq \frac{|t|}{a} \quad (101)$$

This means that the with between allowed K values is  $K_{\text{width}} = \frac{2|t|}{a}$ , and the energy width is then:

$$\Delta\epsilon = \frac{h^2}{2m} \left(K + \frac{2|t|}{a}\right)^2 - \frac{h^2}{2m} K^2 = \frac{h^2}{m} K \frac{2|t|}{a} + \frac{h^2}{2m} \left(\frac{2|t|}{a}\right)^2 = O(|t|) \quad (102)$$

## 2.8

(h) As a concrete example, one often considers the case in which  $v(x) = g\delta(x)$ , where  $\delta(x)$  is the Dirac delta function (a special case of the "Kronig-Penney model"). Show that in this case

$$\cot \delta = -\frac{h^2 K}{mg}, \quad |t| = \cos \delta$$

This model is a common textbook example of a one-dimensional periodic potential. Note, however, that most of the structure we have established is, to a considerable degree, independent of the particular functional dependence of  $|t|$  and  $\delta$  on  $K$ .

### 2.8.1 Solution

We know that in the case of a delta function potential, the width  $a$  of the potential well is 0. We start with the Schrödinger equation and then by enforcing continuity of the right-moving plane wave at either side of the origin with

$$\psi_L(x) = e^{iKx} + re^{-iKx} \quad (103)$$

$$\psi_R(x) = te^{iKx} \quad (104)$$

which leads us to the relationship  $1 + r = t$ . Next, we apply the fundamental theorem of calculus, which gives a relationship between the derivative from both sides of the origin and the area under the curve under the delta function:

$$\psi'_R(0^+) - \psi'_L(0^-) = \frac{2m}{\hbar^2} g\psi(0) \quad (105)$$

So we want to compute the appropriate derivatives:

$$\psi'_L(x) = iK(e^{iKx} - re^{-iKx}) \implies \psi'_L(0^-) = iK(1 - r) \quad (106)$$

$$\psi'_R(x) = iKte^{iKx} \implies \psi'_R(0^+) = iKt \quad (107)$$

and then we can plug in the values to find:

$$iKt - iK(1 - r) = \frac{2m}{\hbar^2} g(1 + r) \implies r = \frac{\frac{mg}{\hbar^2}}{iK - \frac{mg}{\hbar^2}} \implies t = \frac{iK}{iK - \frac{mg}{\hbar^2}} \quad (108)$$

Now, we can define the constant  $\lambda = \frac{mg}{\hbar^2}$  and then we can divide the numerator and denominator by  $iK$  to find:

$$t = \frac{-1}{-1 + \frac{\lambda}{iK}} = \frac{1}{1 + i\frac{\lambda}{K}} = \frac{1 - i\frac{\lambda}{K}}{1 + \frac{\lambda^2}{K^2}} \quad (109)$$

Now that we have defined  $t$  as a complex number, with a real part of 1 and an imaginary part of  $-\frac{\lambda}{K}$ , we know that  $\cot = \frac{\cos}{\sin}$  of an arbitrary complex number is the real part divided by the imaginary part, so we have:

$$\cot \delta = \frac{1}{-\frac{\lambda}{K}} = -\frac{K}{\lambda} = -\frac{h^2 K}{mg} \quad (110)$$

Now, the real part is defined as  $|t| \cos(\delta)$ , but we can factor  $t$ :

$$|t| = \frac{1}{\sqrt{1 + \frac{\lambda^2}{K^2}}} \frac{1 - i\frac{\lambda}{K}}{\sqrt{1 + \frac{\lambda^2}{K^2}}} \quad (111)$$

such that the second factor has unit norm:

$$|t|^2 = \frac{1}{1 + \frac{\lambda^2}{K^2}} \quad (112)$$

We notice that this is identical to  $|t| \cos(\delta)$ , so we have:

$$|t|^2 = |t| \cos(\delta) \implies |t| = \cos(\delta) \quad (113)$$

### 3

#### 3.1

(a) In the free electron case the density of levels at the Fermi energy can be written in the form (Eq. (2.64))  $g(\varepsilon_F) = mk_F/\hbar^2\pi^2$ . Show that the general form (8.63) reduces to this when  $\varepsilon_n(\mathbf{k}) = \hbar^2 k^2/2m$ .

##### 3.1.1 Solution

We start with the general form of the density of states

$$g_n(\mathcal{E}) = \int_{S_n^{(k)}} \frac{dS}{4\pi^3} \frac{1}{|\nabla \varepsilon_n(\mathbf{k})|} \quad (114)$$

Our first task is to avoid with the gradient of the single particle energy for the free electron

$$\nabla \varepsilon_n(\mathbf{k}) = \frac{\partial \varepsilon_n(\mathbf{k})}{\partial k} = \frac{\partial \left( \frac{\hbar^2 k^2}{2m} \right)}{\partial k} = \frac{\hbar^2 k}{m} \quad (115)$$

Then we want to do the integral expression, where we know that the kernel of the integral is independent of the surface parameter  $S$

$$\int_{S_n^{(k)}} dS = 4\pi k_F^2 \quad (116)$$

So, putting all of this together we see that

$$g_n(\varepsilon_F) = \frac{4\pi k_F^2}{4\pi^3} \frac{m}{\hbar^2 k_F} = \frac{mk_F}{\hbar^2 \pi^2} \quad (117)$$

#### 3.2

(b) Consider a band in which, for sufficiently small  $k$ ,  $\varepsilon_n(k) = \varepsilon_0 + (\hbar^2/2) (k_x^2/m_x + k_y^2/m_y + k_z^2/m_z)$  (as might be the case in a crystal of orthorhombic symmetry) where  $m_x, m_y$ , and  $m_z$  are positive constants. Show that if  $\varepsilon$  is close enough to  $\varepsilon_0$  that this form is valid, then  $g_n(\varepsilon)$  is proportional to  $(\varepsilon - \varepsilon_0)^{1/2}$ , so its derivative becomes infinite (van Hove singularity) as  $\varepsilon$  approaches the band minimum. (Hint: Use the form (8.57) for the density of levels.) Deduce from this that if the quadratic form for  $\varepsilon_n(\mathbf{k})$  remains valid up to  $\delta_F$ , then  $g_n(\varepsilon_F)$  can be written in the obvious generalization of the free electron form (2.65):

$$g_n(\varepsilon_F) = \frac{3}{2} \frac{n}{\varepsilon_F - \varepsilon_0}$$

where  $n$  is the concentration of the electrons in the band to the total electronic density.

### 3.2.1 Solution

We start by defining an effective shifted energy  $E = \varepsilon - \varepsilon_0 = (\hbar^2/2) (k_x^2/m_x + k_y^2/m_y + k_z^2/m_z)$ , and then to make the integral simpler, we define  $q_x = \frac{k_x}{\sqrt{m_x}}$ ,  $q_y = \frac{k_y}{\sqrt{m_y}}$ , and  $q_z = \frac{k_z}{\sqrt{m_z}}$ . Then we can write the integral as

$$g_n(\varepsilon) = \int \frac{d\mathbf{k}}{4\pi^3} \delta(\varepsilon - \varepsilon_n(\mathbf{k}))$$

, we know  $d\mathbf{k} = dk_x dk_y dk_z = \sqrt{m_x m_y m_z} dq_x dq_y dq_z$ , and then we can write the integral as

$$g_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{4\pi^3} \int dq_x dq_y dq_z \delta(\varepsilon - (\hbar^2/2) (q_x^2 + q_y^2 + q_z^2)) \quad (118)$$

Because the delta function will vanish for all values of  $q_x, q_y, q_z$  that do not satisfy the energy condition, we define  $q = \sqrt{2\varepsilon/\hbar^2}$ , and then we can write the integral as

$$g_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{4\pi^3} \int d\mathbf{q} \delta(\varepsilon - (\hbar^2/2) q^2) \quad (119)$$

Now we make use of the known property of the delta function, which is that  $\int d\mathbf{q} \delta(f) = \int \frac{dS_q}{|\nabla f|}$ , where  $dS_q$  is the surface element of the constant energy surface. First, we evaluate the gradient of the function

$$\nabla f = \frac{\partial f}{\partial q} = -\frac{\hbar^2}{1} q \implies |\nabla f| = \frac{\hbar^2}{1} q \quad (120)$$

And then the surface area of a sphere is  $4\pi q^2$ , so we can write the integral as

$$g_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{4\pi^3} \frac{4\pi q^2}{\hbar^2 q} = \frac{\sqrt{m_x m_y m_z}}{\pi^2 \hbar^2} q = \frac{\sqrt{m_x m_y m_z}}{\pi^2 \hbar^3} \sqrt{2E} \quad (121)$$

So, we have shown the desired result of the density of states being proportional to  $\sqrt{\varepsilon - \varepsilon_0}$ . Now, we want to find the derivative of the density of states, which is

$$g'_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{2\pi^2 \hbar^3} \frac{1}{\sqrt{2E}} \propto \frac{1}{\sqrt{\varepsilon - \varepsilon_0}} \quad (122)$$

And thus we observe the van Hove singularity because as the energy gap becomes small, the derivative of the density of states will go to infinity. Now we know that the number of electrons can be found by integrating the density of states up to the Fermi level:

$$n = \int_{\varepsilon_0}^{\varepsilon_F} g_n(\varepsilon) d\varepsilon = \frac{\sqrt{2m_x m_y m_z}}{\pi^2 \hbar^3} \int_{\varepsilon_0}^{\varepsilon_F} \sqrt{\varepsilon - \varepsilon_0} d\varepsilon \quad (123)$$

We can make a  $u$  substitution where  $u = \varepsilon - \varepsilon_0$ , so  $du = d\varepsilon$ , and then we can write the integral as

$$n = \frac{\sqrt{2m_x m_y m_z}}{\pi^2 \hbar^3} \int_0^{\varepsilon_F - \varepsilon_0} \sqrt{u} du = \frac{\sqrt{2m_x m_y m_z}}{\pi^2 \hbar^3} \frac{2}{3} (\varepsilon_F - \varepsilon_0)^{3/2} \quad (124)$$



And then rearrange to notice the equality

$$\frac{\sqrt{2m_x m_y m_z}}{\pi^2 \hbar^3} = \frac{3n}{2(\varepsilon_F - \varepsilon_0)^{3/2}} \quad (125)$$

so we can write the density of states as

$$g_n(\varepsilon_F) = \frac{3n}{2(\varepsilon_F - \varepsilon_0)} \quad (126)$$

### 3.3

(c) Consider the density of levels in the neighborhood of a saddle point, where  $\varepsilon_n(k) = \varepsilon_0 + (\hbar^2/2)(k_x^2/m_x + k_y^2/m_y - k_z^2/m_z)$  where  $m_x, m_y$ , and  $m_z$  are positive constants. Show that when  $\varepsilon \approx \varepsilon_0$ , the derivative of the density of levels has the form

$$\begin{aligned} g'_n(\varepsilon) &\approx \text{constant}, & \varepsilon > \varepsilon_0 \\ &\approx (\varepsilon_0 - \varepsilon)^{-1/2}, & \varepsilon < \varepsilon_0 \end{aligned}$$

#### 3.3.1 Solution

We previously found that the density of levels can be expressed as:

$$g_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{4\pi^3} \int dq_x dq_y dq_z \delta(E - (\hbar^2/2)(q_x^2 + q_y^2 - q_z^2)) \quad (127)$$

where we defined  $E = \varepsilon - \varepsilon_0$ . We know that the delta function will vanish for all values of  $q_x, q_y, q_z$  suggesting  $E = \frac{\hbar^2}{2}(q_x^2 + q_y^2 - q_z^2)$ . Therefore, if we have  $s \equiv \frac{2E}{\hbar^2}$ , this implies that  $q_z = \pm\sqrt{q_x^2 + q_y^2 - s}$ . Next, we want to evaluate the integral over  $q_z$  and by the properties of the delta function within the integrand, we know that

$$\int dq_z \delta(q_x^2 + q_y^2 - q_z^2 - s) = \frac{1}{2|q_z|} \Big|_{\pm\sqrt{q_x^2 + q_y^2 - s}} \quad (128)$$

But we need to multiply by a factor of 2 to account for both roots, so we are left with

$$\int dq_z \delta(q_x^2 + q_y^2 - q_z^2 - s) = \frac{1}{|q_z|} \Big|_{\sqrt{q_x^2 + q_y^2 - s}} \quad (129)$$

But we know that the expression under the square root must not be negative, so we split into 2 cases:  $s < 0$  and  $s > 0$ . For the case of  $s > 0$ , we have  $q_x^2 + q_y^2 \geq s$ . Therefore, we can approximate the integral as

$$g_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{4\pi^3} \int_{q_x^2 + q_y^2 \geq s} dq_x dq_y \frac{1}{\sqrt{q_x^2 + q_y^2 - s}} \quad (130)$$

Now, we can make a substitution  $q_x = r \cos(\theta)$  and  $q_y = r \sin(\theta)$ , so we have  $dq_x dq_y = r dr d\theta$ , and  $q_x^2 + q_y^2 = r^2$ , so the integral becomes

$$g_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{4\pi^3} \int_0^{2\pi} d\theta \int_{\sqrt{s}}^R \frac{r dr}{\sqrt{r^2 - s}} = \frac{\sqrt{m_x m_y m_z}}{2\pi^2} \int_{\sqrt{s}}^R \frac{r dr}{\sqrt{r^2 - s}} \quad (131)$$

Note that because we are integrating over a surface, the upper bound of the integral will be the radius  $R$  and for the denominator to be real, we must have  $r > \sqrt{s}$ . Performing the integral gives

$$g_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{2\pi^2} \left[ \sqrt{r^2 - s} \right]_{\sqrt{s}}^R = \frac{\sqrt{m_x m_y m_z}}{2\pi^2} \left[ \sqrt{R^2 - s} - 0 \right] \quad (132)$$

But we know that the energy gap  $s$  is very small, so  $\sqrt{R^2 - s} \approx R$ , so we have

$$g_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{2\pi^2} R \implies g'_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{2\pi^2} \propto \text{constant} \quad (133)$$

This is the result for the case of  $s > 0$ , where  $\varepsilon > \varepsilon_0$ . For the case of  $s < 0$ , we have  $q_x^2 + q_y^2 \leq s$ , we follow a similar procedure, but we now know that  $r^2 - s > 0$ , so we can write the integral as

$$g_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{4\pi^3} \int_0^{2\pi} d\theta \int_0^R \frac{r dr}{\sqrt{r^2 - s}} = \frac{\sqrt{m_x m_y m_z}}{2\pi^2} \int_0^R \frac{r dr}{\sqrt{r^2 - s}} \quad (134)$$

where now the lower bound is 0. Performing the integral gives

$$g_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{2\pi^2} \left[ \sqrt{r^2 - s} \right]_0^R = \frac{\sqrt{m_x m_y m_z}}{2\pi^2} \left[ \sqrt{R^2 - s} - \sqrt{-s} \right] \quad (135)$$

We make the same approximation as before, and then the derivative of the density of states will be

$$g'_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{2\pi^2} \left[ -\frac{1}{2\sqrt{R^2 - s}} + \frac{1}{2\sqrt{-s}} \right] = \frac{\sqrt{m_x m_y m_z}}{2\pi^2} \left[ -\frac{1}{2R} + \frac{1}{2\sqrt{-s}} \right] \quad (136)$$

Notice that because the gap is so small, the first term will have a much smaller contribution than the second term, so we can approximate the derivative as

$$g'_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{2\pi^2} \left[ \frac{1}{2\sqrt{-s}} \right] \propto (\varepsilon_0 - \varepsilon)^{-1/2} \quad (137)$$

In order to find the density of states, we will need to evaluate this integral but then right after we will take the derivative to determine  $g'_n(\varepsilon)$ , which implies that  $g'_n(\varepsilon)$  is a constant. For the case of  $s < 0$ , the integral over  $q_z$  will be

$$\int dq_z \delta(q_x^2 + q_y^2 - q_z^2 - s) = \frac{1}{\sqrt{q_x^2 + q_y^2 + s}} \quad (138)$$

and then the density of states will be

$$g_n(\varepsilon) = \frac{\sqrt{m_x m_y m_z}}{4\pi^3} \int_{q_x^2 + q_y^2 - s} dq_x dq_y \frac{1}{\sqrt{q_x^2 + q_y^2 + s}} \quad (139)$$

## 4 Problem 4: Ashcroft & Mermin, Chapter 9, Problem 1

### 1. Nearly Free Electron Fermi Surface Near a Single Bragg Plane

To investigate the nearly free electron band structure given by (9.26) near a Bragg plane, it is convenient to measure the wave vector  $\mathbf{q}$  with respect to the point  $\frac{1}{2}\mathbf{K}$  on the Bragg plane. If we write  $\mathbf{q} = \frac{1}{2}\mathbf{K} + \mathbf{k}$ , and resolve  $\mathbf{k}$  into its components parallel ( $k_{\parallel}$ ) and perpendicular ( $k_{\perp}$ ) to  $\mathbf{K}$ , then (9.26) becomes

$$\varepsilon = \varepsilon_{K/2}^0 + \frac{\hbar^2}{2m}k^2 \pm \left( 4\varepsilon_{K/2}^0 \frac{\hbar^2}{2m}k_{\perp}^2 + |U_K|^2 \right)^{1/2}$$

It is also convenient to measure the Fermi energy  $\varepsilon_F$  with respect to the lowest value assumed by either of the bands given by (9.36) in the Bragg plane, writing:

$$\varepsilon_F = \varepsilon_{K/2}^0 - |U_K| + \Delta$$

so that when  $\Delta < 0$ , no Fermi surface intersects the Bragg plane.

### 4.1

(a) Show that when  $0 < \Delta < 2|U_K|$ , the Fermi surface lies entirely in the lower band and intersects the Bragg plane in a circle of radius

$$\rho = \sqrt{\frac{2m\Delta}{\hbar^2}}$$

#### 4.1.1 Solution

We know that the lower band will have an energy of

$$\varepsilon = \varepsilon_{K/2}^0 + \frac{\hbar^2}{2m}k^2 - \left( 4\varepsilon_{K/2}^0 \frac{\hbar^2}{2m}k_{\perp}^2 + |U_K|^2 \right)^{1/2} \quad (140)$$

Equating the definition for the Fermi energy to this expression

$$\varepsilon_{K/2}^0 - |U_K| + \Delta = \varepsilon_{K/2}^0 + \frac{\hbar^2}{2m}k^2 - \left( 4\varepsilon_{K/2}^0 \frac{\hbar^2}{2m}k_{\perp}^2 + |U_K|^2 \right)^{1/2} \quad (141)$$

Canceling out the common energy factor and then rearranging gives

$$\frac{\hbar^2}{2m}k^2 + |U_K| - \Delta = \left( 4\varepsilon_{K/2}^0 \frac{\hbar^2}{2m}k_{\perp}^2 + |U_K|^2 \right)^{1/2} \quad (142)$$

Then, we square both sides to get

$$\left( \frac{\hbar^2}{2m}k^2 + |U_K| - \Delta \right)^2 = 4\varepsilon_{K/2}^0 \frac{\hbar^2}{2m}k_{\perp}^2 + |U_K|^2 \quad (143)$$

Now we know that at the Bragg plane,  $k_i = 0$ , so we can simplify the equation to

$$\left( \frac{\hbar^2}{2m} k^2 + |U_K| - \Delta \right)^2 = |U_K|^2 \quad (144)$$

Expanding the left side while defining  $S = |U_K| - \Delta$  gives

$$\left( \frac{\hbar^2}{2m} \right)^2 k^4 + 2 \left( \frac{\hbar^2}{2m} \right) k^2 S + (S^2 - |U_K|^2) = 0 \quad (145)$$

This is a quadratic equation in  $k^2$ , so first we redefine  $E_k \equiv \frac{\hbar^2}{2m} k^2$  and then we recognize a further simplification of

$$S^2 - |U_K|^2 = \Delta^2 - 2\Delta|U_K| + |U_K|^2 - |U_K|^2 = \Delta^2 - 2\Delta|U_K| \quad (146)$$

So in full form, our quadratic equation is

$$E_k^2 + 2(|U_K| - \Delta) E_k + \Delta^2 - 2\Delta|U_K| = 0 \quad (147)$$

Now we use the quadratic formula to solve

$$E_k = \frac{-2(|U_K| - \Delta) \pm \sqrt{4(|U_K| - \Delta)^2 - 4\Delta^2 + 8\Delta|U_K|}}{2} \quad (148)$$

$$= -(|U_K| - \Delta) \pm \sqrt{|U_K|^2 - 2\Delta|U_K| + \Delta^2 - \Delta^2 + 2\Delta|U_K|} \quad (149)$$

$$= -(|U_K| - \Delta) \pm \sqrt{|U_K|^2} \quad (150)$$

$$(151)$$

We need to choose the physical solution, so we get

$$E_k = -(|U_K| - \Delta) + |U_K| = \Delta \implies \frac{\hbar^2}{2m} k^2 = \Delta \implies k = \sqrt{\frac{2m\Delta}{\hbar^2}} \quad (152)$$

At the Bragg plane we have  $k_{\parallel} = k$ , so the radius of the circle where the Fermi surface intersects the Bragg plane is  $\rho = k = \sqrt{\frac{2m\Delta}{\hbar^2}}$ .

## 4.2

(b) Show that if  $\Delta > |2U_K|$ , the Fermi surface lies in both bands, cutting the Bragg plane in two circles of radii  $\rho_1$  and  $\rho_2$  (Figure 9.6), and that the difference in the areas of the two circles is

$$\pi (\rho_2^2 - \rho_1^2) = \frac{4m\pi}{\hbar^2} |U_K|$$

(The area of these circles can be measured directly in some metals through the de Haas-van Alphen effect (Chapter 14), and therefore  $|U_K|$  can be determined directly from experiment for such nearly free electron metals.)

### 4.2.1 Solution

Again, we can write out 9.26 as

$$\varepsilon = \varepsilon_{K/2}^0 + \frac{\hbar^2}{2m} k^2 \pm \left( 4\varepsilon_{K/2}^0 \frac{\hbar^2}{2m} k_i^2 + |U_K|^2 \right)^{1/2} \quad (153)$$

and since this is at the Bragg plane,  $k_i = 0$ , so we can simplify to

$$\varepsilon = \varepsilon_{K/2}^0 + \frac{\hbar^2}{2m} k_{||}^2 \pm |U_K| \quad (154)$$

Now, we set this equal to the expression for the Fermi energy

$$\varepsilon_{K/2}^0 - |U_K| + \Delta = \varepsilon_{K/2}^0 + \frac{\hbar^2}{2m} k_{||}^2 \pm |U_K| \implies \Delta - |U_K| = \frac{\hbar^2}{2m} k_{||}^2 \pm |U_K| \quad (155)$$

First, we will treat the case of the lower band, which has the minus sign, so we have

$$\Delta - |U_K| = \frac{\hbar^2}{2m} k_{||}^2 - |U_K| \implies \frac{\hbar^2}{2m} k_{||}^2 = \Delta \implies k_{||} = \sqrt{\frac{2m\Delta}{\hbar^2}} \quad (156)$$

Next, we do the same for the upper band, which has the plus sign, so we have

$$\Delta - |U_K| = \frac{\hbar^2}{2m} k_{||}^2 - |U_K| \implies \frac{\hbar^2}{2m} k_{||}^2 = \Delta - 2|U_K| \implies k_{||} = \sqrt{\frac{2m(\Delta - 2|U_K|)}{\hbar^2}} \quad (157)$$

So, the radii of the two circles are  $\rho_1 = \sqrt{\frac{2m\Delta}{\hbar^2}}$  and  $\rho_2 = \sqrt{\frac{2m(\Delta - 2|U_K|)}{\hbar^2}}$ , and the difference in the areas of the two circles is then

$$\pi(\rho_2^2 - \rho_1^2) = \pi \left( \frac{2m(\Delta - 2|U_K|)}{\hbar^2} - \frac{2m\Delta}{\hbar^2} \right) = \pi \left( \frac{4m|U_K|}{\hbar^2} \right) \quad (158)$$

## 5 Problem 5: Ashcroft & Mermin, Chapter 9, Problem 3

3. Effect of Weak Periodic Potential at Places in  $k$ -Space Where Bragg Planes Meet Consider the point  $W$  ( $k_w = (2\pi/a)(1, \frac{1}{2}, 0)$ ) in the Brillouin zone of the fcc structure shown (see Figure 9.14). Here three Bragg planes ( $(200)$ ,  $(111)$ ,  $(11\bar{1})$ ) meet, and accordingly the free electron energies

$$\begin{aligned} \varepsilon_1^0 &= \frac{\hbar^2}{2m} k^2 \\ \varepsilon_2^0 &= \frac{\hbar^2}{2m} \left( \mathbf{k} - \frac{2\pi}{a}(1, 1, 1) \right)^2 \\ \varepsilon_3^0 &= \frac{\hbar^2}{2m} \left( \mathbf{k} - \frac{2\pi}{a}(1, 1, \bar{1}) \right)^2 \\ \varepsilon_4^0 &= \frac{\hbar^2}{2m} \left( \mathbf{k} - \frac{2\pi}{a}(2, 0, 0) \right)^2 \end{aligned}$$

are degenerate when  $\mathbf{k} = k_w$ , and equal to  $\varepsilon_w = \hbar^2 \mathbf{k}_w^2 / 2m$ . Figure 9.14 First Brillouin zone for a face-centered cubic crystal.

## 5.1

(a) Show that in a region of  $k$ -space near W, the first-order energies are given by solutions to <sup>15</sup>

$$\begin{vmatrix} \varepsilon_1^0 - \varepsilon & U_1 & U_1 & U_2 \\ U_1 & \varepsilon_2^0 - \varepsilon & U_2 & U_1 \\ U_1 & U_2 & \varepsilon_3^0 - \varepsilon & U_1 \\ U_2 & U_1 & U_1 & \varepsilon_4^0 - \varepsilon \end{vmatrix} = 0$$

where  $U_2 = U_{200}$ ,  $U_1 = U_{111} = U_{11\bar{1}}$ , and that at W the roots are

$$\varepsilon = \varepsilon_W - U_2 \text{ (twice)}, \quad \varepsilon = \varepsilon_W + U_2 \pm 2U_1$$

### 5.1.1 Solution

As shown in lecture, the Hamiltonian for this weak potential has the matrix of elements

$$\varepsilon_i^0 - \varepsilon \text{ if } i = j \quad (159)$$

$$U_{\mathbf{G}_i - \mathbf{G}_j} \text{ if } i \neq j \quad (160)$$

Since we have 4 energy levels, the matrix becomes

$$\begin{bmatrix} \varepsilon_1^0 - \varepsilon & U_1 & U_1 & U_2 \\ U_1 & \varepsilon_2^0 - \varepsilon & U_2 & U_1 \\ U_1 & U_2 & \varepsilon_3^0 - \varepsilon & U_1 \\ U_2 & U_1 & U_1 & \varepsilon_4^0 - \varepsilon \end{bmatrix} \quad (161)$$

At the point W, we have that all  $\varepsilon_i^0 = \varepsilon_W$ , so we want to evaluate this determinant

$$\det(H) = 0 \implies \begin{vmatrix} x & U_1 & U_1 & U_2 \\ U_1 & x & U_2 & U_1 \\ U_1 & U_2 & x & U_1 \\ U_2 & U_1 & U_1 & x \end{vmatrix} = 0 \quad (162)$$

where  $x = \varepsilon_W - \varepsilon$ . Using symbolic algebra to compute the roots, we get

$$(-U_2 + x)^2 (-2U_1 + U_2 + x) (2U_1 + U_2 + x) = 0 \quad (163)$$

which gives the desired roots of  $\varepsilon = \varepsilon_W - U_2$  twice and  $\varepsilon = \varepsilon_W + U_2 \pm 2U_1$ .

## 5.2

(b) Using a similar method, show that the energies at the point U ( $\mathbf{k}_U = (2\pi/a)(1, \frac{1}{4}, \frac{1}{4})$ ) are

$$\varepsilon = \varepsilon_U - U_2, \quad \varepsilon = \varepsilon_U + \frac{1}{2}U_2 \pm \frac{1}{2}(U_2^2 + 8U_1^2)^{1/2}$$

where  $\varepsilon_U = \hbar^2 \mathbf{k}_U^2 / 2m$ .

### 5.2.1 Solution

First we want to find the deconvolved energies at the point  $U$ . This can be determined by considering where

$$|k_U|^2 = |k_U - G|^2 \implies 2\mathbf{k}_U \cdot \mathbf{G} - \mathbf{G}^2 = 0 \quad (164)$$

If we set  $\mathbf{G} = \frac{2\pi}{a}(h, k, l)$  and plug in the values for  $\mathbf{k}_U = (2\pi/a)(1, 1/4, 1/4)$ , we get

$$2 \left( h + \frac{k}{4} + \frac{l}{4} \right) = (h^2 + k^2 + l^2) \quad (165)$$

This becomes true only for the pairing (1,1,1) and (2,0,0), so the possible reciprocal lattices vectors are  $\mathbf{G}_1 = \frac{2\pi}{a}(1, 1, 1)$  and  $\mathbf{G}_2 = \frac{2\pi}{a}(2, 0, 0)$ . This implies that we want to consider to degenerate states for  $\mathbf{k}_U$ ,  $\mathbf{k}_U - \mathbf{G}_1$  and  $\mathbf{k}_U - \mathbf{G}_2$ . We can write the Hamiltonian matrix as

$$H = \begin{bmatrix} \varepsilon_1^0 - \varepsilon & U_1 & U_2 \\ U_1 & \varepsilon_2^0 - \varepsilon & U_1 \\ U_2 & U_1 & \varepsilon_3^0 - \varepsilon \end{bmatrix} \quad (166)$$

Then we have the same deal as before where all of the  $\varepsilon_i^0 = \varepsilon_U$ , so we can write the determinant as

$$\det(H) = 0 \implies \begin{vmatrix} x & U_1 & U_2 \\ U_1 & x & U_1 \\ U_2 & U_1 & x \end{vmatrix} = 0 \quad (167)$$

where  $x = \varepsilon_U - \varepsilon$ . Using symbolic algebra to compute the roots, we get

$$(-U_2 + x)(-2U_1^2 + U_2x + x^2) = 0 \quad (168)$$

which gives the desired roots of  $\varepsilon = \varepsilon_U - U_2$  and then solving the quadratic equation in the second parenthesis gives  $\varepsilon = \varepsilon_U + \frac{1}{2}U_2 \pm \frac{1}{2}(U_2^2 + 8U_1^2)^{1/2}$ .