

## ECE344, Midterm Exam 1: Solutions October 19, 2009

1. **Problem:** The workfunction of a material is defined as the energy required to remove an electron from the material and is usually indicated by the Greek letter  $\chi$  (pron: 'khi').

The workfunction of gold is  $\chi_{Au} = 4.9$  eV, the workfunction of cesium is  $\chi_{Cs} = 1.9$  eV.

a. Calculate the maximum wavelength (in nm) of light required to remove an electron from Au via the photoelectric effect.

b. Do the same for Cs.

**Solution:** Since

$$\lambda = \frac{c}{\nu} = \frac{2\pi\hbar c}{h\nu} = \frac{2\pi\hbar}{E}, \quad (1)$$

or, expressing the energy in eV:

$$\lambda = \frac{2\pi\hbar}{eE}, \quad (2)$$

using  $E = \chi_{Au}$  we have  $\lambda = 253$  nm and using  $E = \chi_{Cs}$  we get  $\lambda = 653$  nm.

2. **Problem:** Electrons in a 'swarm' at thermal equilibrium have each an average kinetic energy of  $(3/2)k_B T$ .

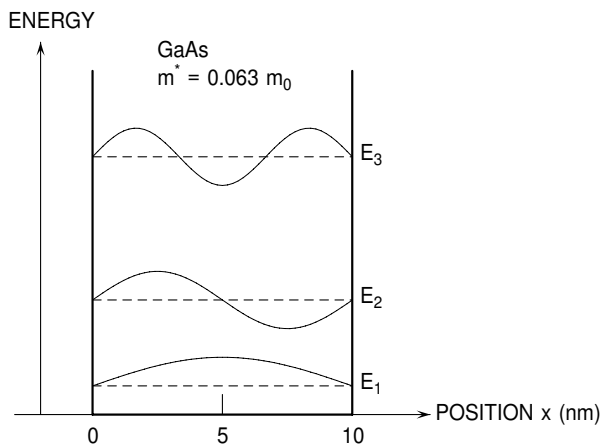
a. Calculate the value of  $(3/2)k_B T$  in eV.

b. Calculate the velocity (in cm/s) of an electron having a kinetic energy equal to the average energy  $(3/2)k_B T$ . This is the so-called 'thermal velocity'.

c. Calculate its average de Broglie wavevavelength in cm. This is called the 'thermal wavelength'.

**Solution:** Plugging the appropriate numbers, we get  $E_{th} = (3/2)k_B T = 38.77$  meV. From  $E_{th} = (1/2)mv_{th}^2$  we get for the thermal velocity  $v_{th} = \sqrt{2E_{th}/m} = 1.16 \times 10^7$  cm/s. Finally, from  $\lambda_{th} = h/(mv_{th})$  we get for the thermal wavelength  $\lambda_{th} = 6.23$  nm.

3. **Problem:** Consider a quantum well of GaAs (which is like the one-dimensional 'box' we have considered in class): Electrons are assumed to be in one dimension, assume the  $x$  axis to be along this direction, the potential is zero in the interval



**Fig. 1: Figure relative to problem 3**

$(0, L)$ , infinite outside, so the the wavefunction of the electrons must vanish at  $x = 0$  and  $x = L$ .

Using the effective mass of GaAs,  $m^* = 0.063m_0$  (where  $m_0 = 9.1 \times 10^{-31}$  Kg is the electron mass in vacuum), and assuming  $L = 10$  nm, calculate the energy (in eV) of the three lowest-energy levels in the quantum well.

Figure 1 illustrates schematically the potential well, the position of the first 3 energy levels in the well, and the corresponding wavefunctions.

**Solution:** The energy levels  $E_n$  of the particle in the box are given by:

$$E_n = \frac{\hbar^2}{2m^*} \left( \frac{n\pi}{L} \right)^2 = \frac{\hbar^2 \pi^2}{2m^* L^2} n^2. \quad (3)$$

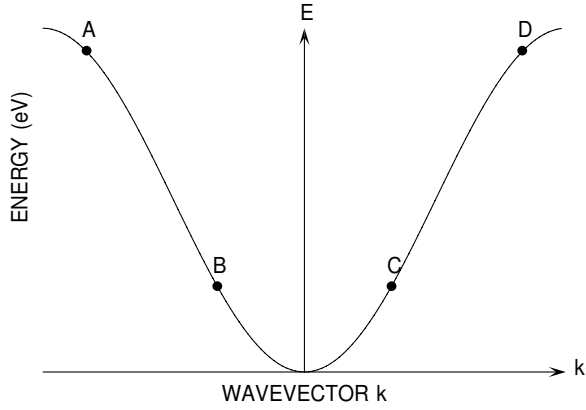
Using the values for  $L$  and  $m^*$  specified by the problem, we get  $E_n = 38n^2$  meV. So,  $E_1 = 38$  meV,  $E_2 = 152$  meV, and  $E_3 = 342$  meV.

4. **Problem:** Figure 2 shows a band-diagram for some ideal material along one particular direction in  $\mathbf{k}$ -space.

Indicate the sign of the effective mass and of the velocity (along the  $x$ -axis in Fig. 2) for electrons with wavevectors  $\mathbf{k}$  at each of the points labeled  $A$ ,  $B$ ,  $C$ , and  $D$ .

**Solution:** Recall that for our ‘isotropic’ case:

$$m^* = \frac{\hbar^2}{d^2 E / dk^2} \quad (4)$$



**Fig. 2: Figure relative to problem 4**

and

$$v = \frac{dE/dk}{\hbar}, \quad (5)$$

so that positive (negative) curvature implies a positive (negative) effective mass, a positive (negative) slope a positive (negative)  $x$ -component of the velocity. Thus:

At point A:  $m^* < 0$ ,  $v < 0$

At point B:  $m^* > 0$ ,  $v < 0$

At point C:  $m^* > 0$ ,  $v > 0$

At point D:  $m^* < 0$ ,  $v > 0$ .

Some confusion might have been caused by associating the momentum  $p = m^*v$  with  $\hbar k$ . This is correct only near the bottom of the band (positive effective mass) at  $k = 0$ , while near the top of the band (negative effective mass) the momentum is  $p \approx \hbar(k - k_0)$ , where  $k_0$  is the location of the band maximum.

5. **Problem:** Calculate the probability that a single state at the bottom of the conduction band will be occupied at 300 K in Si, Ge, and GaAs. Assume that the Fermi level is at midgap (that is:  $E_F = E_i = E_{mg}$  and use for the values of the gap  $E_g = 1.12$  eV (for Si), 0.64 eV (for Ge), 1.42 eV (for GaAs).

**Solution:** The probability of a state at the bottom of the conduction band to be occupied is given by the Fermi-Dirac distribution at  $E - E_F = E_{mg}$ . Thus we must evaluate

$$P = \left[ 1 + \exp\left(\frac{E_{mg}}{k_B T}\right) \right]^{-1}$$

for the three cases given, obtaining  $3.91 \times 10^{-10}$  for Si,  $4.21 \times 10^{-6}$  for Ge, and  $1.18 \times 10^{-12}$  for GaAs.

6. **Bonus problem.** Consider the conduction band of a crystal whose energy-vs-wavevector relation  $E(k)$  along some direction in  $\mathbf{k}$ -space is given by the following expression:

$$E(k) = E_0 + E_1 \cos[\alpha(k - k_0)] .$$

Note that this band has a minimum at  $k = k_0$ , pretty much like Si along the line from the symmetry point  $\Gamma$  to the symmetry point  $X$ .

Calculate the electron effective mass near the minimum of the band in terms of the parameters of the problem,  $E_0$ ,  $E_1$ ,  $k_0$  and  $\alpha$ . (Not all of them will necessarily appear in the final answer).

**Solution:** From Eq. (4) and from the fact that

$$\begin{aligned} \frac{dE(k)}{dk} &= -\alpha E_1 \sin[\alpha(k - k_0)] , \\ \frac{d^2E(k)}{dk^2} &= -\alpha^2 E_1 \cos[\alpha(k - k_0)] , \end{aligned}$$

which is  $-E_1\alpha^2$  at  $k = k_0$ , we have  $m^* = -\hbar/(\alpha^2 E_1)$ .