## ECE344 Fall08 HOMEWORK 2 Energy Band Theory and Carrier Densities

## 1 DOS

- 1. For a 3D semiconductor, develop an expression for the total number of available states/ $cm^3$  in the conduction band between energies  $E_c$  and  $E_c + \gamma k_B T$ , where  $\gamma$  is an arbitrary constant. Do the same for the valence band between energies  $E_v \gamma k_B T$  and  $E_v$ .
- 2. Using the effective mass approximation, derive the expression of the density of state (DOS) g(E) for a 2D and 1D electron gas. You will use the fact that  $E_1$  is the minimum potential energy for the electrons ( $E_1 > E_c$  due to quantization effects in systems with low-dimensionalities) such that

$$E = \frac{\hbar^2 k^2}{2m^*} + E_1$$

Also for 2D, we give the following density of state in k-space (per unit of Volume):

$$g(k)dk = 2 * \frac{S}{(2\pi)^2} 2\pi k dk,$$

and for 1D:

$$g(k)dk = 2 * \frac{L}{(2\pi)} 2dk$$

S for (2D) Surface and L for (1D) Length.

- 3. Calculate the number of states per unit of energy in a 3D 100nm by 100nm by 10nm piece of Silicon  $(m^* = 1.08m_0, m_0; \text{mass of the electron}) 100meV$  above the conduction band edge. Write the result in units of  $eV^{-1}$ . Do the same for a 100nm by 100nm 2D silicon sheet and 100nm 1D silicon wire (we consider  $E_1 = 40meV$ ).
- 4. For non-degenerate semiconductors, derive analytically the expressions of the effective density of states for a 2D and 1D system ( $N_c^{2D}$  and  $N_c^{1D}$ ).
- 5. In the general case (Fermi-Dirac distribution), derive analytically the expression of electron density for a 2D electron gas (you will give the expression in function of  $N_c^{2D}$ ).

## **2** Distribution function

- 1. Evaluate the approximation errors if one uses the Maxwell-Boltzmann statistics instead of the Fermi-Dirac statistics, for the following energies:  $E - E_F = 2.2k_BT$ ,  $E - E_F = 2.9k_BT$ ,  $E - E_F = 4.6k_BT$ . Comment.
- 2. Under equilibrium conditions and T > 0K, what is the probability of an electron state being occupied if it is located at the Fermi level ?, (ii) If  $E_F$  is positioned at  $E_c$ , determine (numerical answer) the probability of finding an electron at  $E_c + k_B T$ .

## 3 Energy band

In the energy band diagram below, indicate the usual positioning of the following energy levels: (i)  $E_i$ , intrinsic Fermi level, (ii)  $E_{F_1}$ , Fermi level for N-type non-degenerate semiconductor with  $N_D = 10^{15}/cm^3$ , (iii)  $E_{F_2}$ , Fermi level for N-type non-degenerate semiconductor with  $N_D = 10^{17}/cm^3$ , (iv)  $E_{vac}$ , vacuum energy level, (v)  $E_G$ , energy bandgap. Add comments as necessary to forestall any misinterpretation of your graphical answer.

*E<sub>c</sub>*\_\_\_\_\_

 $E_v$  \_\_\_\_\_