## HW #1

## Due Feb. 26 (Thu) in class

Please submit **Hard Copy**. Please remember to put down **your name** and **SID**.

- 1. We have derived the expressions for the 3-d (bulk) and 2-d (quantum well) electron density of states in class. Follow the same procedure, derive the 1-d electron density of state function for a "quantum wire" with dimension of  $L_x$  and  $L_y$  in the x and y direction, and unconfined in the z direction.
  - a. Derive the 1-D electron density of state function,  $\rho_{1D}(E)$ .
  - b. Find the first three energy levels for  $L_x = L_y = 5$  nm, assume the electron effective mass  $m_e^* = 0.1 \cdot m_0$ , where  $m_0$  is the free electron mass.
  - c. Find the electron concentration if the Fermi level is located just below the second energy level (as you find in Part b.). For simplicity, assume T = 0 Kevin.
- 2. Consider a "quantum box" (also called quantum dot or QD) with dimensions of  $L_x \times L_y \times L_z$ .
  - a. Derive the 0-D electron density of state function,  $\rho_{0D}(E)$ .
  - b. Find the first three energy levels for  $L_x = L_y = L_z = 5$  nm, assume the electron effective mass  $m_e^* = 0.1 \cdot m_0$ , where  $m_0$  is the free electron mass.
  - c. Find the electron concentration if the Fermi level is located just below the second energy level (as you find in Part b.). For simplicity, assume T = 0 Kevin.
- 3. Refer to the diagram on the right. Under biased condition, both conduction and valence bands are populated. The electron distribution in conduction band is described by Fermi-Dirac distribution,  $f_C(E_2)$ , with quasi-Fermi energy  $F_C$ . The electron distribution in valence band is described by Fermi-Dirac distribution,  $f_V(E_1)$ , with quasi-Fermi energy  $F_V$ . Here,  $E_1$  and  $E_2$  are related by an optical transition (i.e., they have the same k). For optical matrix element, use

$$E_z$$
 $E_z$ 
 $E_z$ 
 $f_C(E_z)$ 
 $f_C(E_z)$ 
 $f_V(E_1)$ 

$$\left| \hat{e} \cdot \vec{P}_{cv} \right|^2 = \frac{m_0}{6} E_p \text{ with } E_p = 25.7 \text{ eV}$$

- a. Use the energy reference below (i.e,  $E_V$  = 0 and  $E_C$  =  $E_g$ , the bandgap energy), find  $E_1$  and  $E_2$  as functions of the photon energy,  $h\omega$ .
- b. Derive  $f_{\mathcal{C}}(E_2(\hbar\omega))$  as a function of  $\hbar\omega$ .
- c. Derive  $f_{\nu}(E_{1}(\hbar\omega))$  as a function of  $\hbar\omega$ .
- d. Assuming  $E_g = 1$  eV,  $F_C F_V = 1.2$  eV,  $m_e^* = 0.1 m_0$ ,  $m_h^* = 0.4 m_0$ . Calculate and plot the emission probability  $f_e(\hbar\omega) = f_C(E_2(\hbar\omega)) \cdot \left[1 f_V(E_1(\hbar\omega))\right]$  for photon energies from 0.8 to 1.5 eV. Plot for two temperatures: T = 0 and T = 300 K.
- e. Repeat part d) for the Fermi inversion factor:  $f_g(\hbar\omega) = f_C(E_2(\hbar\omega)) f_V(E_1(\hbar\omega))$
- f. Plot the gain spectra for T = 0 and T = 300 K for the condition given in d).
- g. Plot the spontaneous emission spectra for T = 0 and T = 300 K for the condition given in d).