## ECE609 Spring07 HOMEWORK 2 Energy Band Theory and Semiconductor Fundamentals

## **1** Energy Band Theory (10pts)

- We consider a crystal with a periodic potential. The first and second form of the Bloch theorem are given by (i)  $\Psi_k(\mathbf{r} + \mathbf{R}) = e^{i\mathbf{k}\mathbf{R}}\Psi_k(\mathbf{r})$ , where  $\mathbf{R}$  has the periodicity of the crystal, and (ii)  $\Psi_k(\mathbf{r}) = e^{i\mathbf{k}\mathbf{r}}u_k(\mathbf{r})$  with  $u_k(\mathbf{r} + \mathbf{R}) = u_k(\mathbf{r})$ . From (ii) you will show (i). From (i) you will show (ii).
- The Schrödinger equation is given by:

$$-\frac{\hbar^2}{2m}\Delta_r\Psi(r) + U(r)\Psi(r) = E\Psi(r)$$

where U(r) is periodic. Using the second form of the Bloch theorem, derive the partial differential equation for  $u_k$ .

## 2 Energy band theory using the Linear Combination of Atomic Orbitals (LCAO) method (55pts)

We consider an infinite chain of atoms (going from  $-\infty$  to  $\infty$ ) where l is the distance between the atoms. For a given atom n, only one energy state  $v_n(x)$  is available. We will not consider the overlap between the different states (orthogonal tight-binding approach), in addition, the basis function  $\{v_n\}$  satisfies the normalization condition:

$$\int_{-\infty}^{+\infty} v_n(x)v_p(x) = \delta_{n,p}.$$

We would like to solve the Schrödinger equation  $H\Psi = E\Psi$ .

- 1. Write  $\Psi(x)$  as a linear combination of basis functions  $v_n(x)$  (atomic orbital- one by atom). We will introduce the coefficients  $c_n$ .
- 2. Show that  $c_n$  satisfies the following equation (we consider only the interactions between the first neighbors):

$$E_0 c_n - A c_{n+1} - A c_{n-1} = E c_n,$$

where you will give the expression of  $E_0$  (on site energy) and A (coupling term).

- 3. We consider solutions of this form:  $c_n = \exp(iknl)$ , where k belongs to the first Brillouin zone. Find and plot the dispersion relation. Comment. From the graph, comment on what would happen if the tunneling effect between atoms increases.
- 4. Using the fact that the state  $v_n$  can be obtained by translation nl of the state  $v_0$  (i.e.  $v_n(x) = v_0(x ql)$ ), demonstrate the first form of the Bloch theorem.
- 5. We set  $\Psi_k(x) = \exp(ikx)u_k(x)$ , which condition  $u_k$  should satisfy ?
- 6. In the atom chain, where is the electron associated with the state  $\Psi_k$ ?

- 7. We consider a finite chain L with N number of atoms such as N is very large. Using Born-Von Karman boundary conditions (periodic B.C.), which are the possible values for k? How many k states are available in the first brillouin zone (give the result in fonction of N, also we will not consider the spin factor). How can we approximate this expression if the number of atoms is very large (N >> 1).
- 8. Calculate and plot the group velocity of the electrons which is given by

$$V_G = \frac{1}{\hbar} \left( \frac{dE(k)}{dk} \right).$$

What is happening if the energy of the electron goes to  $E_0 + 2A$ ?

## **3** Carrier densities (35pts)

- 1. Derive the expression of the density of state (DOS) g(E) for a 2D and 1D electron gas. We will use the fact that  $E_n$  are the quantized energies inside the conduction band due to the confined direction(s). We will also consider that  $E_1 < E < E_2$ .
- 2. Calculate the number of states per unit energy in a 100nm by 100nm by 10nm piece of silicon ( $m^* = 1.08m_0, m_0$ : mass of the electron) 100 meV 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 silicon wire (we consider  $E_1 = 40meV$ ).
- 3. For non-degenerate semiconductors, derive analytically the expressions of the effective density of states for a 2D and 1D system ( $N_c^{2D}$  and  $N_c^{1D}$ ).
- 4. In the general case (Fermi-Dirac distribution), derive analytically the expression of electron density for a 2D electron gas (we will give the expression in function of  $N_c^{2D}$ )