

Tight-Binding Method: single atom

I. Motivation:  $H_{at} \psi_n = E_n \psi_n$  single atom w/  
 $H = H_{at} + \Delta U(r)$  with  $\Delta U(r)$  small  
 see Fig. 10.2

There is a set of approximate solutions:

$$\langle r | n, R \rangle = \psi_n(r - R).$$

Individually these do not satisfy Bloch's Thm;  
 however, there is a linear combination which  
 does:

$$\psi_{nk}(r) = \sum_{R'} e^{ik \cdot R'} \psi_n(r - R')$$

$$\begin{aligned} \psi_{nk}(r + R) &= \sum_{R'} e^{ik \cdot R'} \psi_n(r - R' + R) \quad ; \quad R'' = R' - R \\ & \quad \quad \quad R' = R'' + R \\ &= \sum_{R''} e^{ik \cdot (R'' + R)} \psi_n(r - R'') \\ &= e^{ik \cdot R} \psi_{nk}(r) // \end{aligned}$$

II. Derivation: Look for  $\psi_{nk}(r)$  of the form  
 (of tight binding) depends on k

$$|\psi\rangle = \sum_R e^{ik \cdot R} \sum_n b_n |n, R\rangle$$

$$= \sum_{R, n} e^{ik \cdot R} b_n |n, R\rangle$$

$$= \sum_n b_n |n, \vec{0}\rangle + \sum_{R \neq 0, n} e^{ik \cdot R} b_n |n, R\rangle$$

Solve for the  $b_n$ :

only small for w.f. centered about the origin

$$\begin{aligned}
 H|\psi\rangle &= (H_{\text{lat}} + \Delta U)|\psi\rangle \\
 &= H_{\text{lat}} \sum b_n |n, \vec{0}\rangle \\
 &+ \Delta U \sum_n b_n |n, \vec{0}\rangle \\
 &+ H_{\text{lat}} \sum'_R e^{ik \cdot R} b_n |n, \vec{R}\rangle \quad \text{\(\sum'\) means } R \neq 0 \\
 &+ \Delta U \sum_{n,R} e^{ik \cdot R} b_n |n, \vec{R}\rangle \\
 &= \epsilon \sum b_n |n, \vec{0}\rangle \\
 &+ \epsilon \sum'_{n,R} e^{ik \cdot R} b_n |n, \vec{R}\rangle
 \end{aligned}$$

Act on the LHS with  $\langle m, \vec{0} |$ . All the  $H_{\text{lat}}$  become  $E_m$ :

$$\begin{aligned}
 (E - E_m) b_m &= - (E - E_m) \sum_n b_n \left( \sum'_R e^{ik \cdot R} \langle m, \vec{0} | n, \vec{R} \rangle \right) \\
 &+ \sum_n b_n \langle m, \vec{0} | \Delta U | n, \vec{0} \rangle \\
 &+ \sum_n b_n \left( \sum'_R e^{ik \cdot R} \langle m, \vec{0} | \Delta U | n, \vec{R} \rangle \right)
 \end{aligned}$$

} Small

III. Application to single s-level:

$$\begin{aligned}
 (E - E_s) &= - (E - E_s) \left( \sum'_R e^{ik \cdot R} \langle s, \vec{0} | s, \vec{R} \rangle \right) \leftarrow \alpha(R) \\
 &+ \langle s, \vec{0} | \Delta U | s, \vec{0} \rangle \leftarrow -\beta \\
 &+ \left( \sum'_R e^{ik \cdot R} \langle s, \vec{0} | \Delta U | s, \vec{R} \rangle \right) \leftarrow -\gamma(R)
 \end{aligned}$$

$$\Rightarrow E - E_s = - \frac{\beta + \sum'_R e^{ik \cdot R} \gamma(R)}{1 + \sum'_R e^{ik \cdot R} \alpha(R)} \approx -\beta - \sum'_R e^{ik \cdot R} \gamma(R)$$

Make the further approximation, that the sum is only over nearest neighbors:  $\gamma(R) \rightarrow \gamma$ .

$$E_s = E_s - \beta - \gamma \sum_{R \text{ n.n.}} e^{i\mathbf{k} \cdot \mathbf{R}}$$

A. Simple cubic lattice:  $\mathbf{R} = \pm a \hat{x}, \pm a \hat{y}, \pm a \hat{z}$

$$E = E_s - \beta - 3\gamma (\cos(k_x a) + \cos(k_y a) + \cos(k_z a))$$

For small  $k$  this is approximately

$$E = E_s - \beta - 6\gamma + \gamma a^2 (k_x^2 + k_y^2 + k_z^2).$$

B. BCC lattice:  $\mathbf{R} = \frac{a}{2} (\pm \hat{x} \pm \hat{y} \pm \hat{z})$

$$E = E_s - \beta - 8\gamma \cos\left(\frac{k_x a}{2}\right) \cos\left(\frac{k_y a}{2}\right) \cos\left(\frac{k_z a}{2}\right)$$

C. FCC lattice:  $\mathbf{R} = \frac{a}{2} (\pm \hat{x} \pm \hat{y}), \frac{a}{2} (\pm \hat{y} \pm \hat{z}), \frac{a}{2} (\pm \hat{z} \pm \hat{x})$

$$E = E_s - \beta - 4\gamma \left( \cos\left(\frac{k_x a}{2}\right) \cos\left(\frac{k_y a}{2}\right) + \cos\left(\frac{k_y a}{2}\right) \cos\left(\frac{k_z a}{2}\right) + \cos\left(\frac{k_z a}{2}\right) \cos\left(\frac{k_x a}{2}\right) \right)$$

Remarks on tight binding:

1. More atomic levels  
3x3 p  
5x5 d  
Plotted in direction of symmetry
2. Localized level  $\rightarrow$  narrow band  
Too wide  $\rightarrow$  T.B. not work
3. Traveling wave even though constructed of localized levels
4. Lattice w/ basis  $\rightarrow$  even only "s" bands have multiple bands (graphene example)
5. Spin-orbit coupling  $\rightarrow$  spin & spatial together
6. Indep. electron (discuss later)

Always of form  $\sum_R e^{ik \cdot R} \psi_n(r-R) = \psi_{nk}(r)$

IV. Wannier functions: Periodicity in k-space

$\psi_{n(k+k)}(r) = \psi_{nk}(r) \rightarrow \psi_{nk}(r) = \sum_R f_{n,r}(R) e^{iR \cdot k}$   
(periodic in k)

where  $f_{n,r}(R) = \frac{1}{V} \int_{1BZ} d^3k e^{-iR \cdot k} \psi_{nk}(r)$ .  $V = \text{volume in } 1BZ(k\text{-space})$

$f_{n(r+R_0)}(R+R_0) = \frac{1}{V} \int_{1BZ} d^3k e^{-i(R+R_0) \cdot k} \psi_{nk}(r+R_0)$   
 $= \frac{1}{V} \int_{1BZ} d^3k e^{-i(R+R_0) \cdot k} e^{ik \cdot R_0} \psi_{nk}(r)$   
 $= f_{n,r}(R)$

$\rightarrow f_{n,r}(R) = \phi_n(r-R)$

$\psi_{nk}(r) = \sum_R e^{ik \cdot R} \phi_n(r-R) //$