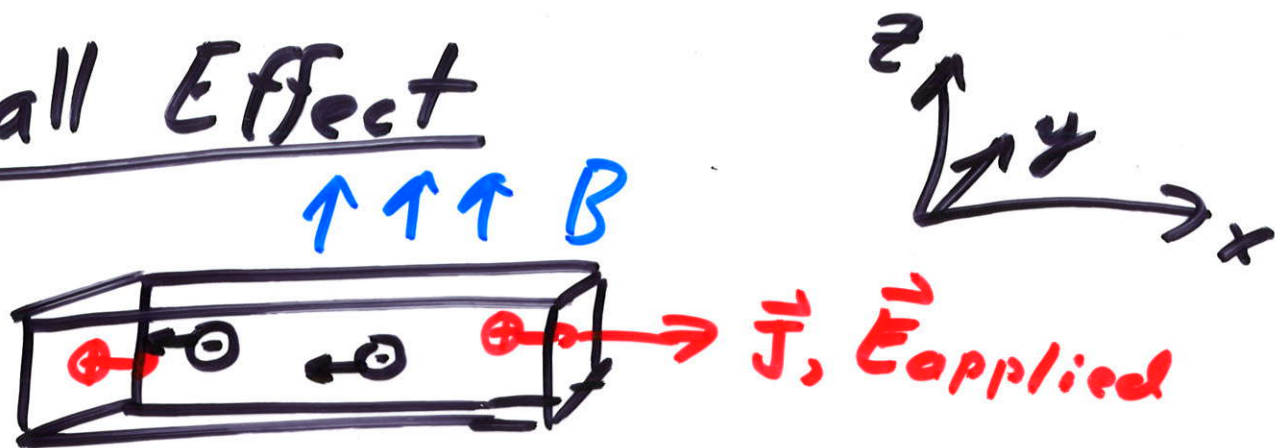


# Hall Effect



Ohm's Law  $\vec{J} = \sigma \vec{E}$ ,  $\sigma = \frac{n q^2 \tau}{m}$

Lorentz force  $\vec{F} = q(\vec{E} + \vec{v} \times \vec{B})$

+++ front  
--- back

$\vec{E}_{Hall}$  into page

--- front  
+++ back  
 $\vec{E}_{Hall}$  out of page

$\vec{F} \neq m\vec{a}$ ,  $\vec{F} = \frac{m\vec{v}}{\tau}$

x)  $\frac{m v_x}{\tau} = q(E_x + B_z v_y)$

y)  $\frac{m v_y}{\tau} = q(E_y - B_z v_x)$

z)  $\frac{m v_z}{\tau} = q E_z$

$E_{\text{Hall}} = E_y$  builds up until no more charge can flow in the  $y$ -direction

$\Rightarrow v_y = 0$  eventually

$$\Rightarrow E_y = B v_x = B \frac{q E_x \tau}{m}$$

Hall Coefficient

$$R_H \equiv \frac{E_y}{J_x B_z} = \frac{1}{nq}$$

$$\frac{R_H < 0}{\text{electrons}} \\ \hline R_H > 0 \\ \text{holes}$$

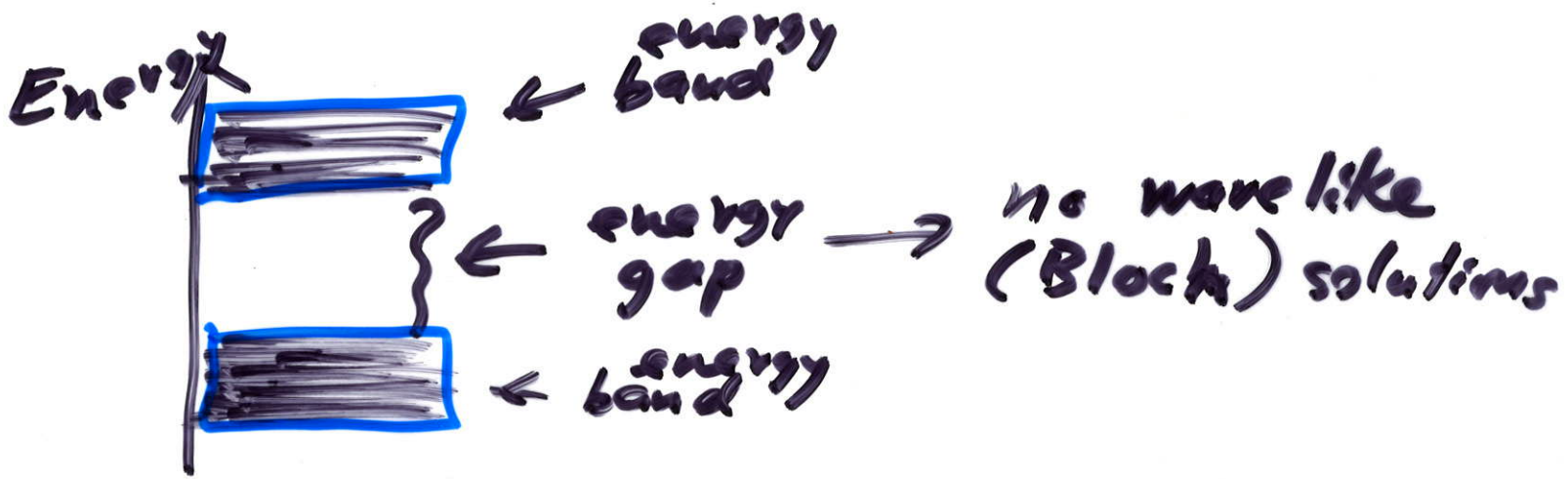
- Quantum Hall Effect
- fractional QHE

# Energy Bands

Last chapter - Free electron Fermi gas

Nearly free electron Fermi gas

Now electrons will interact (weakly) with the crystal lattice.

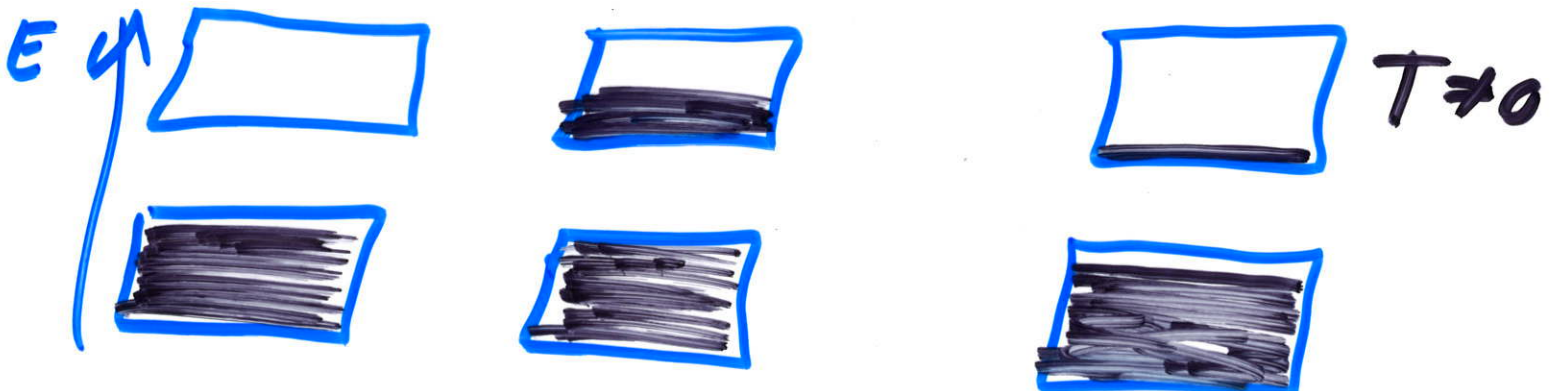


horizontal direction means nothing

Insulator

Metal

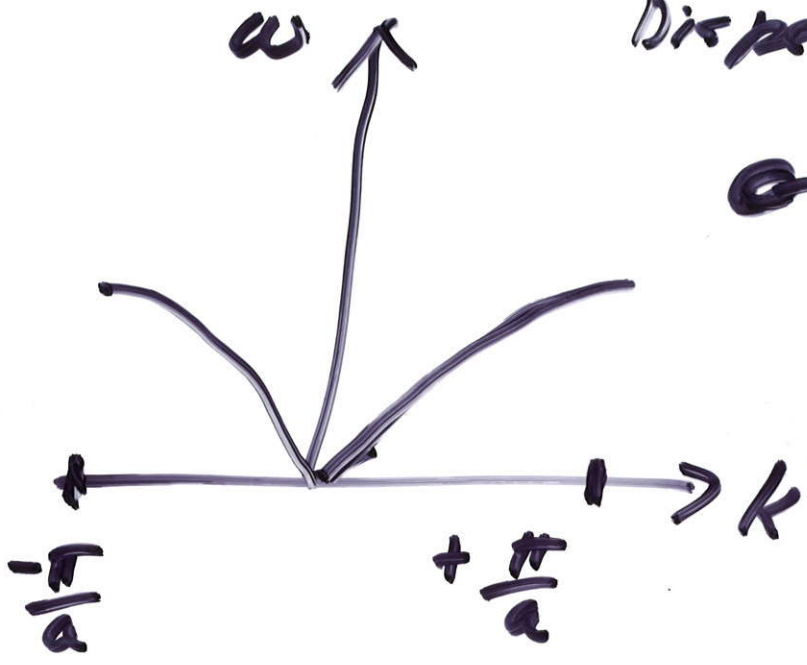
Semiconductor



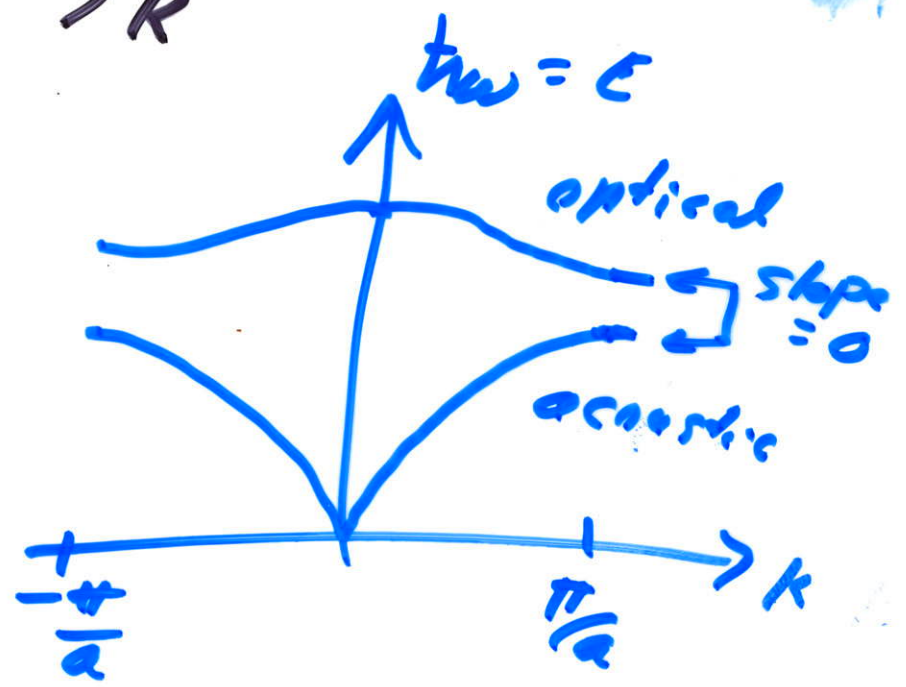
Have seen band gap before Phonons

Dispersion Relation

1-atom basis



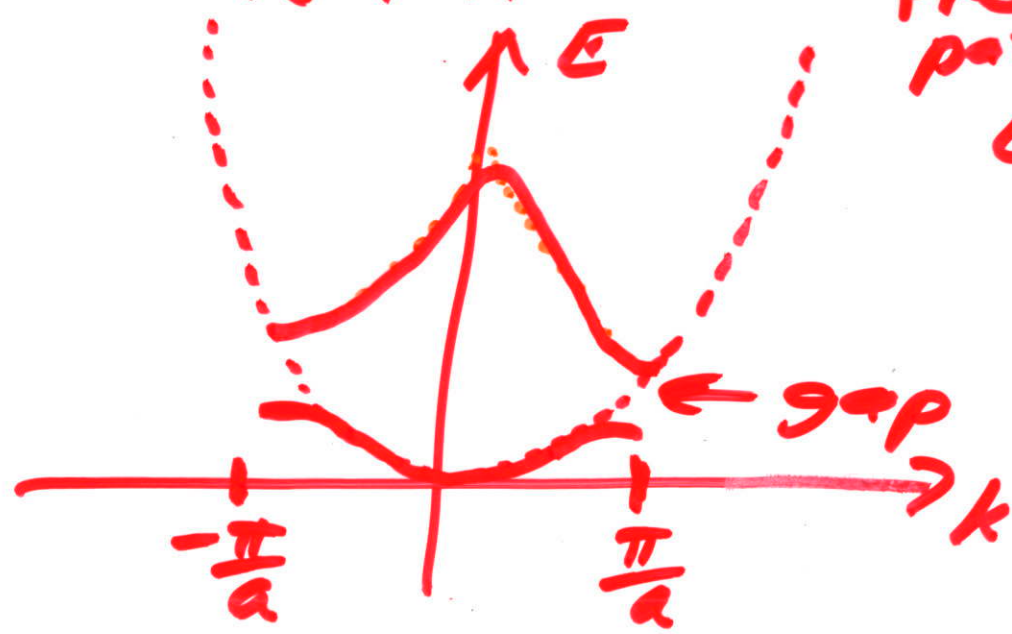
2-atom basis



Electrons

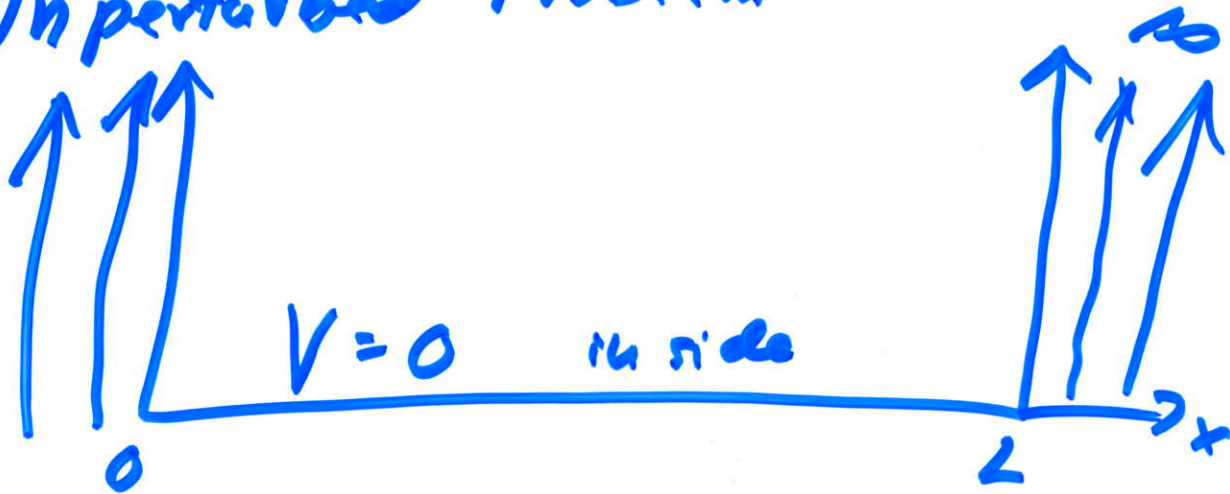
free electrons particle in a box

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

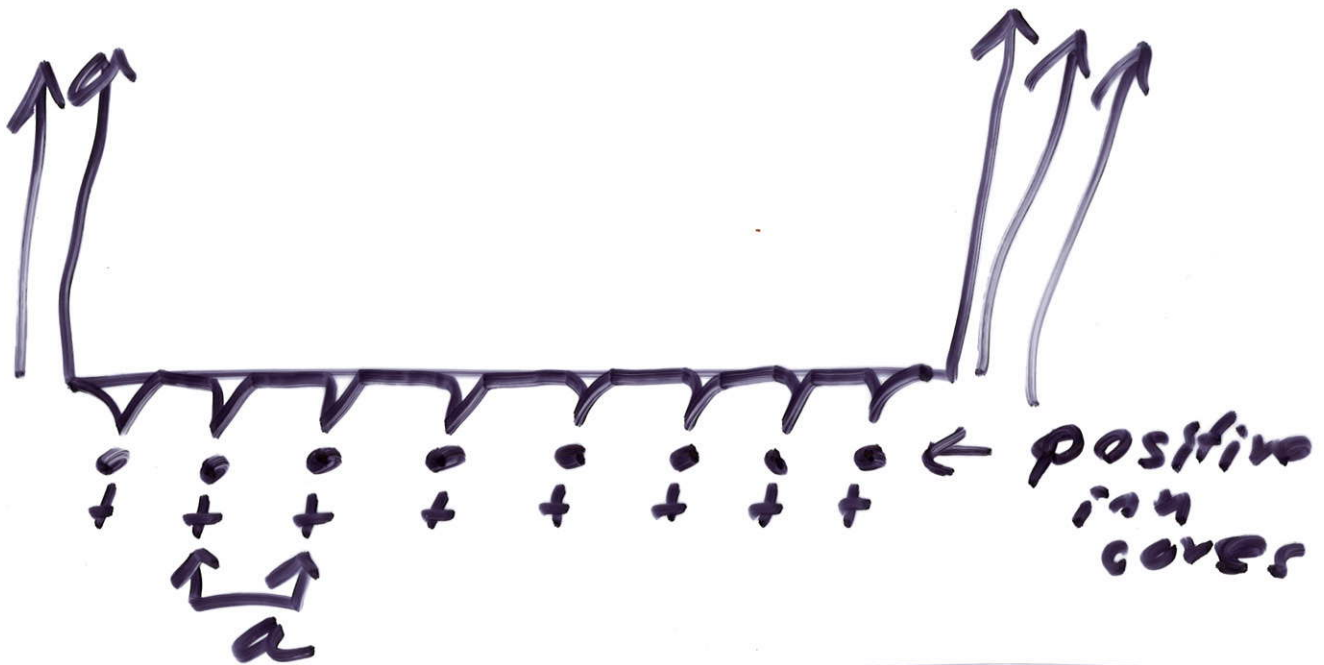




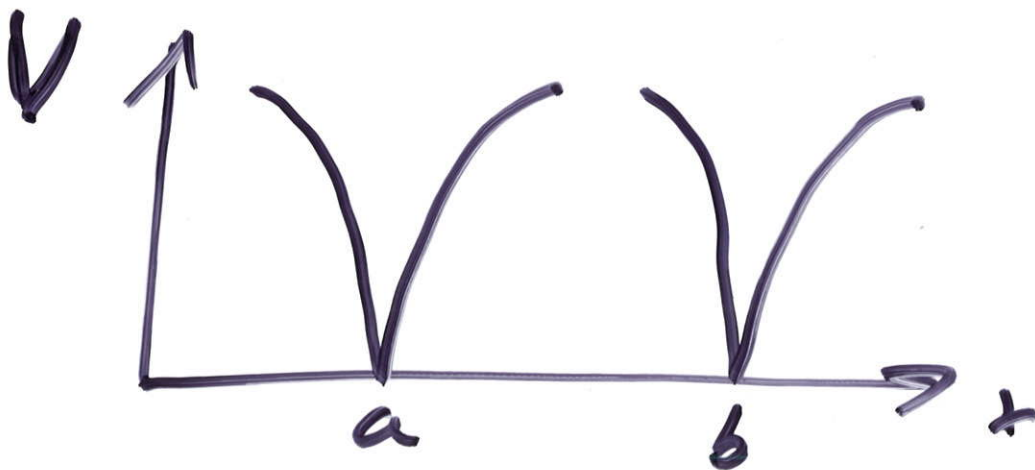
# Unperturbed Problem



# Perturbation



e.g. Two Hydrogen atoms  $H_2$  molecule



Electron wave function - overall must be antisymmetric

$$\Psi = \Psi_{\text{space}} \cdot \Psi_{\text{spin}} \quad \text{--- } 1/3.6 \text{ eV} \quad \left[ \begin{array}{l} \text{anti} \\ \text{binding} \\ \text{binding} \end{array} \right.$$

$$\Psi = \underbrace{\left\{ \Psi_a(\vec{r}_1) \Psi_b(\vec{r}_2) + \Psi_a(\vec{r}_2) \Psi_b(\vec{r}_1) \right\}}_{\text{space - symmetric}} \cdot \text{binding}$$

$$\frac{1}{\sqrt{2}} \left( \begin{array}{c} \uparrow \\ a \end{array} \begin{array}{c} \downarrow \\ b \end{array} - \begin{array}{c} \downarrow \\ a \end{array} \begin{array}{c} \uparrow \\ b \end{array} \right) \quad \begin{array}{l} S=0 \\ M_S=0 \end{array}$$

spin - antisymmetric  
singlet

$$\Psi = \left\{ \Psi_a(\vec{r}_1) \Psi_b(\vec{r}_2) - \Psi_a(\vec{r}_2) \Psi_b(\vec{r}_1) \right\} \cdot \text{anti-binding}$$

space - antisymmetric

• (triplet)

$$\begin{array}{c} \uparrow \uparrow \\ a \ b \\ S=1 \\ M_S=+1 \end{array}$$

$$\begin{array}{c} \downarrow \downarrow \\ a \ b \\ S=1 \\ M_S=-1 \end{array}$$

$$\frac{1}{\sqrt{2}} \left( \begin{array}{c} \uparrow \downarrow \\ a \ b \end{array} + \begin{array}{c} \downarrow \uparrow \\ a \ b \end{array} \right)$$


S=1  
M\_S=0

$N = 2$  Hydrogen atoms  
 $\Rightarrow$  2 energy levels

---

$N \sim N_A \Rightarrow N_A$  energy levels

 } bands

 } gaps

---

At the Brillouin zone boundaries  
 $\rightarrow$  standing waves due to  
Bragg reflection.

Traveling waves  $\psi \propto e^{\pm ikx}$

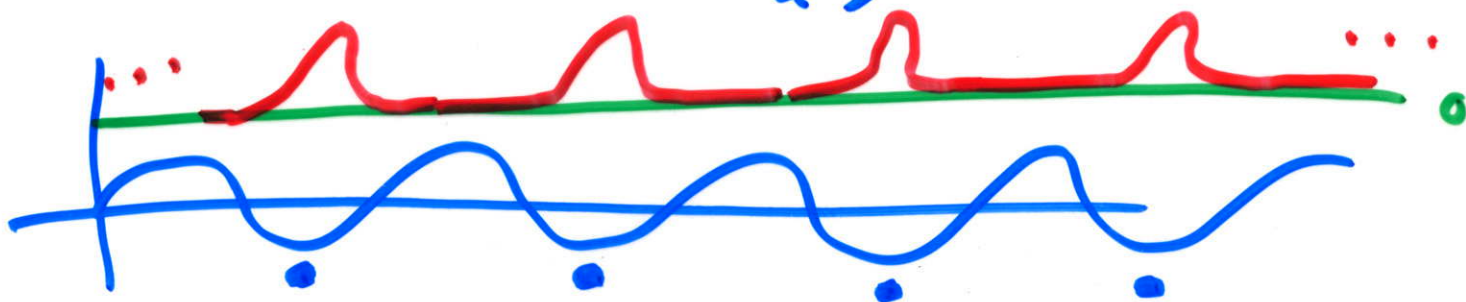
standing waves at B.Z.

$$\psi_+ \propto e^{ikx} + e^{-ikx} = 2 \cos(kx) = 2 \cos\left(\frac{\pi x}{a}\right)$$

$$\psi_- \propto e^{ikx} - e^{-ikx} = 2i \sin(kx) = 2i \sin\left(\frac{\pi x}{a}\right)$$



$$U(x) = U \cos\left(\frac{2\pi x}{a}\right)$$



(charge density)

$$\rho_+ = |\psi_+|^2 = \psi_+^* \psi_+ \propto \cos^2\left(\frac{\pi x}{a}\right)$$

$$\rho_- = |\psi_-|^2 \propto \sin^2\left(\frac{\pi x}{a}\right)$$

Traveling wave  $\psi \propto e^{+ikx}$

$$\rho \propto \psi^* \psi = e^{ikx} \cdot e^{-ikx} = 1$$

energy

$$E_+ = \frac{\int_0^L U(x) \rho_+(x) dx}{\int_0^L \rho_+(x) dx} = +\frac{U}{2}$$

$$E_- = \frac{\int_0^L U(x) \rho_-(x) dx}{\int_0^L \rho_-(x) dx} = -\frac{U}{2}$$