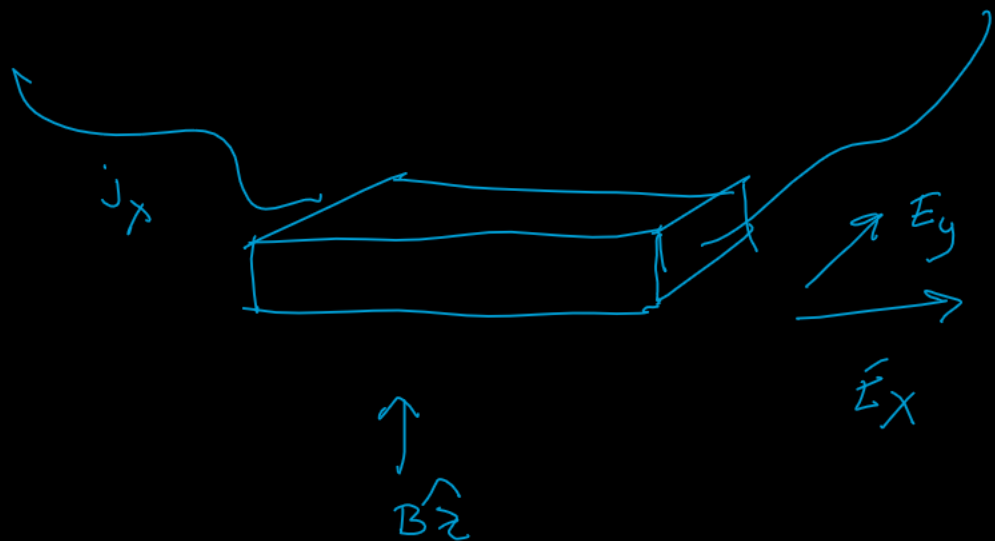
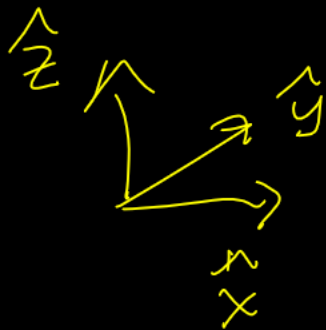


Lecture 5

Monday, January 25, 2021

HALL EFFECT



HALL FIELD IS E_y (COMES FROM \vec{B})

STEADY STATE:

$$V_x = -\frac{eE_x \tau}{m} - \omega_c \tau V_y$$

$$\omega_c = \frac{eB}{mc} \quad \text{CYCLOTRON FREQ.}$$

$$V_y = - \frac{e E_y \tau}{m} + \omega_c \tau V_x$$

$$V_y = 0 \text{ (NO CURRENT)}$$

$$\frac{+e E_y \tau}{m} = \omega_c \tau V_x = \frac{-e E_x \tau (\omega_c \tau)}{m}$$

HALL COEFFICIENT

$$R_H \equiv \frac{E_y}{j_x B} = - \frac{e \tau}{m c} \frac{E_x}{j_x} = - \frac{e \tau}{m c \sigma} = \frac{-e \tau}{m c} \left(\frac{m}{n e^2 \tau} \right)$$

$$R_H = \frac{-1}{n e c}$$

① MEASURES "SIGN" OF
CHARGE CARRIER

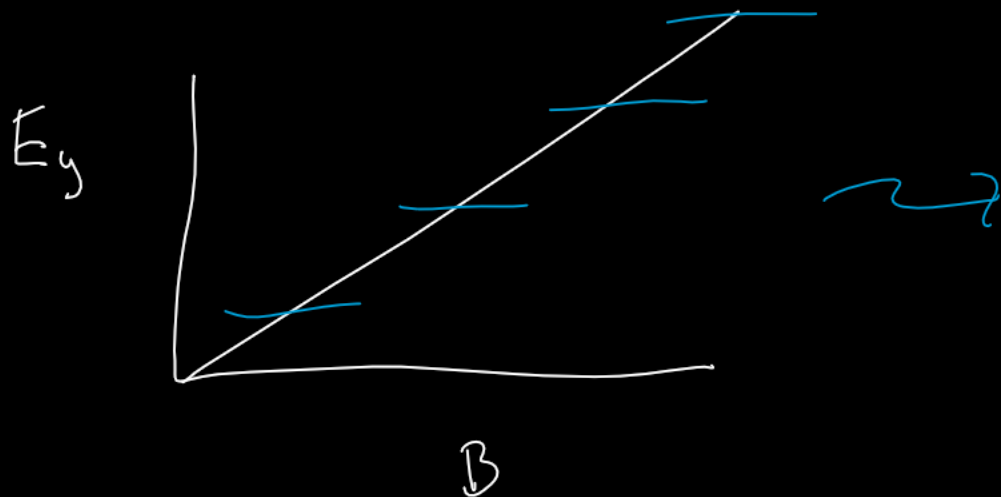
$$\left(\frac{1}{n(-e)c} \right)$$

...

$$\left(\frac{1}{n(-e)c} \right)$$

② DOES NOT DEPEND ON τ , m

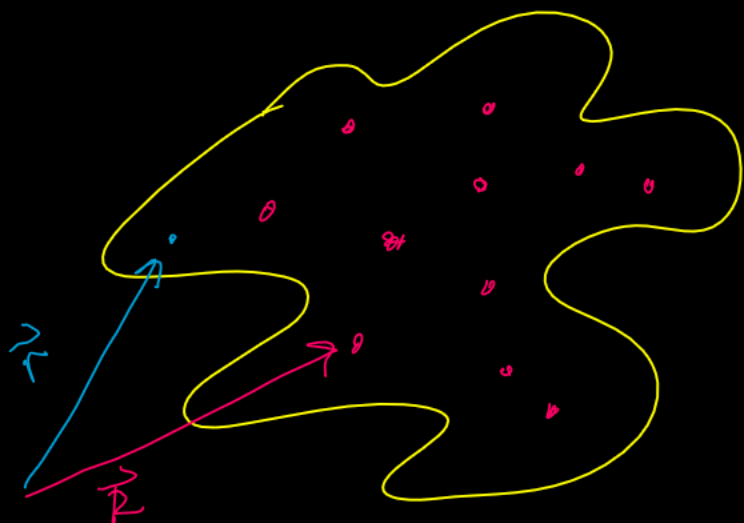
③ R_H MEASURES n WHICH COULD DEPEND ON T



LOW T , PURE
MATERIALS

"QUANTUM HALL EFFECT"

CHAP. 1



• NUCLEI + CORE ELECTRONS

$$U(\vec{r}) = \sum_{\{\vec{R}\}} \varphi(\vec{r} - \vec{R})$$

$$H = \sum_i H_i$$

↑
1-PARTICLE

$$H_i = \frac{-\hbar^2 \nabla_i^2}{2m} + \underbrace{\varphi(\vec{r}_i)}_{e^- \text{-ION POTENTIAL}}$$

$U(\vec{r})$

① LIQUID (1-CONFIGURATION)

② GLASSY (AMORPHOUS)

③ IONS ARE IN CRYSTAL (HIGH SYMMETRY)

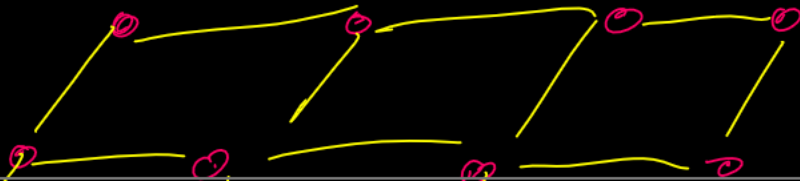
$$H(\vec{P}, \vec{r} + \vec{R}) = H(\vec{P}, \vec{r}) \quad \forall \vec{R} \ni \vec{R} = n_1 \vec{a}_1 + n_2 \vec{a}_2 + n_3 \vec{a}_3$$

$$n_1, n_2, n_3 \in \mathbb{Z}$$

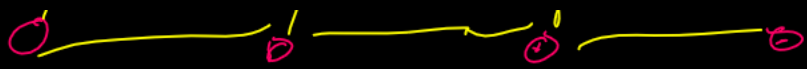
$\vec{a}_1, \vec{a}_2, \vec{a}_3$ ARE LINEAR
IND. VECTORS

$$\left\{ \vec{R} : \vec{R} = n_1 \vec{a}_1 + n_2 \vec{a}_2 + n_3 \vec{a}_3, n_1, n_2, n_3 \in \mathbb{Z} \right\}$$

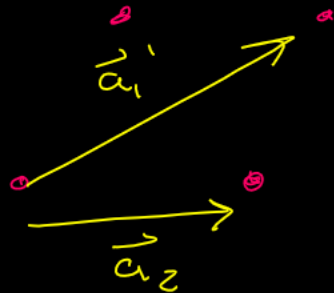
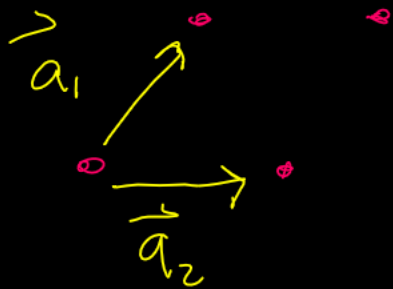
"LATTICE" (BRAVAIS LATTICE)



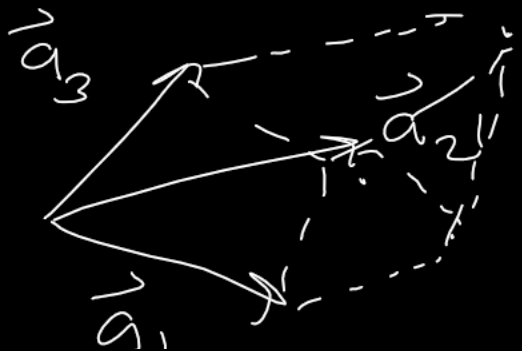
PERIODIC ARRANGEMENT OF POINTS
IN SPACE



$\{ a_i \}$ ARE NOT UNIQUE (PRIMITIVE VECTORS)



FILL ALL SPACE BY TRANSLATING A "PRIMITIVE" CELL

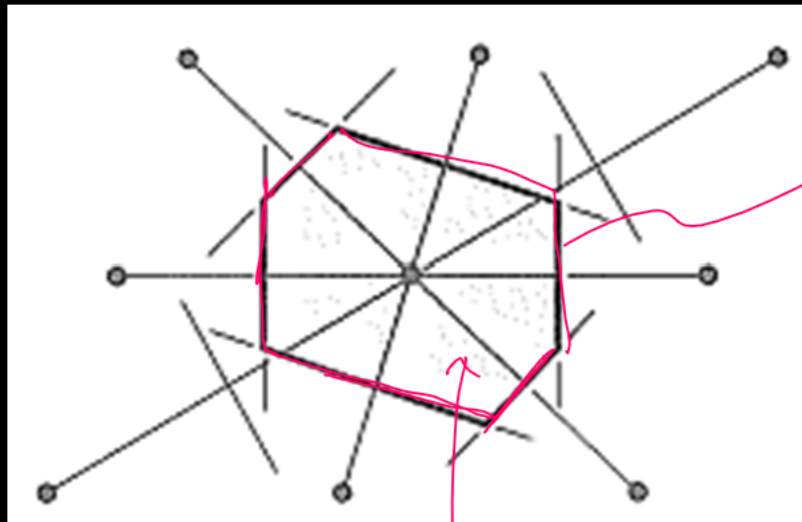


$$V_{||} = \vec{a}_1 \cdot \vec{a}_2 \times \vec{a}_3$$

4) -
PARALLELEPIPED

"SPECIAL UNIT CELL"

WIGNER-SEITZ CELL



UNIT CELL

L-BISECTOR

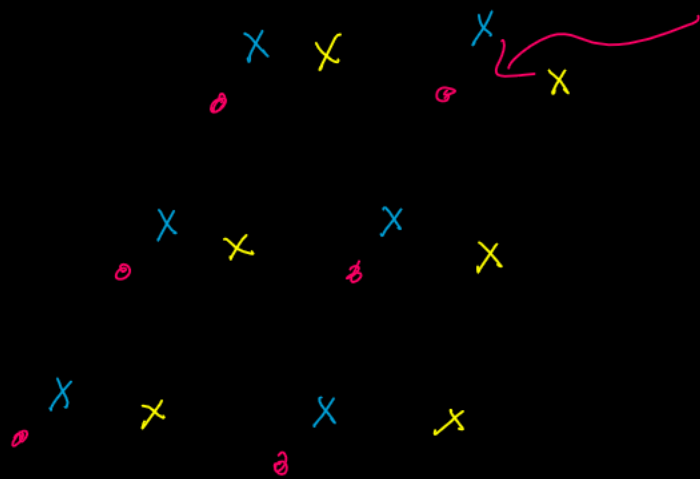
CRYSTAL = LATTICE + "BASIS"
STRUCTURE

BASIS = $\{ \vec{r}_i \}$

$$\vec{r}_j = x_j \vec{a}_1 + y_j \vec{a}_2 + z_j \vec{a}_3$$

BASIS $\approx \{ \vec{r}_i \}$
"ATOMS"

$$\vec{r}_i = x_i \vec{a}_1 + y_i \vec{a}_2 + z_i \vec{a}_3$$
$$0 \leq x_i, y_i, z_i < 1$$



LATTICE POINTS

$$\{ \text{red}, \text{blue} \}$$

2-D \Rightarrow 5 LATTICE TYPES

3-D \Rightarrow 14 LATTICE TYPES

CRYSTALS MORE COMPLICATED.