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DEVICES
EE 207

Some Devices →

BJT, MOSFET, Diode, JFET, Relay, Thyristor,
 UJFET, HEMT, Memory Devices

Digital CKts.: - 2 ~ 3 Giga Hz.

Purely based on mosfet

4G Speed → 5 - 10 GHz
 5G Speed → 5 - 20 GHz

e.g. neuromorphic Computing,
 10T Devices --

Midgap - 30%
 Endgap - 70%
 Quizzes - graded

SEMICONDUCTORS

Proper Covalent Bond

Group IV S.C. Si:Si {Equally Shared e⁻ b/w both atoms}

Isotype Semiconductor

Group III-V S.C. GaN ← An-isotype Semiconductor
 (distinct atoms) {e⁻ more drawn towards N}

Covalent Bond / Polar Bond

BRANAI LATTICE :- atoms arranged in a certain Config'

Eg → Simple Cubic, FCC (face centered cubic)

every corner

BCC

every corner + 1 @ every face

1 @ centre of cube

Basis → 1 silicon.

Note → Silicon is FCC + FCC

fused diagonally

= Diamond Cubic Structure

GaN = Hexagonally Closed Packed Structure (HCP)

$\text{Si} \rightarrow$ Crystal Structure \rightarrow Lattice (Diamond cubic) + Basis (Si atom)

↳ Driving force for formation of Crystal
↳ minimization of energy

Cast, Amorphous < Poly Cryst. < Crystal

ENERGY BAND DIAGRAM

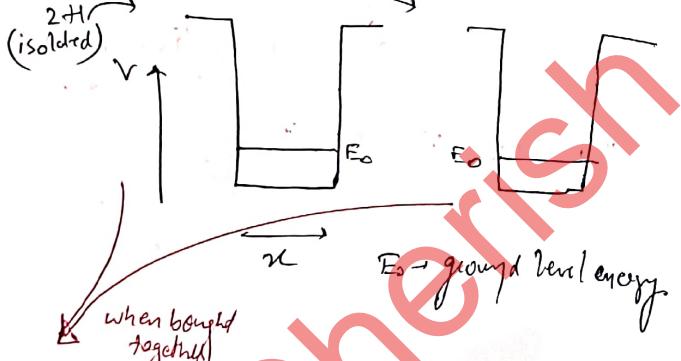
why is energy quantized?

$$\text{Duality Principle } (mv\lambda = n\lambda)$$

Waves \longleftrightarrow matter

Schrodinger Eqⁿ (hypothesis)

(+))



lattice of Energy levels

↳ Crystalline
Crystal ordering in micro.
Poly-Crystalline like grains
Amorphous - in
Chopped

Current-Voltage

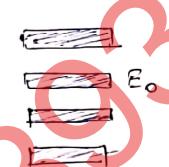
↓
flow of e⁻

e⁻ in a Crystal
Structure

in a Crystal Structure;

order of $10^{23} \sim 24$ atoms / cm³

as we bring more & more atoms, a band starts to form due to splitting

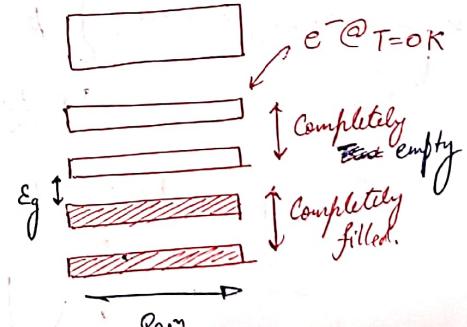


Energy Band Structure

(Pauli's principle)

if we magnify a band; we start seeing distinct energy levels; but separation is very small.

Si \rightarrow Unbalance e⁻



@ OK; all e⁻ are held by covalent bond, no. electrons (e⁻). available for conduction

$E_g \rightarrow$ Bandgap.

Semiconductor $\rightarrow E_g$ is small

Insulator $\rightarrow E_g$ is large ($E_g > 4eV$)

Metal $\rightarrow E_g = 4eV$ half filled band at 0K
Partially

In metal @ 0 Kelvin:

atom empty = conduction
mosi

Partially filled band @ 0 Kelvin
Topmost filled band = Valence Band

metal $\rightarrow e^-$'s are free.

When a e^- is moving in a metal cut freely
 $m_e = 9.11 \times 10^{-31} \text{ kg}$
 $F_{\text{int}} + F_{\text{ext}}$

$$p = \hbar k$$
$$m' \quad p = \hbar k$$
$$\frac{\hbar^2}{2m'} k^2$$

free e^- :

$$p = mv$$
$$E = \frac{1}{2} m v^2 \quad | \quad E = \frac{p^2}{2m_e}$$

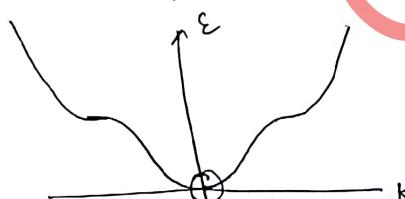
$$p = \hbar k \rightarrow \text{De Broglie wavelength } \left\{ p = \frac{h}{\lambda} \right\}$$
$$k = \frac{2\pi}{\lambda}$$

Dispersion Relation
 $a/k/a$

By (1) & (2);

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

$E-k$ relation in a Semiconductor



$$\text{Expanding } E \text{ about encircled pt; } \boxed{k_0=0}$$
$$E = E_0 + \frac{\partial E}{\partial k} k + \frac{1}{2} \frac{\partial^2 E}{\partial k^2} k^2 + \dots$$

Hence; for Semiconductors;

$$(\text{for S.C.}) \quad E = E_0 + \frac{1}{2} \frac{\partial^2 E}{\partial k^2} k^2$$

$$V_g = \frac{1}{\hbar} \cdot \frac{dE}{dk} \text{ of free}$$

$$(\text{for metal}) ; \quad E = \hbar^2 k^2 = \frac{\partial^2 E}{2m_e} k^2$$

$E-k$ curve parabolic
 $\therefore \frac{\partial^2 E}{\partial k^2} = \text{const}$

for silicon

$$m_{eff} = 0.2 m_e$$

$$m_{eff}/m^* = \hbar^2 \cdot \left(\frac{\partial^2 E}{\partial k^2} \right)^{-1}$$

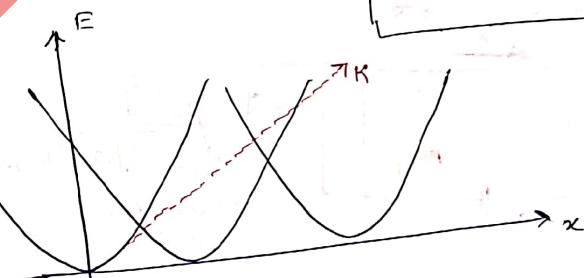
If m^* is less, inertia will be less

$$|E| = V/L$$

$$F = -q E$$

$$= -q \frac{V}{L} = m^* \frac{dV_L}{dt}$$

free on e^- in SC



Band gap E_c (Conduction Band Edge) $|$ Band gap vs pos.
 E_v (Valence Band Edge)

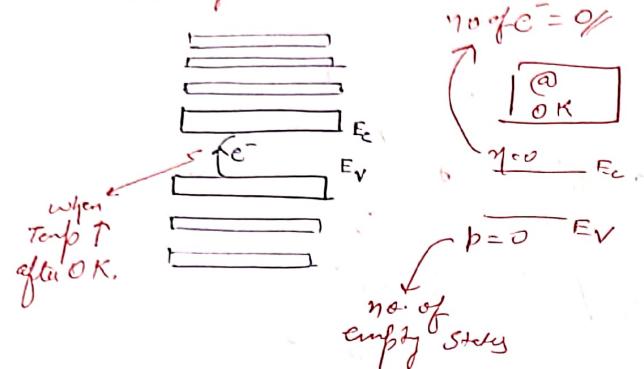
x.

E_c $|$ without bias;
 E_v

$$\text{Free } e^-$$
$$E = E_0 + \frac{\hbar^2 k^2}{2m_e} (\text{free } e^-)$$
$$PE \rightarrow \frac{\hbar^2 k^2}{2m_e} \rightarrow KE$$

$$\Delta^2 \frac{\partial E}{\partial k}$$

Concept of holes



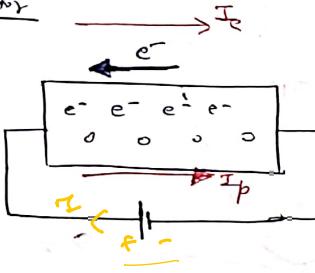
$T = 10 \text{ K}$

$e^{-} \rightarrow E_C \rightarrow 10^{11} / \text{cm}^3$

$\epsilon \downarrow E_V \rightarrow 10^{22} / \text{cm}^3$

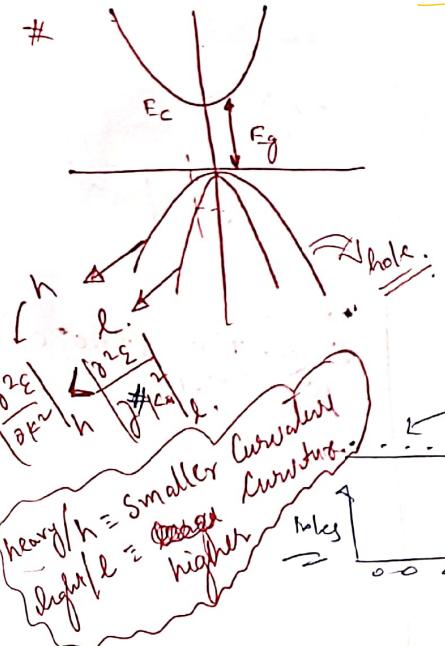
in principle every e^- below E_V also goes up to & above E_C but statistically; only e^- in E_V goes up to E_C

In Semiconductor

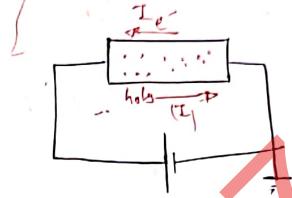


$I = I_c + I_p$

for e^- in valence band
in conduction band

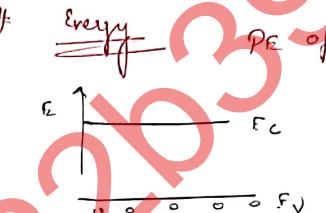


#



if u. charge Temp; E_C & E_V changes
Hence does the E_F change in meV/ $^{\circ}\text{C}$.

#



Energy PE of system \uparrow where e^- move from E_V to E_C
& hence minimization of energy is not driving force.

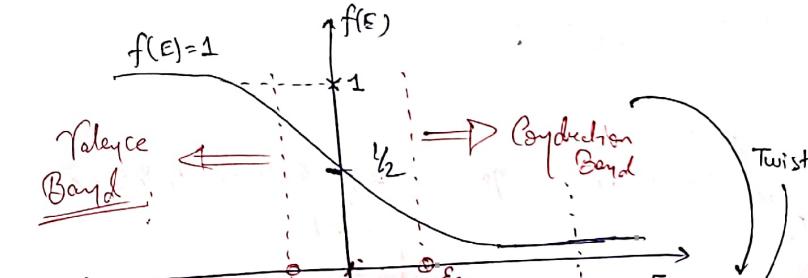
e^- jump from E_C to E_V for randomness (Entropy)

Probability of occupation of an energy level of E for any e^-

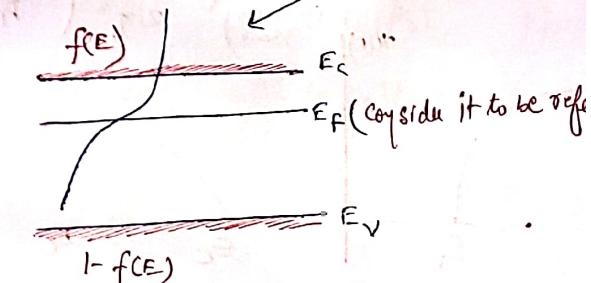
$$f(E) = \frac{1}{1 + e^{(E - E_F)/k_B T}}$$

Fermi-Dirac Statistics
 E_F = Fermi energy

E_F const.
 k_B



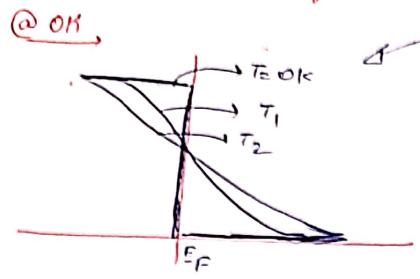
Prob. of occ. is high in vicinity of $E_F (\neq 0)$



$f(E)$ also tells den. of band ($f(E) \propto$)

...

Prob. of occup. of holes = $1 - f(E)$

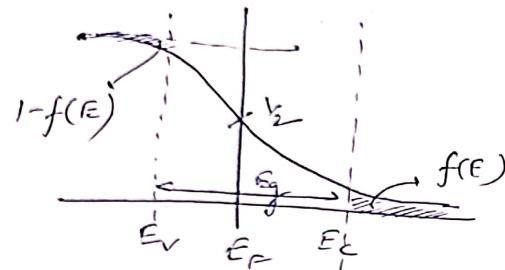


for fermions (e^-) {spin: $\pm \frac{1}{2}$ }

$$T_2 > T_1 > 0K$$

\Rightarrow E_F also changes with

Temp.,

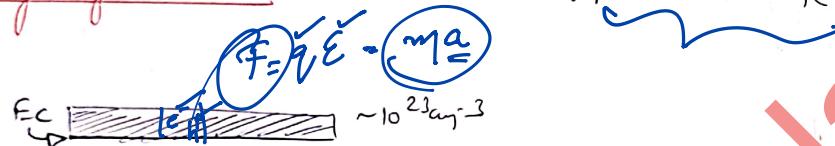


$$\mu_L = 1200$$

$$\mu_P = 500$$

$$\begin{aligned} N_C &= 4.7 \times 10^{17} \\ N_V &= 7 \times 10^{18} \text{ cm}^{-3} \\ E_F &= 1.42 \text{ eV} \\ n_L &= 4.7 \times 10^{17} \exp\left(-\frac{(E_F - \mu_L)}{kT}\right) \\ n_P &= 4.7 \times 10^{17} \exp\left(-\frac{(E_F - \mu_P)}{kT}\right) \end{aligned}$$

Density of States \rightarrow unit = Fermi energy / per unit volume

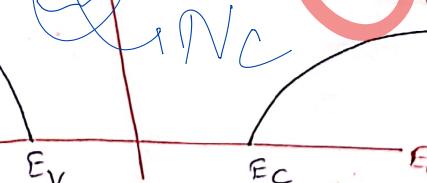


$$N_c = \text{density of energy levels} / \text{vol.}$$

$$N_c = \int_{E_C}^{E_V} N_c(E) dE = \eta \cdot \text{no. of energy levels by vol. } E_C \text{ to } E_V / \text{vol.}$$

$$N_c = \text{Density of Energy levels} / \text{vol.} \cdot \frac{1}{2\pi^2} \left(\frac{2m^*}{\hbar^2}\right)^{3/2} \sqrt{E - E_C}$$

$$f(E) \cdot N_c(E)$$



@ E_C ; density = 0

$$N_V = \frac{1}{2\pi^2} \left(\frac{2m_h^*}{\hbar^2}\right)^{3/2} \sqrt{E_V - E} \quad (m_e^* + m_h^*)$$

\Rightarrow Densities of States are independent of Temperature

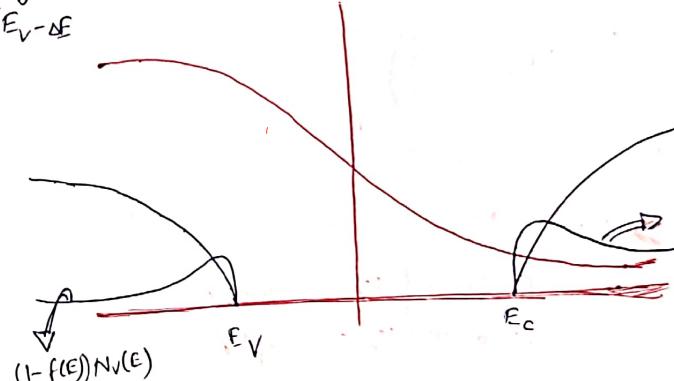
$$\int_{E_C}^{E_C + \Delta E} f(E) \cdot N_C(E) dE = \int dE = n$$

no. of levels in the " dE " interval

Prob. of occ. of that " dE " \times no. of levels

"no. of e^- "

$$\int_{E_V - \Delta E}^{E_V} (1 - f(E)) \cdot N_V(E) dE = p.$$



$$f(E) \cdot N_C(E)$$

$$E - E_F = 21eV$$

$$E - E_F = 72eV$$

$$\text{Boltzmann approximation} \approx \frac{1}{1 + e^{(E - E_F)/k_B T}} \cdot e^{-(E - E_F)/k_B T}$$

$$f(E) \approx$$

$$e^{-(E - E_F)/k_B T}$$

$$+ 1$$

$$f \uparrow$$

$f(\xi)$ after F is very big

$$\therefore f(E) \approx \frac{f_{c\rightarrow\infty}(\infty)}{E_c + \omega_r} e^{-j\omega_r t} f(E) N_c(t) dE$$

$$\eta = \left(\frac{n_c}{N} \right) \cdot e^{-(E_c - E_F)/k_B T}$$

$\text{Prefactor} = \text{Effective Density of States}$

$$n_c = \frac{m^* k T}{2\pi \hbar^2}^{3/2} \Rightarrow n_c \propto T^{3/2}$$

$$n = n_c \exp \left(- \frac{(E_c - E_f)}{kT} \right)$$

$$b = N_V' \exp \left(- \left(\frac{E_F - E_V}{kT} \right) \right)$$

if $(E - E_F) > 3k_B T$ Thermal Voltage

$$\frac{-(E-E_F)}{k_B T} \quad : \quad \begin{array}{l|l} T = 300K & k_B T = 26 meV \\ T = 600K & k_B T \approx 52 meV \end{array}$$

$$\frac{1}{1 + 20.08} \approx \frac{1}{20.08}$$

$$\text{Unit of } \frac{N_c(E)}{1} = \text{cm}^{-3} \text{ eV}^4$$

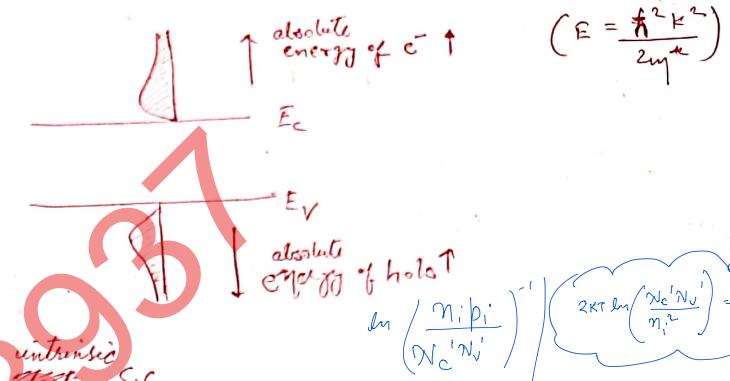
of allowed states per unit vol. per unit Energy

$$f(\varepsilon) \uparrow$$

4

$$\eta = \left(\frac{n_c}{n} \right) \cdot e^{-(E_C - E_F)/k_B T}$$

Δ
Prefactor = Effective Density of States



$$n_i = p_i$$

$$n_i p_i = n_c n_v \exp\left[-\frac{E_C - FV}{k_B T}\right]$$

$$b_i = n_i = \sqrt{n_c n_v} \exp \left[-\frac{E_g}{2k_b T} \right]$$

$$E_g \uparrow \Rightarrow n_i \downarrow$$

if $n_i \neq p_i$ (extrinsic Semiconductors)

$$n_F = N_C' N_V' \exp\left[-\frac{E_C - E_V}{k_B T}\right]$$

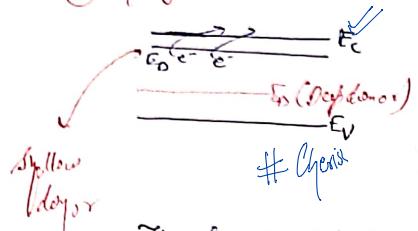
$$\left(\text{Either } n_i \text{ or } p_i \text{ is dominant} \right) \boxed{m_p = m_i^2} \rightarrow \begin{array}{l} \text{Law of mass action} \\ \text{always } (n_i = p_i \text{ or } n_i \neq p_i) \end{array}$$

In Intrinsic; $n_i = p_i$ { EHP } e-hole pair
 for Si; $n_i \approx 1.1 \times 10^{10} \text{ cm}^{-3}$

Extraipic \leftrightarrow n-type {e⁻} / Donor level \checkmark

p-type {h⁺} / Acceptor level

① n-type



The free e⁻ rotates in the crystal, for that e⁻;

$E_C - E_D$ = Ionization energy.

if $E_C - E_D$ is $\downarrow\downarrow$; e⁻ will find it very easy to jump to E_C to Conduction Band

K_BT = Reference

$E_C - E_D \approx K_B T \rightarrow$ Shallow donor

$E_C - E_D \ggg K_B T \rightarrow$ Deep donor $\xrightarrow{P, As}$

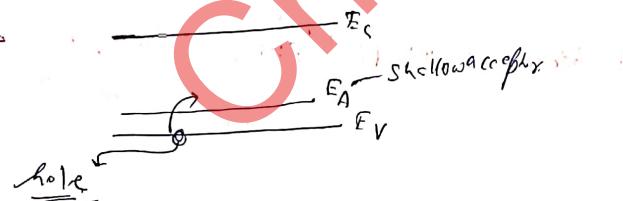
For Sulfur

Such type of e⁻ are K/a
Shallow Donors.

Donor becomes more charged in the process

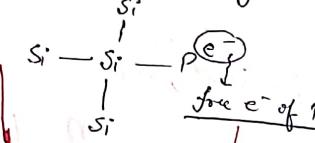
Note, In b/w E_C & E_V; One can only have the Donor level having e⁻. Rest no e⁻ may be b/w these 2 states

② p-type



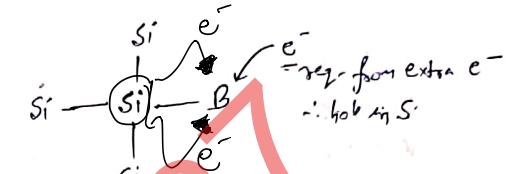
$$\frac{E}{P} \text{ Valency} = 4$$

Substitutional Impurity



Ionization Energy
is the energy for movement of e⁻ to move from Donor level

to Conduction Band



N_D = Donor density

N_A = Acceptor "

Charge neutrality

$$n + N_A = p + N_D$$

$$n_i = 1.1 \times 10^{10} \text{ cm}^{-3}$$

$$N_D \approx 10^3 \text{ cm}^{-3}$$

$$N_A \approx 10^{17} \text{ cm}^{-3}$$

$$p = \frac{n_i^2}{n} \quad n = \frac{n_i^2}{p}$$

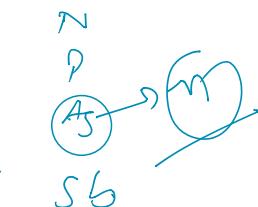
n



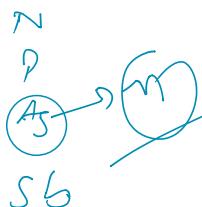
H



C



N



As

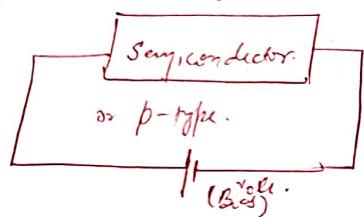
Sb

n-type

$n \gg p$
 $E_F - E_C$ is large
 $E_V - E_F$ is small.



n-type



The highest energy level that an electron can occupy at the absolute zero temperature is known as the Fermi Level. The Fermi level lies between the valence band and conduction band because at absolute zero temperature, the electrons are all in the lowest energy state

Cherish Jain

Extrinsic S.C.

$$n + N_A^- = p + N_D^+$$

$$n \approx N_D - N_A, \quad N_D - N_A \gg n_i$$

$$N_c = (\) \sqrt{E_F E_C}, \quad E_F > E_C$$

$$m_{\text{eff}} = 2m_B$$

$$E_C$$

$$E_F$$

$$E_V$$

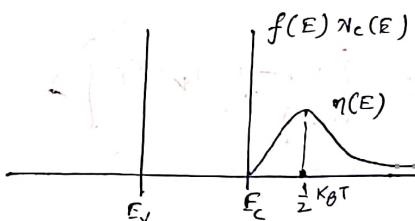
Similarly

$$\frac{E_C}{E_F} \quad \frac{E_F}{E_V}$$

\uparrow as E_F moves up

as E_F moves down; Boltzmann approximation becomes valid

$$\langle E \rangle = \left(\frac{\int E n(E) dE}{n} \right) = \text{avg energy of an } e^-$$



Avg energy of e^- = $\frac{3}{2} K_B T$

Energy / deg. of freedom, no. of deg. of freedom

Silicon

$$E_g = 1.1 \text{ eV}$$

$$N_c' = 10 \text{ cm}^{-3}$$

$$N_V' = 2 \times 10^{22} \text{ cm}^{-3}$$

$$n_i' = \sqrt{N_c' N_V'} \cdot \exp \left(- \frac{E_g}{2 K_B T} \right)$$

$$= 10^{22} \sqrt{2} \exp \left(- \frac{1.1}{2 \times 0.026} \right)$$

$$= 9.192 \times 10^{22} \text{ cm}^{-3}$$

Why ??

Why ??



$$\eta = N_c' \exp\left(\frac{E_F - E_C}{k_B T}\right)$$

$$E_F - E_C = k_B T \ln\left(\frac{n}{N_c'}\right)$$

$$= 0.026 \ln\left(\frac{9.19 \times 10^{12}}{10^{22}}\right) \quad \text{when } m_e = m_p \\ \epsilon_F = E_C + \epsilon_V$$

if S.C. is intrinsic ; $E_F = \frac{E_{FI}}{2}$

$N_c k$ if $n_i = p_i$

 $\epsilon_F = \frac{\epsilon_C + \epsilon_V}{2} \quad \left\{ \begin{array}{l} E_{FI} = E_C = E_V - E_{FI} \\ \text{Intrinsic Fermi level} \end{array} \right.$

$E_g = 1.1eV$
 $N_c' = 10^{22} \text{ cm}^{-3}$
 $N_V = 2 \times 10^{22} \text{ cm}^{-3}$
 $\eta_i = 9.19 \times 10^{12} \text{ cm}^{-3}$

$N_D = 10^{17} \text{ cm}^{-3}$

$N_A \approx$

$\approx 10^{12} \text{ cm}^{-3}$

$\eta \approx N_D - N_A = 10^{17} \gg \eta_i$

$\eta = N_c' \exp\left(\frac{E_F - E_C}{kT}\right)$

$E_F - E_C = k_B T \ln\left(\frac{n}{N_c'}\right) = 0.026 \ln\left(\frac{10^{17}}{10^{22}}\right) = -0.29$

Close to E_C

if $N_D = 10^{12}$

$E_F - E_C$ can't be done using Boltzmann factor

* * * if $N_D - N_A \approx N_c'$
then Boltzmann not valid

Effect of Temp on E_F

$N_c' = 10^{22} \text{ cm}^{-3} \quad T = 300K$

$N_c \sim T^{3/2}$

$\eta = 10^{17} \text{ cm}^{-3}$

$T = 400K$
 E_F
 ϵ_V

$n \uparrow \Leftarrow T \uparrow$

$\text{as } T \uparrow; \quad \epsilon_F \rightarrow \frac{\epsilon_C + \epsilon_V}{2}$

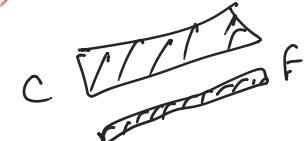
$\eta = N_c' \exp\left(-\left(\frac{\epsilon_C - \epsilon_F}{kT}\right)\right)$

$\eta = N_c' \cdot \exp\left(\frac{E_F - E_C}{k_B T}\right)$

$n_i = 10^{12}$
 $\eta = 10^{17}$ (for large T)

ρ_{FI} $\rho = \frac{n_i^2}{n} \cdot \frac{10^{24}}{10^{12}} = 10^7 \text{ cm}^{-3}$

$\eta + N_A = N_D + \rho^+$



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Transport

$$\left(\frac{J = I}{\text{Area}} \right)$$

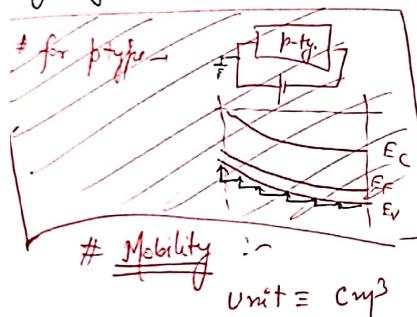
$$n(x) = n_c' \exp\left(\frac{E_F - E_C}{K_B T}\right)$$

if $n(x)$ is constant &
 n_c' is material const
then $E_F(x) - E_C(x)$ is also
constant

position of $e^- \rightarrow \text{random}$

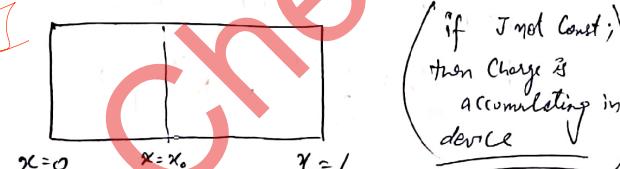
but net motion from
 $x=L$ to $x=0$ in a

ZigZag path.



$J(x) = \text{Constant}$, for const. Current in same surface area obj

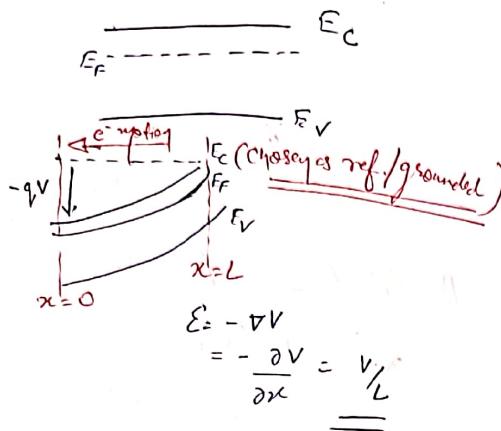
$$\frac{J = I}{A} \Rightarrow J = I$$



now ($A=1$) assume then

$$J = \frac{dQ}{dt} = qnV$$

Drift vel = v (avg speed of e^-)



$$\text{Now; } V = IR$$

$$R = \frac{rl}{A}; J = \frac{I}{A}$$

$$\text{Now; } V = JAx \cdot \frac{rl}{A}$$

$$\Rightarrow \frac{V}{L} = \frac{J}{\sigma}; \sigma = \frac{1}{l}$$

$$\Rightarrow J = \sigma E$$

another form of Ohm's law

$$\beta \approx \sigma E$$

$$\left(\frac{\text{drift Current}}{\text{area.}} \right) = \frac{q_n v}{\text{Area}} = J_{\text{drift}}$$

e^- in vol. in drift trar.
by e^- in 1 sec.

Order of mag
metals $\eta 10^{22} \text{ cm}^{-3}$
S-C $\eta 10^{17-19} \text{ cm}^{-3}$

Mobility

$$v = \mu E \quad \mu = \frac{v}{E} \quad (\text{cm}^2/\text{Volt}\cdot\text{sec.})$$

mobility

mobility Order

$$\mu = 0.5 - 1 \text{ cm}^2/\text{V}\cdot\text{sec} \Rightarrow \text{metal}$$

$$\mu = 500 \text{ cm}^2/\text{V}\cdot\text{sec} \Rightarrow \text{Graphene}$$

$$\mu = \frac{qz}{m^*}$$

$$\text{Now; } J = qnV \Rightarrow J = qn \mu E$$

$$\text{Now; } F = m^* a = qE = m^* \left(\frac{dv}{dt} \right) = m^* \frac{v}{\tau}$$

effective mass: $\langle v \rangle = \text{avg. velocity} = v_{\text{drift}}$
 $\tau = \text{scattering time}$ (Time b/w 2 collisions)

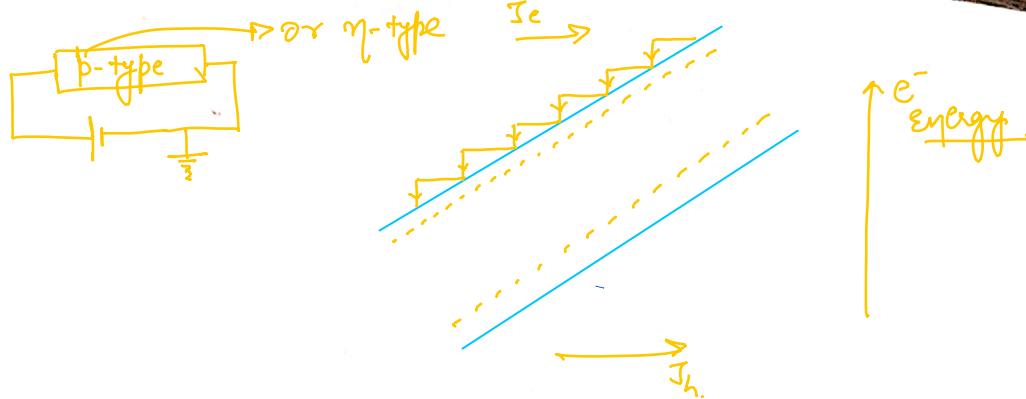
Hence; $\mu = \frac{q\tau}{m^*}$ $(\mu \propto \frac{1}{m^*})$

$V\tau = \text{dist b/w 2 collision}$

Ballistic Transport

all e^- travel without any scattering

TRANSISTER'S ($\Rightarrow \tau_{\text{long}}$) are semi-Ballistic
transport type



$$\begin{aligned} J_p &= q \mu_p \mu_p \epsilon \\ J_n &= q n \mu_n \epsilon \quad \left| \begin{array}{l} J_2 J_n + J_p \\ = q (\mu_p + \eta \mu_n) \epsilon \\ = \sigma \epsilon \end{array} \right. \end{aligned}$$

DIFFUSION

Driving force = $\frac{\partial n}{\partial x}$ → Conc gradient

Carrier Transport

Diffusion (due to EEF)

Drift (due to Conc diff.)

$$V = \mu \epsilon ; \mu = \frac{e \tau}{m^*}$$

$$J_n = q n \mu_n \epsilon$$

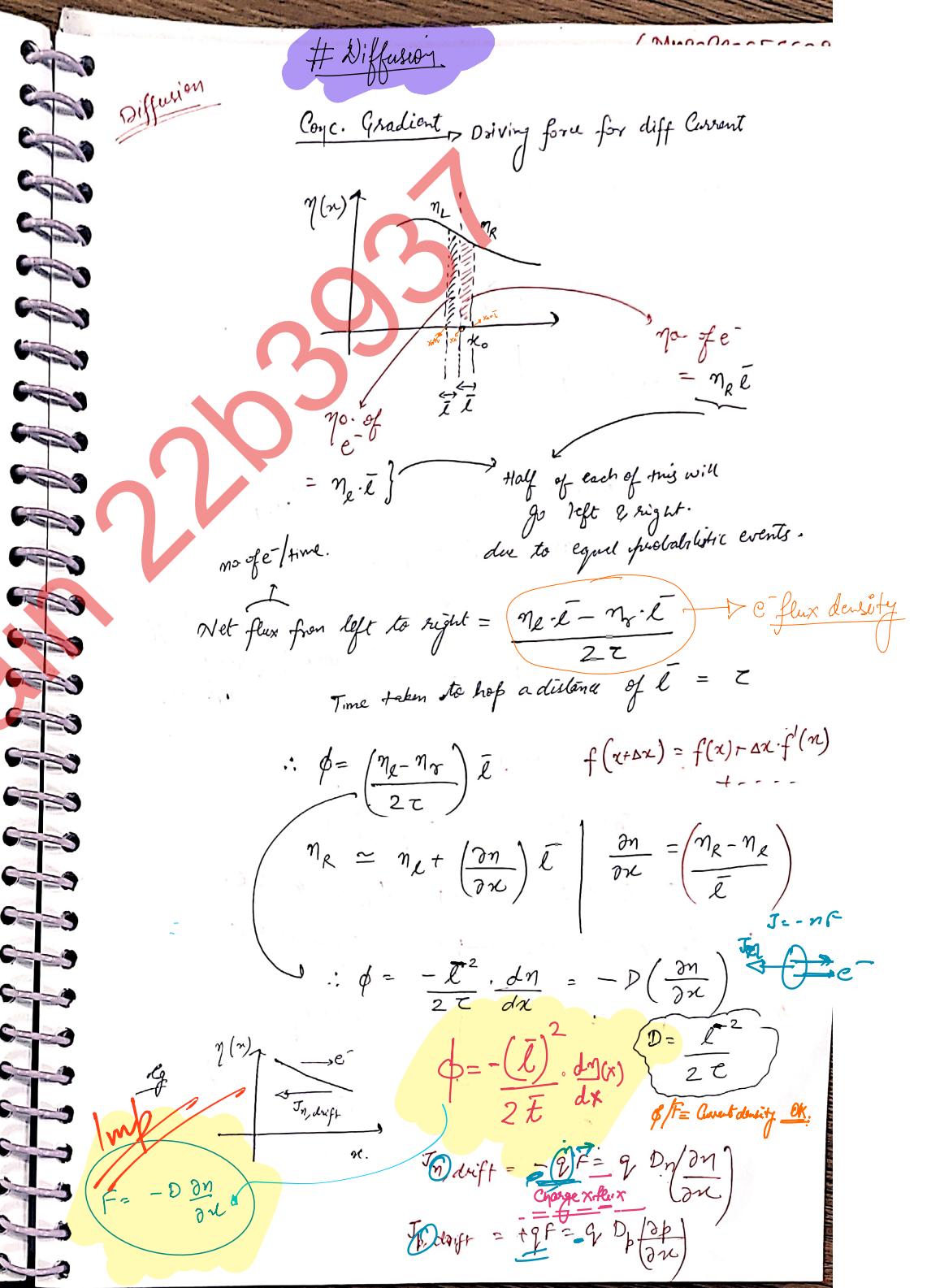
$$J_p = q \mu_p \mu_p \epsilon$$

$$J = J_n + J_p$$

we want $\uparrow \mu \Rightarrow \uparrow \text{Current}$

$\tau \uparrow$ $m^* \downarrow$

$\epsilon = V_L$



So now; accounting both diffusion & drift:-

$$J_n = q \mu_n E + q D_n \frac{\partial n}{\partial x}$$

$$J_p = q \mu_p E - q D_p \frac{\partial p}{\partial x}$$

$$\boxed{J_{\text{tot.}} = J_n + J_p}$$

$$\frac{D}{\mu} = \frac{k_B T}{q}$$

without any bias in a S-C.
assuming $J=0 + \infty$

$$\left. \begin{array}{l} J_n = 0 \\ J_p = 0 \end{array} \right\}$$

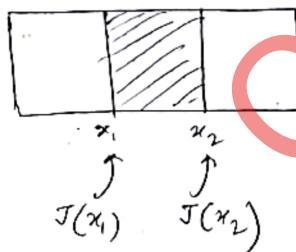
By applying Boltzmann's app.
Relation (Book)

By Einstein's relation on J_n

$$J_n = q \mu_n \left(\eta E + \frac{k_B T}{q} \frac{\partial n}{\partial x} \right)$$

$E_c - E_F < 2-3 k_B T \rightarrow$ Degenerate S-C. (Boltzian not valid)

$E_c - E_F > 2-3 k_B T \rightarrow$ non-degenerate S-C



$$\text{Steady State} \quad J(x_1) = J(x_2) \quad (\text{eqm})$$

But S-C devices
don't operate in
Condition :-

large for
n/p

Einstein's
Relation.

$$\frac{D}{\mu} = \frac{k_B T}{q}$$

only for systems
in which
 $E_v - E_F > 2k_B T$

$$J = 0$$

(eqm)

non-eq^m:

$$F(x_2) - F(x_1) = - \left(\frac{\partial \eta}{\partial t} (\eta \Delta x) \right)$$



Now; ~~$J_{\text{tot.}} = J_n + J_p$~~ $J = -q F$

$$\frac{J(x_2) - J(x_1)}{\Delta x} = q$$

$$\boxed{J_n = q \mu_n \left(\eta E + \frac{k_B T}{q} \frac{\partial n}{\partial x} \right)}$$

$$\boxed{J_p = q \mu_p \left(\eta E - \frac{k_B T}{q} \frac{\partial p}{\partial x} \right)}$$

$$\boxed{J_n = q \mu_n \left[\eta E + \frac{KT}{q} \frac{dn}{dx} \right]}$$

$$\boxed{J_p = q \mu_p \left[\eta E - \frac{KT}{q} \frac{dp}{dx} \right]}$$

$$\boxed{J_{\text{tot.}} = J_n + J_p}$$