

② Carrier densities: Complement

① Other expressions for n and p

For an intrinsic SC, we have $n = p = n_i$ and let us define $E_i = E_F$ the Fermi-level of an intrinsic SC.

For non-degenerate SC, it comes

$$\begin{cases} n_i = N_c e^{(E_i - E_c)/\beta} \\ p_i = N_v e^{(E_v - E_i)/\beta} \end{cases} \Rightarrow \begin{cases} N_c = n_i e^{(E_c - E_i)/\beta} \\ N_v = p_i e^{(E_i - E_v)/\beta} \end{cases}$$

we replace N_c, N_v in the expression for n and p. in general then

$$n = N_c e^{(E_F - E_c)/\beta} = n_i e^{(E_F - E_i)/\beta} \tag{2.19a}$$

$$p = p_i e^{(E_i - E_F)/\beta} \tag{2.19b}$$

Consequence \Rightarrow $n p = n_i^2$ (2.2e)

(b) How to calculate n_i ?

$$n_i = N_c N_v e^{-\frac{(E_c - E_v)\beta}{2}} = N_c N_v e^{-\frac{E_g \beta}{2}}$$

\Rightarrow [2.21] $n_i = \sqrt{N_c N_v} e^{-\frac{\beta E_g}{2}}$ $\beta = \frac{1}{k_B T}$

E_g is the energy bandgap.

$\Rightarrow n_i$ depends on the temperature (see Fig [2.20])

(c) Doped SC (Extrinsic SC)

if the SC is doped by both donors and acceptors, the material is however charge neutral

[2.24] $n + N_A^- = p + N_D^+$

$N_A^- \equiv N_A$
 $N_D^+ \equiv N_D$ } assume total ionization of dopant atoms.

N_A = density of acceptors
 N_D = " " donors-

Let us find a new expression for n and p in function of N_A, N_D, n_i .

$$p = \frac{n_i^2}{n} \Rightarrow \frac{n_i^2}{n} + N_D - n - N_A = 0.$$

$$n^2 - n(N_D - N_A) - n_i^2 = 0 \quad [\text{quadratic equation}]$$

$$n = \frac{N_D - N_A}{2} + \sqrt{\left(\frac{N_D - N_A}{2}\right)^2 + n_i^2} \quad [2.29a]$$

$$p = \frac{n_i^2}{n} = \frac{N_A - N_D}{2} + \sqrt{\left(\frac{N_A - N_D}{2}\right)^2 + n_i^2} \quad [2.29b]$$

(d) N-type and P-type SC

- { N-type SC has more free e^- than h^+
- { P-type SC has more free h^+ than e^-

a sc containing donor atoms in majority is called N-type sc.

a sc containing acceptor atoms in majority is called P-type sc.

N-type

P-type

$N_A \approx 0$

$N_D \approx 0$

$n \ll p$

$p \gg n$

$n \approx N_D$
$p \approx \frac{n_i^2}{N_D}$

$p \approx N_A$
$n \approx \frac{n_i^2}{N_A}$

[2.30]

[2.31]

(e) where is E_c ?

$n = p$ for intrinsic semiconductor.

$$\Rightarrow N_c e^{\beta(E_i - E_c)} = N_v e^{\beta(E_v - E_i)}$$

Solving for E_i .

$$E_i = \frac{E_c + E_v}{2} + \frac{k_B T}{2} \ln\left(\frac{N_v}{N_c}\right)$$

however $\frac{N_v}{N_c} = \left(\frac{m_p^*}{m_n^*}\right)^{3/2}$ [2.35]

at cones \Rightarrow $E_i = \frac{E_c + E_v}{2} + \frac{3}{4} k_B T \ln\left(\frac{m_p^*}{m_n^*}\right)$ [2.36]

if $T=0$ or $m_p^* = m_n^* \Rightarrow E_i$ is exactly in the middle of the bandgap.

(in ~~the~~ general case it will be very close to the middle).

Q) where is E_F ?

for extrinsic semiconductors ($E_F \neq E_i$)
from the expression of n and p we get.

[2.37]

$$E_F - E_i = k_B T \ln\left(\frac{n}{n_i}\right) = -k_B T \ln\left(\frac{p}{n_i}\right)$$

for N-type SC $n \approx N_d$

$$E_F - E_i = k_B T \ln\left(\frac{N_d}{n_i}\right) \quad [2.38a]$$

for P-type SC

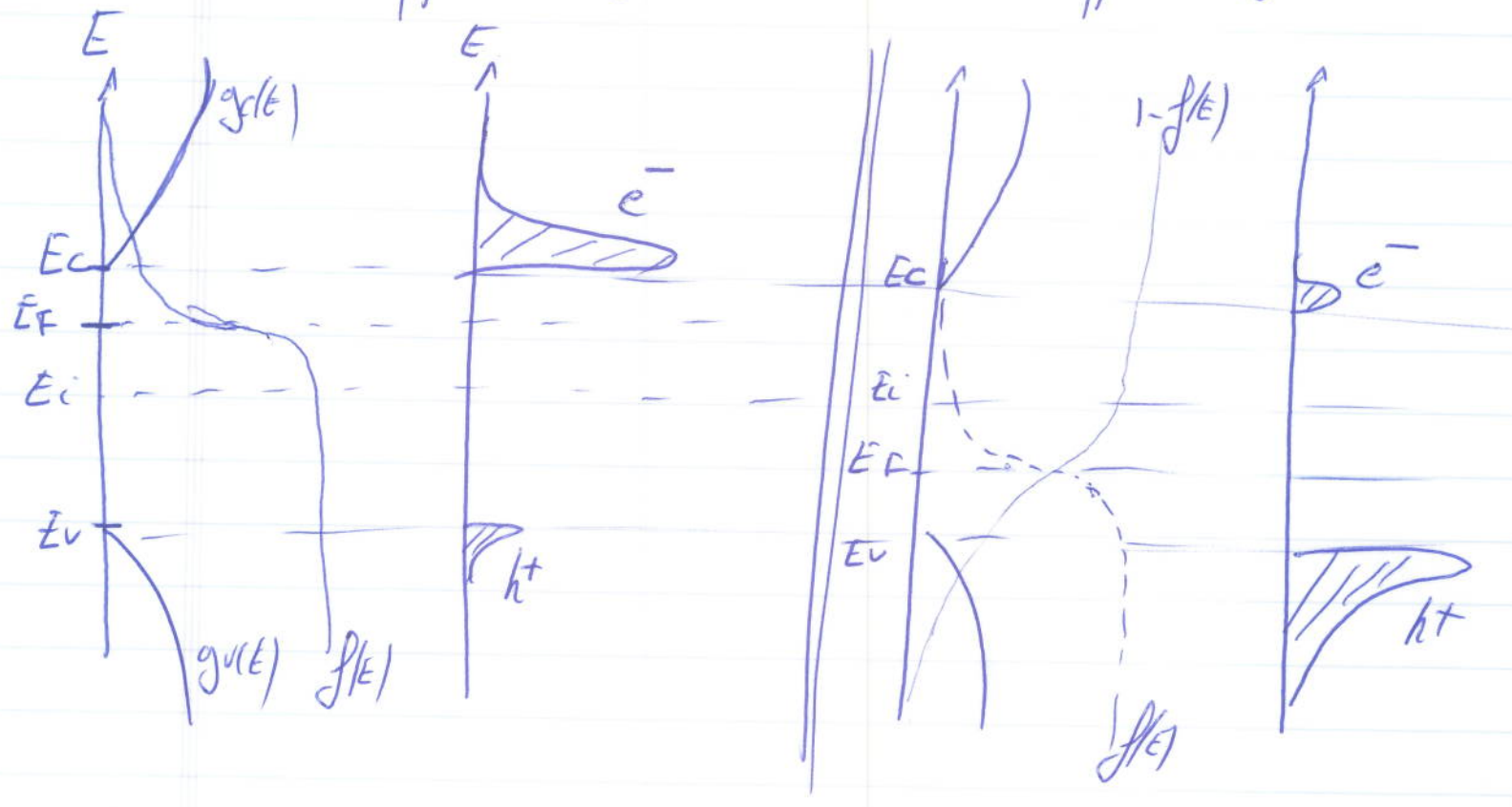
$p \approx N_A$ $E_F - E_i = -k_B T \ln\left(\frac{N_A}{n_i}\right)$ [2.38b]

So $E_F \neq E_i \Rightarrow$ the Fermi-level adjusts itself to ensure charge neutrality in the doped SC.

see Fig 2-21

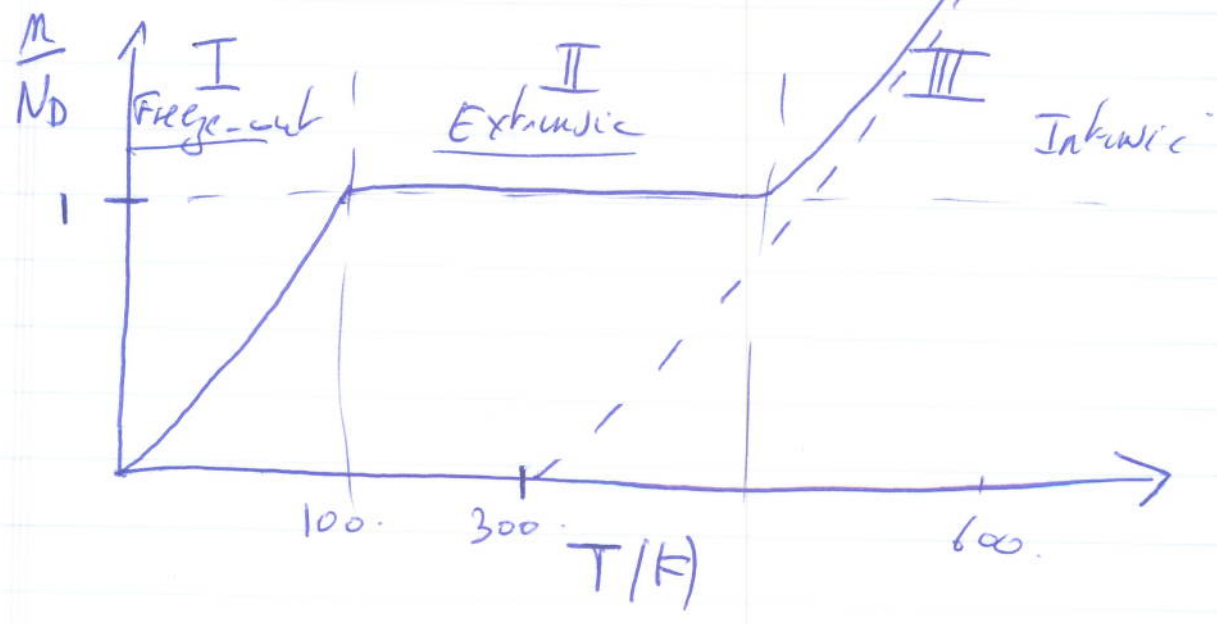
N-type Fig 2.16a

P-type Fig 2.16c

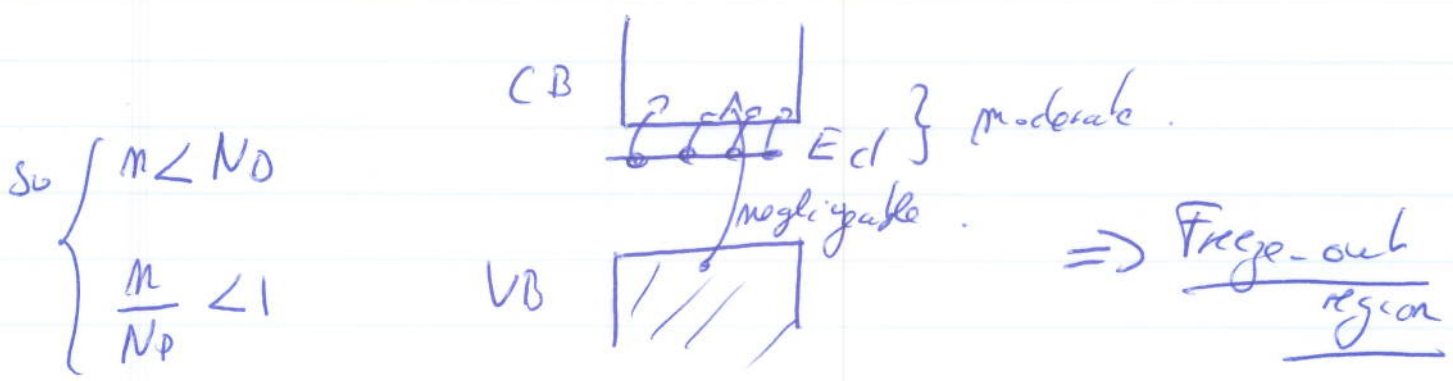


[257] (g) temperature dependence

Suppose a N-type SC Fig 2.22



Region I; at low Temperature, the thermal energy is not enough to ionize all the donor impurities. Some e^- are frozen at the donor energy level.



Region II at moderate T ; the condition of complete ionization is reached $n = N_D$ $\frac{n}{N_D} = 1$.

The number of e^- going from VB to CB is still negligible as compared to N_D . \Rightarrow Extrinsic region.

Region III if T is very large, the intrinsic carrier concentration n_i (e^- going from VB \rightarrow CB) is comparable to N_D and beyond this point SC becomes intrinsic \Rightarrow intrinsic region

$$\begin{array}{l} n > N_D \\ \frac{n}{N_D} > 1 \end{array}$$

(h) Band-bending

[Under equilibrium conditions, the Fermi-level E_F inside a structure is constant and unique.

$\Rightarrow E_F$ appears as a horizontal line on equilibrium band diagrams.

However E_c, E_v may vary with position \Rightarrow band bending

Example Non uniformly ~~doped~~ doped SC

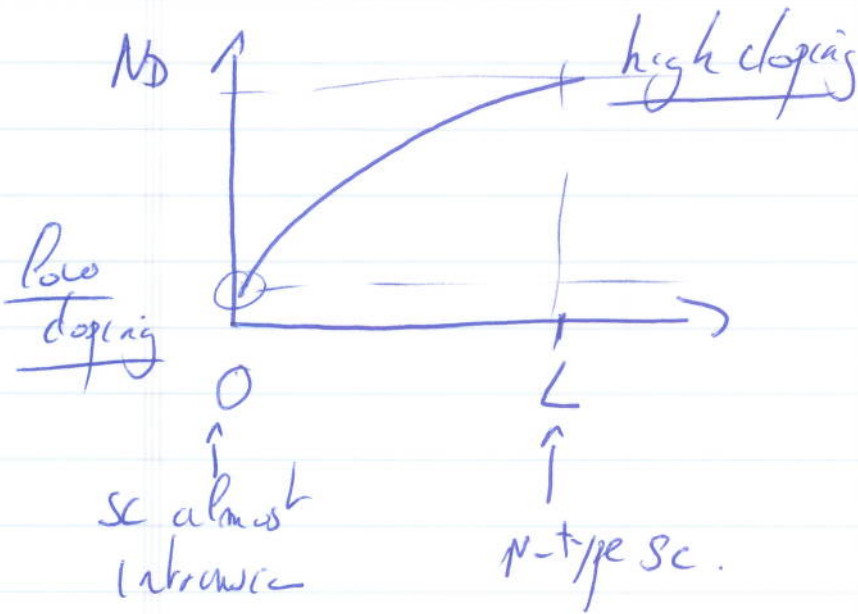


Fig 3.14a

energy band diagram (Recipe)

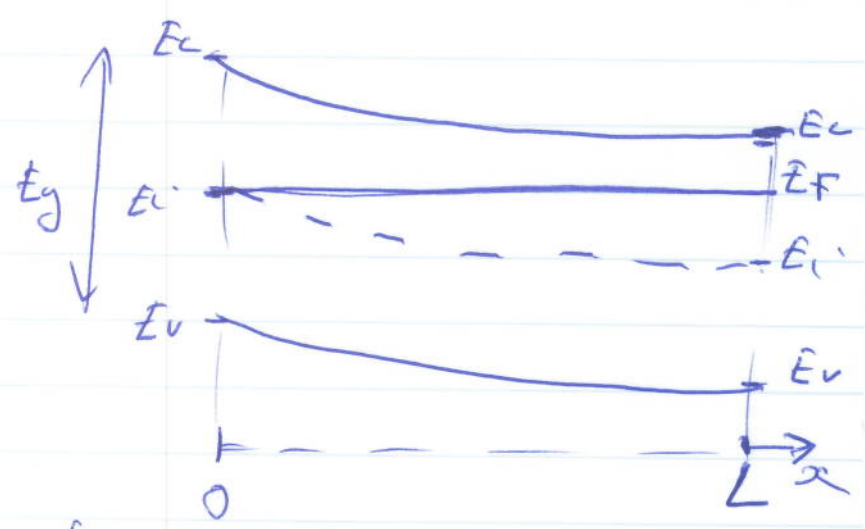
① plot E_F

② if $x \ll L$, SC is almost intrinsic so $E_F \approx E_i$ in the middle of bandgap

③ if $x \gg 0$, SC is N-type, so $E_F > E_i$

④ E_g is constant.

Fig 3.14b



⇒ In equilibrium situation, E_F is constant and unique, $E_c(x); E_v(x); E_i(x)$ may vary with position x .