

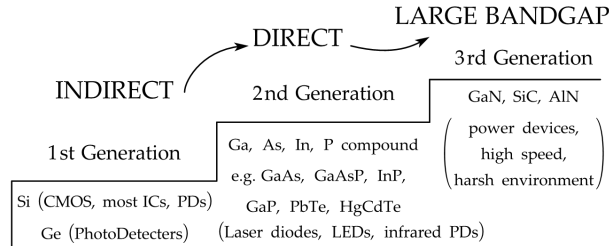
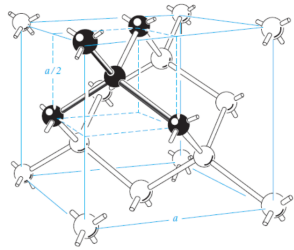
Semiconductor Midterm CheatSheet (chapter1-4)

	$e = 1.602 \times 10^{-19} \text{C}$
Boltzmann's constant	$k = 8.62 \times 10^{-5} \text{eV/K} = 1.38 \times 10^{-23} \text{J/K}$
ppm/ppb/ppt	partpermillion/billion/trillion

Si $E_g(\text{Si}) = 1.1 \text{eV}$ $5 \times 10^{22} \text{atoms/cm}^3$ $n_i \approx 1.5 \times 10^{10} \text{cm}^{-3}$ $\mu_n = 1350 \text{cm}^2/\text{V-s}, \mu_p = 480 \text{cm}^2/\text{V-s}$	Ge $E_g(\text{Ge}) = .67 \text{eV}$ $n_i = 2.4 \times 10^{13} \text{cm}^{-3}$ $\mu_n = 3900 \text{cm}^2/\text{V-s}, \mu_p = 1900 \text{cm}^2/\text{V-s}$
--	---

three types of solids: Crystalline(silicon wafer), Polycrystalline(Poly-Si), Amorphous(glass)
Primitive Cell, Unit Cell, Miller Indices (*hkl*) Diamond Lattice, Zinc Blende Structure,

Wurtzite Structure. packing fraction = $\frac{V_{\text{atom}} \cdot N_{\text{atom}}}{V_{\text{cell}}} = \frac{\frac{4}{3}\pi(\frac{d_{\text{atom,min}}}{2})^3 \cdot N_{\text{atom}}}{(\text{lattice constant})^3}$



Czochralski Growth (CZ) for Si:

$$k_d = \frac{C_s}{C_l} = \frac{[\text{impurity in solid}]}{[\text{impurity in liquid}]}, \text{MASS} = \frac{[\text{number of atoms}]}{(\text{Avogadro's constant})} (\text{atomic mass})$$

Liquid Encapsulated Czochralski Growth (LEC), Bridgman Growth (BR) for GaAs

Atom structure

Heisenberg Uncertainty Principle: $(\Delta x)(\Delta p_x) \geq \hbar/2$, $(\Delta E)(\Delta t) \geq \hbar/2$ The probability sum of finding a particle over the entire space $\int_{-\infty}^{\infty} \Psi^* \Psi dx dy dz = 1$. Average value $\langle Q \rangle = \int_{-\infty}^{\infty} \Psi^* Q_{\text{op}} \Psi dx dy dz$, where Q_{op} is defined as,

classical variable	quantum operator
$x, f(x)$	$x, f(x)$
momentum $p(x)$	$\frac{\hbar}{j} \frac{\partial}{\partial x}$
energy E	$-\frac{\hbar}{j} \frac{\partial}{\partial t}$

Schrödinger equation $E_k(\frac{p^2}{2m}) + E_p(V) = E_{\text{total}}(E)$, by replacing p and E with operators, we can derive $-\frac{\hbar^2}{2m} \frac{\partial^2 \psi(x)}{\partial x^2} \phi(t) + V(x) \psi(x) \phi(t) = -\frac{\hbar}{j} \psi(x) \frac{\partial \phi(t)}{\partial t}$.

Quantum Numbers for $\Psi(r, \theta, \phi) = R_n(r) \Theta_l(\theta) \Phi_m(\phi)$		
n	principal quantum number, shell	$1, 2, \dots$
l	angular momentum quantum number, subshell	$0, 1, \dots, n-1$ ($s, p, d, f, g, \dots$)
m	magnetic quantum number, orientation of orbitals (within the subshells)	$0, \pm 1, \dots, \pm(n-1)$
s	electron spin	$\pm \frac{1}{2}$

Pauli exclusion principle No two electrons in an interacting system with the same n, l, m, s . The only two electrons with same n, l, m must have opposite spins

Atomic number (Z)	Element	Number of electrons										Shorthand notation
		1s	2s 2p	3s 3p	3d	4s	4p	4d	4f	5s	5p	
1	H	1										1s ¹
2	He	2										1s ²
3	Li		1							1s ² 2s ¹		
4	Be		2							1s ² 2s ²		
5	B		2	1						1s ² 2s ² 2p ¹		
6	C		2	2						1s ² 2s ² 2p ²		
7	N		2	3						1s ² 2s ² 2p ³		
8	O		2	4						1s ² 2s ² 2p ⁴		
9	F		2	5						1s ² 2s ² 2p ⁵		
10	Ne		2	6						1s ² 2s ² 2p ⁶		
11	Na				1					[Ne] 3s ¹		
12	Mg				2					3s ²		
13	Al				2	1				3s ² 3p ¹		
14	Si				2	2				3s ² 3p ²		
15	P				2	3				3s ² 3p ³		
16	S				2	4				3s ² 3p ⁴		
17	Cl				2	5				3s ² 3p ⁵		
18	Ar				2	6				3s ² 3p ⁶		
19	K						1			[Ar] 4s ¹		
20	Ca						2			4s ²		
21	Sc				1					3d ¹ 4s ²		
22	Ti				2					3d ² 4s ²		
23	V				3					3d ³ 4s ²		
24	Cr				5	1				3d ⁵ 4s ¹		
25	Mn				5	2				3d ⁵ 4s ²		
26	Fe				6					3d ⁶ 4s ²		
27	Co				7					3d ⁷ 4s ²		
28	Ni				8					3d ⁸ 4s ²		
29	Cu				10	1				3d ¹⁰ 4s ¹		
30	Zn				10	2				3d ¹⁰ 4s ²		
31	Ga				10	2	1			3d ¹⁰ 4s ² 4p ¹		
32	Ge				10	2	2			3d ¹⁰ 4s ² 4p ²		
33	As				10	2	3			3d ¹⁰ 4s ² 4p ³		
34	Se				10	2	4			3d ¹⁰ 4s ² 4p ⁴		
35	Br				10	2	5			3d ¹⁰ 4s ² 4p ⁵		
36	Kr				10	2	6			3d ¹⁰ 4s ² 4p ⁶		

n	l	m	s/h	Allowable states in subshell	Allowable states in complete shell
1	0	0	$\pm \frac{1}{2}$	2	2
2	0	0	$\pm \frac{1}{2}$	2	8
	1	-1	$\pm \frac{1}{2}$	6	
		0	$\pm \frac{1}{2}$		
		1	$\pm \frac{1}{2}$		
3	0	0	$\pm \frac{1}{2}$	2	18
	1	-1	$\pm \frac{1}{2}$	6	
		0	$\pm \frac{1}{2}$		
		1	$\pm \frac{1}{2}$		
	2	-2	$\pm \frac{1}{2}$	10	
		-1	$\pm \frac{1}{2}$		
		0	$\pm \frac{1}{2}$		
1		$\pm \frac{1}{2}$			
	2	$\pm \frac{1}{2}$			

Silicon: $1s^2 2s^2 2p^6 3s^2 3p^2$

Energy Band and Carriers

Octet Rule, LCAO(antisymmetric/anti-bonding, parallel spin, high energy; symmetric/bonding); hybridization (Pauli exclusion principle, split energy levels $\rightarrow$ continuous energy bands) **E-k diagram**: parabolic $E = \frac{\hbar^2}{2m^*} k^2 + E_c$, $k = \frac{2\pi}{\lambda}$

Energy Levels and Concentrations

carrier concentrations = (Fermi distribution)(available states density)

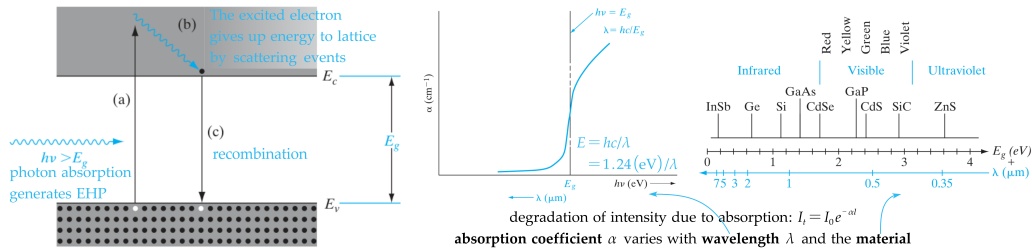
$$n_0 = \int_{E_c}^{\infty} f(E) N(E) dE = N_c f(E_c) \approx N_c e^{-(E_c - E_F)/kT}$$

$$p_0 = \int_0^{E_v} (1 - f(E)) N(E) dE = N_v (1 - f(E_v)) \approx N_v e^{-(E_F - E_v)/kT}$$

$$\text{intrinsic} \begin{cases} n_i = \sqrt{N_c N_v} e^{-E_g/2kT} \\ n_0 p_0 = n_i^2 \end{cases} \quad \text{doped, equilibrium} \begin{cases} n_0 = n_i e^{(E_F - E_i)/kT} \\ p_0 = n_i e^{(E_i - E_F)/kT} \end{cases}$$

$$\text{excess carriers, quasi-Fermi level} \begin{cases} n = n_i e^{(F_n - E_i)/kT} \\ p = n_i e^{(E_i - F_p)/kT} \end{cases} \quad (\text{for majority carriers } F_n/F_p \approx E_i)$$

Excess Carriers



Fluorescence (direct recombination) most energy directly into photon, fast(10^{-10} s)

Phosphorescence (indirect recombination) needs extra (deep) energy level between E_g , gives smaller $h\nu$, slow(s); application: UV excites phosphors to generate white light instead of smooth spectrum

equilibrium no stimulate

steady state can have stimulation inputs, stabilises after a long time

Recombination and Generation Balance

excess carrier concentrations $\delta n(t) = \delta p(t)$, initial values $\Delta n = \Delta p$.

low-level injection (linear)	n-type	p-type
$\frac{d\delta n(t)}{dt} = -\alpha_r(n_0 + p_0)\delta n(t)$	$\delta n(t) = \Delta n e^{-t/\tau}$	
minority carrier lifetime $\tau_n = \frac{1}{\alpha_r(n_0 + p_0)}$	$\tau_n = (\alpha_r n_0)^{-1}$	$\tau_p = (\alpha_r p_0)^{-1}$
$g_{op} = \alpha_r[(n_0 + p_0)\delta n(t) + \delta n^2(t)]$	$\delta n = \delta p = g_{op} \tau_n$	

photoconductive decay:

$$\sigma(t) = q[n(t)\mu_n + p(t)\mu_p]$$

$$\Delta\sigma = qg_{op}(r_n\mu_n + r_p\mu_p) = q(\delta n\mu_n + \delta p\mu_p)$$

Direct Recombination:

net change = thermal generation – recombination

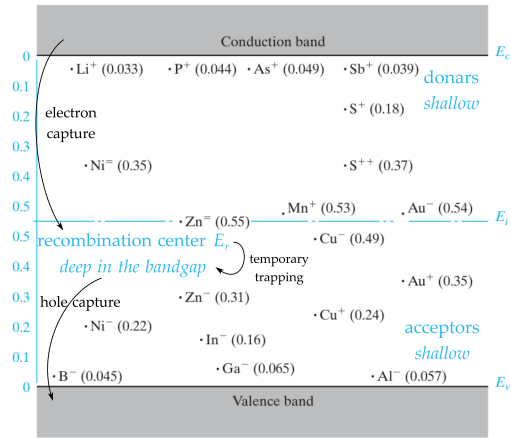
$$\frac{dn(t)}{dt} = \alpha_r n_i^2 - \alpha_r n(t)p(t)$$
$$= -\alpha_r[(n_0 + p_0)\delta n(t) + \delta n^2(t)]$$

steady state EHP G–R balance:

generation = recombination

[equilibrium] $G(T) = \alpha_r n_0 p_0 = \alpha_r n_i^2$

[steady state] $G(T) + g_{op} = \alpha_r n_0 p_0 + \alpha_r[(n_0 + p_0)\delta n(t) + \delta n^2(t)]$



Drift and Diffusion

Trade – off: high $N_a, N_d \Rightarrow$ high σ and low μ_n, μ_p

(at equilibrium)

mean free time $\bar{t}$

mobility μ_n, μ_p

drift current density J_x

conductivity σ

electric force accelerates $\frac{dp_x}{dt} = -nq\mathcal{E}_x$

$J_x = q(n\mu_n + p\mu_p)\mathcal{E}_x = \sigma\mathcal{E}_x$

$R = \frac{L}{wt}\rho = \frac{L}{wt}\frac{1}{\sigma}$

collision decelerates $\frac{dp_x}{dt} = -\frac{p_x}{\bar{t}}$

