

Theory of Semiconductor Devices

(반도체 소자 이론)

Lecture 16

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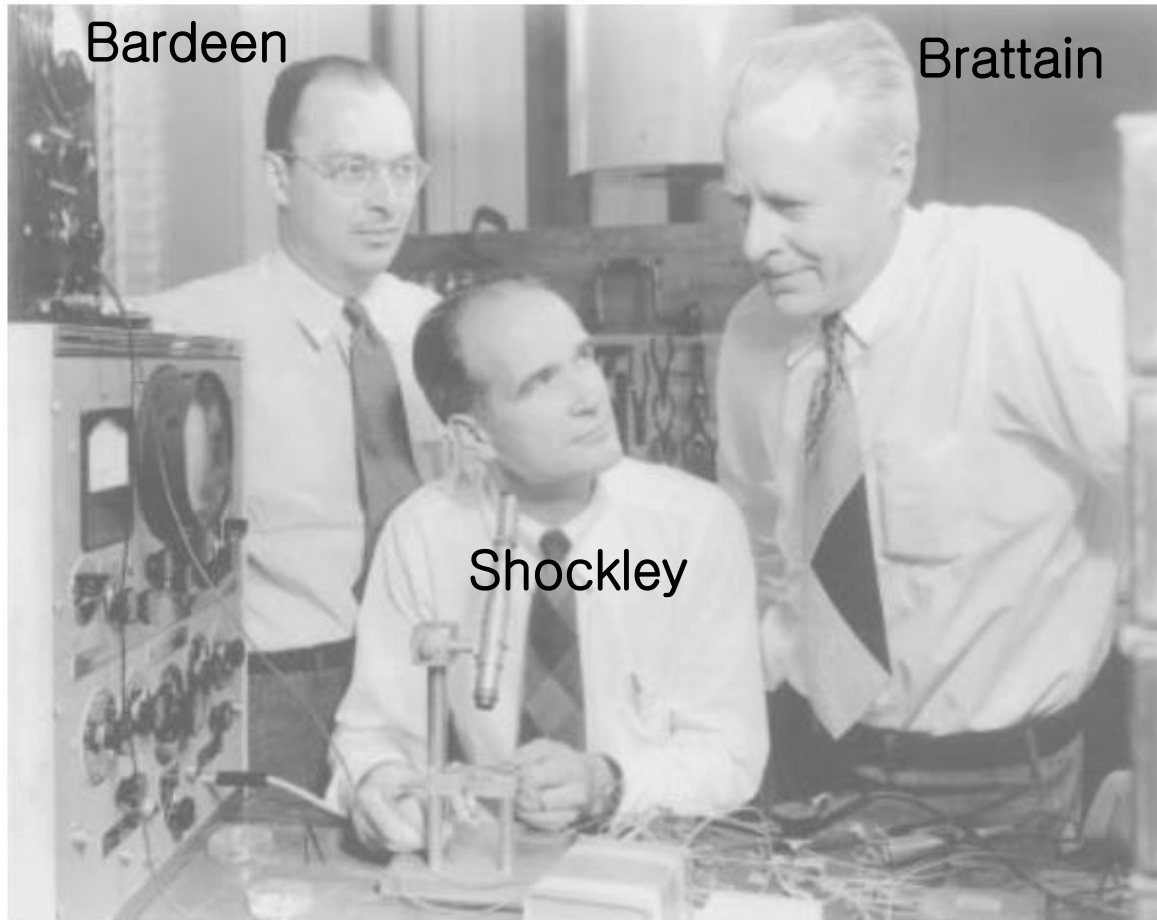
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Introduction

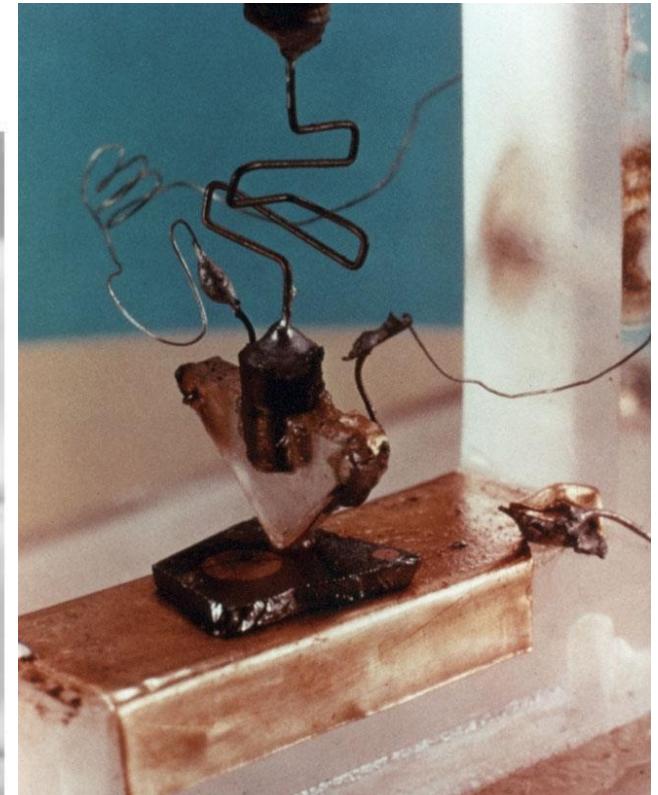
- *p-n junctions* are of great importance both in modern electronic applications and in understanding other semiconductor devices.
- The p-n junction theory serves as the *foundation of the physics of semiconductor devices*.
- The basic theory of *current-voltage characteristics* of *p-n* junctions was established by **Shockley**. This theory was then extended by Sah, Noyce, and Shockley, and by Moll.
- The *basic equations* (previous class) are used to develop the *ideal static* and *dynamic* characteristics of *p-n* junctions.
- Departures *from the ideal characteristics* due to *generation* and *recombination* in the *depletion layer*, *to high injection*, and *to series resistance effects* are then discussed. *Junction breakdown*, especially that due to *avalanche multiplication*, is considered in detail, after which transient behavior and noise performance in *p-n* junctions are presented.
- A *p-n* junction is a *two-terminal device*. Depending on the *doping profile*, *device geometry*, and *biasing condition*, a p-n junction can perform *various* terminal functions which are considered briefly in Section 2.6.
- The chapter closes with a discussion of an important group of devices-the *heterojunctions*, which are junctions formed *between dissimilar semiconductors* (e.g., *n-type GaAs* on *p-type AlGaAs*).

Introduction

Bardeen, Shockley, and Brattain



Nobel Prize (Physics) @1956



The first point-contact Transistor invented.

Bell Telephone Lab.
Dec. 23, 1947

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Depletion Region: Abrupt Junction

Built-in Potential and Depletion-Layer Width

- When the **impurity concentration** in a semiconductor **changes abruptly** from acceptor impurities N_A to donor impurities N_D , as shown in Fig. 1(a), one obtains an **abrupt junction**.
- In particular, if $N_A \gg N_D$ (or vice versa), one obtains a **one-sided abrupt** p^+-n (or n^+-p) junction.
- We first consider the **thermal equilibrium** condition, that is, one **without applied voltage and current flow**. From the current equation of drift and diffusion

$$J_n = 0 = q\mu_n \left(n\mathcal{E} + \frac{kT}{q} \frac{dn}{dx} \right) = \mu_n n \frac{dE_F}{dx} \quad (1)$$

OR

$$\frac{dE_F}{dx} = 0. \quad (2)$$

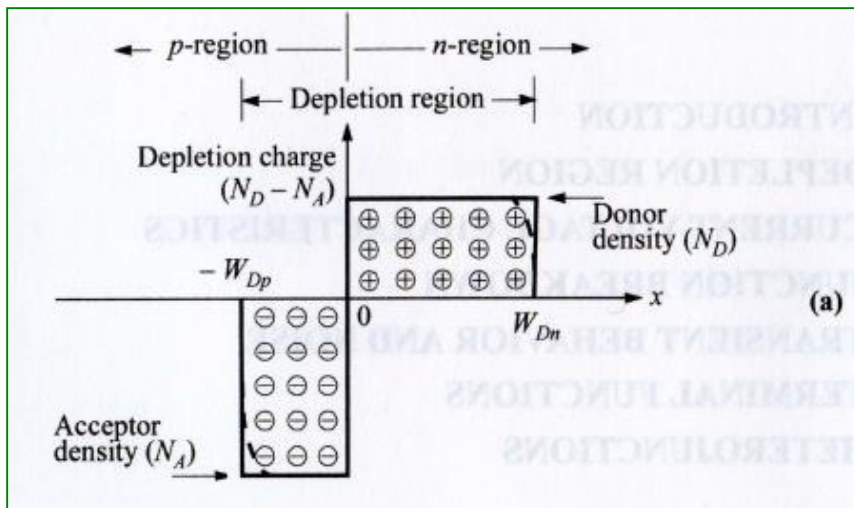


Fig. 1 Abrupt p - n junction in **thermal equilibrium**.
(a) **Space-charge distribution**. Dashed lines indicate corrections to depletion approximation.

Abrupt Junction

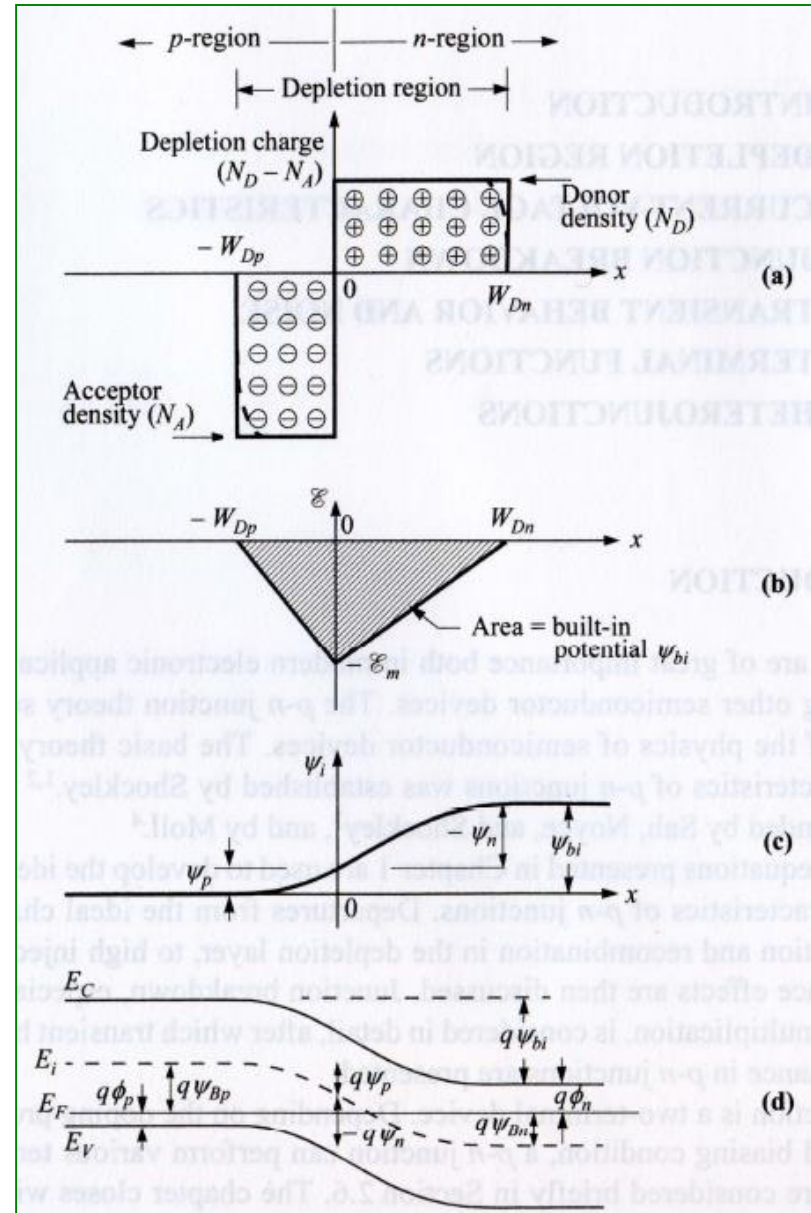
- Similarly,

$$J_p = 0 = \mu_p p \frac{dE_F}{dx} \quad (3)$$

- Thus the condition of **zero net electron and hole currents** requires that the **Fermi level must be constant** throughout the sample. The **built-in potential** ψ_{bi} , or **diffusion potential**, as shown in Fig. 1 b, c, and d, is equal to

$$q\psi_{bi} = E_g - (q\phi_n + q\phi_p) = q\psi_{Bn} + q\psi_{Bp} \quad (4)$$

Fig. 1 Abrupt p - n junction in thermal equilibrium.
(a) Space-charge distribution. Dashed lines indicate corrections to depletion approximation.
(b) Electric-field distribution.
(c) Potential distribution where ψ_{bi} is the built-in potential. (d) Energy-band diagram.



Abrupt Junction

- For nondegenerate semiconductors,

$$\psi_{bi} = \frac{kT}{q} \ln\left(\frac{n_{no}}{n_i}\right) + \frac{kT}{q} \ln\left(\frac{p_{po}}{n_i}\right) \approx \frac{kT}{q} \ln\left(\frac{N_D N_A}{n_i^2}\right) . \quad (5)$$

- Since at equilibrium $n_{no} p_{no} = n_{po} p_{po} = n_i^2$,

$$\psi_{bi} = \frac{kT}{q} \ln\left(\frac{p_{po}}{p_{no}}\right) = \frac{kT}{q} \ln\left(\frac{n_{no}}{n_{po}}\right) . \quad (6)$$

This gives the relationship between carrier densities on either side of the junction.

- If one or both sides of the junction are **degenerate**, *care* has to be taken in calculating the **Fermi-levels** and **built-in potential**.
- Equation (4) has to be used since Boltzmann statistics *cannot* be used to **simplify** the Fermi-Dirac integral.
- Furthermore, **incomplete ionization** has to be considered, i.e., $n_{no} \neq N_D$ and/or $p_{po} \neq N_A$.

Abrupt Junction

- Next, we proceed to calculate the **field** and **potential distribution** inside the **depletion region**.
- To *simplify* the analysis, the **depletion approximation** is used which assumes that the depleted charge has a box profile.
- Since in **thermal equilibrium** the **electric field in the neutral regions** (*far from* the junction at either side) **of the Semiconductor must be zero**, the total negative charge per unit area in the p-side must be precisely equal to the total positive charge per unit area in the n-side:

$$N_A W_{Dp} = N_D W_{Dn} \quad (7)$$

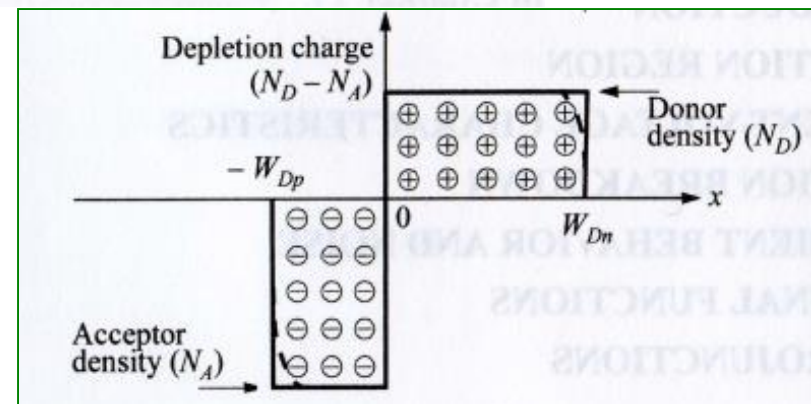
- From the Poisson equation we obtain

$$-\frac{d^2 \psi_i}{dx^2} = \frac{d\mathcal{E}}{dx} = \frac{\rho(x)}{\epsilon_s} = \frac{q}{\epsilon_s} [N_D^+(x) - n(x) - N_A^-(x) + p(x)] \quad (8)$$

- Inside the depletion region, $n(x) \approx p(x) \approx 0$, and **assuming complete ionization**,

$$\frac{d^2 \psi_i}{dx^2} \approx \frac{q N_A}{\epsilon_s} \quad \text{for } -W_{Dp} \leq x \leq 0, \quad (9a)$$

$$-\frac{d^2 \psi_i}{dx^2} \approx \frac{q N_D}{\epsilon_s} \quad \text{for } 0 \leq x \leq W_{Dn}. \quad (9b)$$



Abrupt Junction

- The **electric field** is then obtained **by integrating** the above equations, as shown in Fig. 1b:

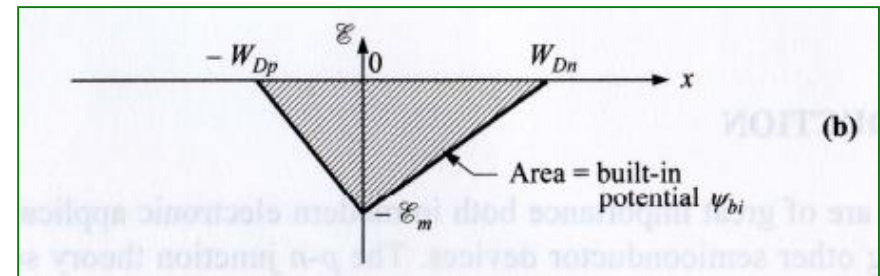
$$\mathcal{E}(x) = -\frac{qN_A(x + W_{Dp})}{\epsilon_s} \quad \text{for } -W_{Dp} \leq x \leq 0, \quad (10)$$

$$\mathcal{E}(x) = -\mathcal{E}_m + \frac{qN_D x}{\epsilon_s} = -\frac{qN_D}{\epsilon_s}(W_{Dn} - x) \quad \text{for } 0 \leq x \leq W_{Dn} \quad (11)$$

where \mathcal{E}_m is the **maximum field** that exists at $x = 0$ and is given by

$$|\mathcal{E}_m| = \frac{qN_D W_{Dn}}{\epsilon_s} = \frac{qN_A W_{Dp}}{\epsilon_s}. \quad (12)$$

-Integrating Eqs. 10 and 11 once again gives the potential distribution $\psi_i(x)$ (Fig. 1c)



$$\psi_i(x) = \frac{qN_A}{2\epsilon_s}(x + W_{Dp})^2 \quad \text{for } -W_{Dp} \leq x \leq 0, \quad (13)$$

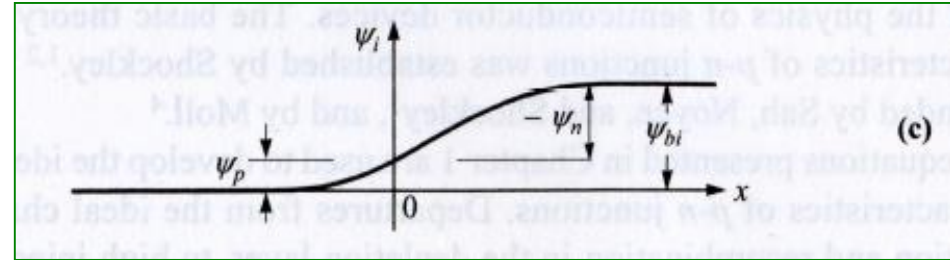
$$\psi_i(x) = \psi_i(0) + \frac{qN_D}{\epsilon_s}\left(W_{Dn} - \frac{x}{2}\right)x \quad \text{for } 0 \leq x \leq W_{Dn}. \quad (14)$$

Abrupt Junction

- With these, the potentials across different regions can be found as:

$$\psi_p = \frac{qN_A W_{Dp}^2}{2\epsilon_s}, \quad (15a)$$

$$|\psi_n| = \frac{qN_D W_{Dn}^2}{2\epsilon_s}, \quad (15b)$$



(ψ_n is relative to the n -type bulk and is thus **negative**.)

$$\psi_{bi} = \psi_p + |\psi_n| = \psi_i(W_{Dn}) = \frac{|\mathcal{E}_m|}{2}(W_{Dp} + W_{Dn}) \quad (16)$$

where \mathcal{E}_m can also be expressed as:

$$|\mathcal{E}_m| = \sqrt{\frac{2qN_A \psi_p}{\epsilon_s}} = \sqrt{\frac{2qN_D |\psi_n|}{\epsilon_s}}. \quad (17)$$

From Eqs. 16 and 7, the depletion widths are calculated to be:

$$W_{Dp} = \sqrt{\frac{2\epsilon_s \psi_{bi}}{q} \frac{N_D}{N_A(N_A + N_D)}} \quad (18a)$$

$$W_{Dn} = \sqrt{\frac{2\epsilon_s \psi_{bi}}{q} \frac{N_A}{N_D(N_A + N_D)}} \quad (18b)$$

$$W_{Dp} + W_{Dn} = \sqrt{\frac{2\epsilon_s}{q} \left(\frac{N_A + N_D}{N_A N_D} \right) \psi_{bi}}. \quad (19)$$

Abrupt Junction

- The following relationships can be further deduced:

$$\frac{|\psi_n|}{\psi_{bi}} = \frac{W_{Dn}}{W_{Dp} + W_{Dn}} = \frac{N_A}{N_A + N_D}, \quad (20a)$$

$$\frac{\psi_p}{\psi_{bi}} = \frac{W_{Dp}}{W_{Dp} + W_{Dn}} = \frac{N_D}{N_A + N_D}. \quad (20b)$$

- For a **one-sided abrupt** junction (p^+-n or n^+-p), Eq. 4 is used to calculate the **built-in potential**.
- In this case, the **majority of the potential variation and depletion region** will be **inside the lightly doped side**.
- Equation 19 reduces to

$$W_D = \sqrt{\frac{2\epsilon_s \psi_{bi}}{qN}} \quad (21)$$

where N is N_D or N_A depending on whether $N_A \gg N_D$ or vice versa, and

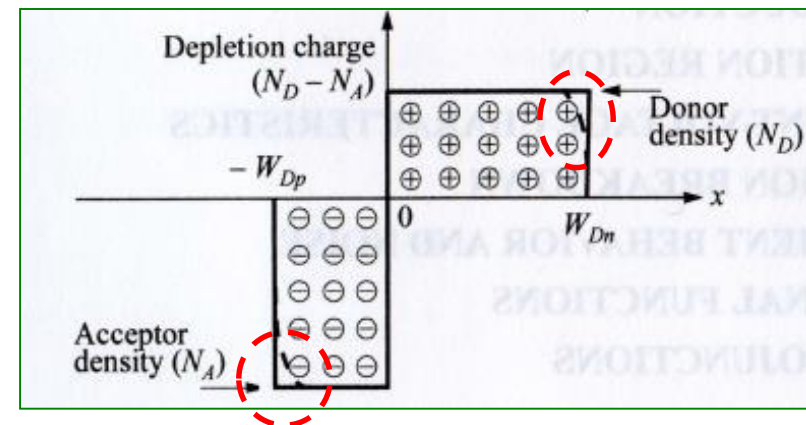
$$\psi_i(x) = |\mathcal{E}_m| \left(x - \frac{x^2}{2W_D} \right). \quad (22)$$

This discussion uses **box profiles** for the depletion charges, i.e., **depletion approximation**.

Abrupt Junction

- A **more accurate** result for the depletion-layer properties can be obtained by **considering the majority-carrier contribution** in addition to the **impurity concentration** in the Poisson equation, that is, $\rho \approx -q[N_A - p(x)]$ on the ***p*-side** and $\rho \approx q[N_D - n(x)]$ on the ***n*-side**.
- The **depletion width** is essentially the **same** as given by Eq. 19, **except** that ψ_{bi} is replaced by $(\psi_{bi} - 2kT/q)$.
- The **correction factor $2kT/q$** comes about because of the **two majority-carrier distribution tails** (electrons in *n*-side and holes in *p*-side, as shown by the dashed lines in Fig. 1 (a) near the edges of the depletion region). Each contributes a correction factor kT/q .
- The depletion-layer width at thermal equilibrium for a one-sided abrupt junction becomes

$$W_D = \sqrt{\frac{2\epsilon_s}{qN} \left(\psi_{bi} - \frac{2kT}{q} \right)} \quad (23)$$



Abrupt Junction

With applied bias

- Furthermore, **when a voltage V is applied** to the junction, the total electrostatic potential variation across the junction is given by $(\psi_{bi} - V)$ where V is **positive for forward bias** (positive voltage on p-region with respect to n-region) and **negative for reverse bias**.
- Substituting $(\psi_{bi} - V)$ for ψ_{bi} in Eq. 23 yields the depletion-layer width as a function of the applied voltage. The results for **one-sided abrupt junctions** in silicon are shown in Fig. 2.
- The net potential at zero bias is near 0.8 V for Si and 1.3 V for GaAs. This **net potential** will be **decreased under forward bias** and **increased under reverse bias**.
- These results can also be used for GaAs since both Si and GaAs have approximately the same static dielectric constants.
- To obtain the depletion-layer width for other semiconductors such as Ge, one must multiply the results of Si by the factor $\sqrt{\epsilon_s(\text{Ge})/\epsilon_s(\text{Si})}$ ($= 1.16$).
- The simple model above can give adequate predictions for most abrupt p - n junctions.

Abrupt Junction

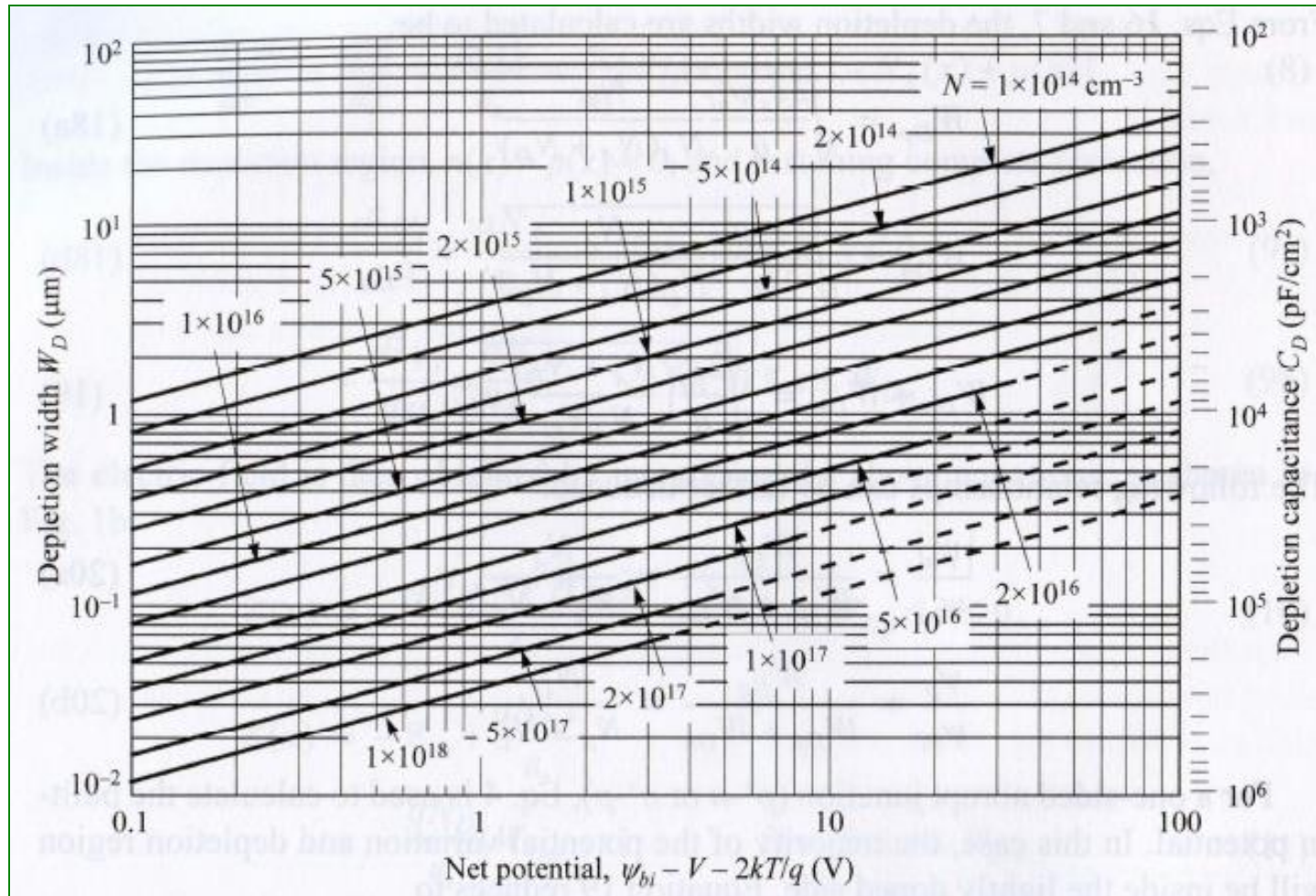


Fig.2 Depletion-layer width and depletion-layer capacitance per unit area as a function of net potential ($\psi_{bi} - V - 2kT/q$) for one-sided abrupt junctions in Si. Doping N is from the lightly doped side. Dashed lines represent breakdown conditions.

Depletion Layer Capacitance

- The **depletion-layer capacitance** per unit area is defined as $C_D = dQ_D/dV = \epsilon_s/W_D$, where dQ_D is the **incremental depletion charge** on each side of the junction (**total charge is zero**) upon an incremental change of the applied voltage dV .
- For one-sided abrupt junctions, the capacitance per unit area is given by

$$C_D = \frac{\epsilon_s}{W_D} = \sqrt{\frac{q\epsilon_s N}{2}} \left(\psi_{bi} - V - \frac{2kT}{q} \right)^{-1/2} \quad (24)$$

where V is positive/negative for forward/reverse bias. The results of the depletion layer capacitance are also shown in Fig. 2.

- Rearrange the above equation leads to:

$$\frac{1}{C_D^2} = \frac{2}{q\epsilon_s N} \left(\psi_{bi} - V - \frac{2kT}{q} \right), \quad (25)$$

$$\frac{d(1/C_D^2)}{dV} = -\frac{2}{q\epsilon_s N}. \quad (26)$$

- It is clear from Eqs. 25 and 26 that by plotting $1/C^2$ versus V , a **straight line** should result from a **one-sided abrupt junction** (Fig. 3). The slope gives the impurity concentration of the substrate (N), and the extrapolation to $1/C^2 = 0$ gives $(\psi_{bi} - 2kT/q)$. Note that, for the forward bias, a diffusion capacitance exists in addition to the depletion capacitance mentioned previously. The diffusion capacitance will be discussed in Section 2.3.4.

Depletion Layer Capacitance

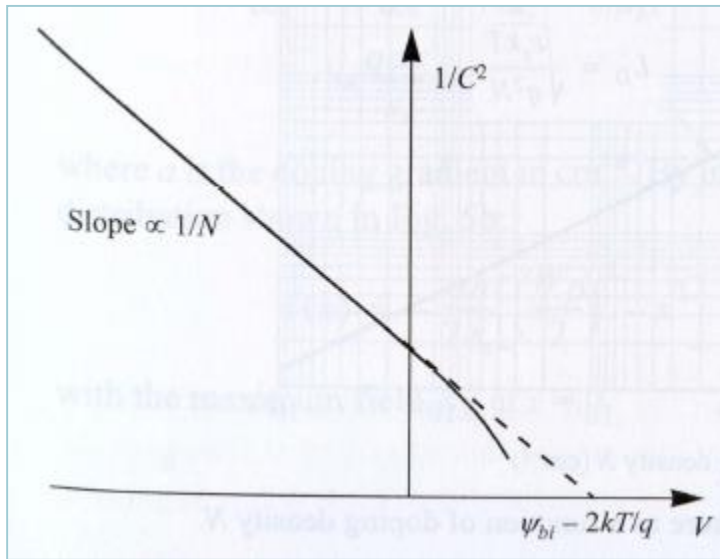


Fig. 3 A $1/C^2-V$ plot can yield the built-in potential and doping density N .

Note that the semiconductor **potential** and the **capacitance-voltage** data are insensitive to changes in the doping profiles that occur in a distance less than a **Debye length**. The Debye length L_D is a characteristic length for semiconductors and is defined as

$$L_D \equiv \sqrt{\frac{\epsilon_s kT}{q^2 N}}. \quad (27)$$

- This **Debye length** gives an idea of the **limit of the potential change in response to an abrupt change in the doping profile**.
- Consider a case where the doping has a small increase of ΔN_D in the background of N_D , the change of potential $\Delta \psi_i(x)$ near the step is given by

$$n = N_D \exp\left(\frac{\Delta \psi_i q}{kT}\right), \quad (28)$$

Depletion Layer Capacitance

$$\begin{aligned} \frac{d^2 \Delta \psi_i}{dx^2} &= -\frac{q}{\epsilon_s} (N_D + \Delta N_D - n) = -\frac{q N_D}{\epsilon_s} \left[1 + \frac{\Delta N_D}{N_D} - \exp\left(\frac{\Delta \psi_i q}{kT}\right) \right] \\ &\approx -\frac{q N_D}{\epsilon_s} \left[1 + \frac{\Delta N_D}{N_D} - \left(1 + \frac{\Delta \psi_i q}{kT} \right) \right] \approx \frac{q^2 N_D}{\epsilon_s kT} \Delta \psi_i \end{aligned} \quad (29)$$

whose solution has a decay length given by Eq. 27.

- This implies that if the doping profile changes *abruptly* in a scale less than the Debye length, this variation has *no effect* and *cannot be resolved*, and that if the depletion width is smaller than the Debye length, the analysis using the *Poisson equation* is *no longer valid*.

- At **thermal equilibrium** the depletion-layer widths of abrupt junctions are about $8L_D$ for Si, and $10L_D$ for GaAs. The Debye length as a function of doping density is shown in Fig. 4 for silicon at room temperature.

- For a doping density of 10^{16} cm^{-3} , the Debye length is 40 nm; for other dopings, L_D will vary as $1/\sqrt{N}$, that is, a reduction by a factor of 3.16 per decade.

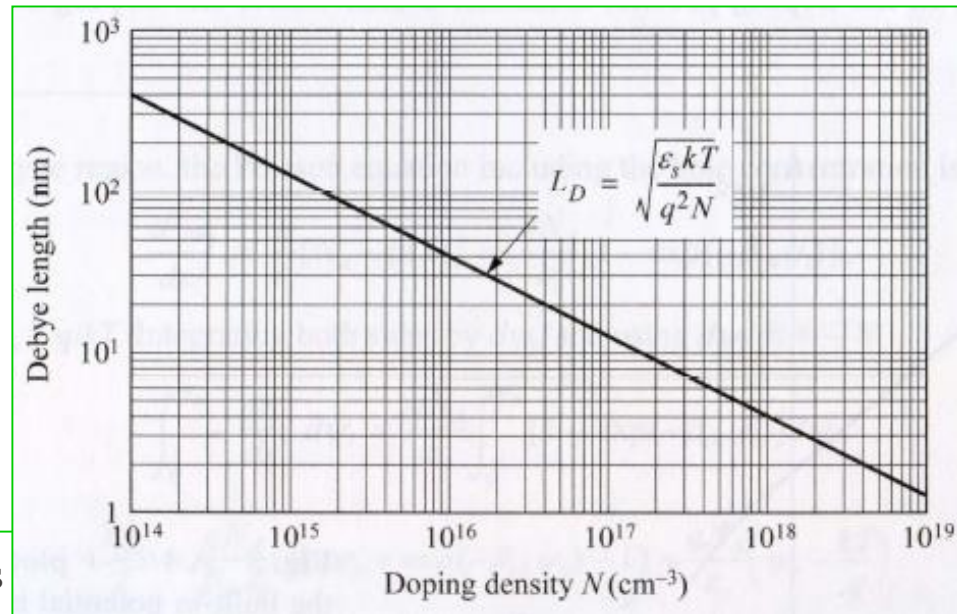


Fig.4 Debye length in Si at room temperature as a function of doping density N .