

Lecture 4

Density of States and Fermi Energy Concepts

Reading:

(Cont'd) Pierret 2.1-2.6



Density of States Concept

Quantum Mechanics tells us that the number of available states in a cubic cm per unit of energy, the density of states, is given by:

$$g_c(E) = \frac{m_n^* \sqrt{2m_n^*(E - E_c)}}{\pi^2 \hbar^3}, E \ge E_c$$

$$g_{\nu}(E) = \frac{m_p^* \sqrt{2m_p^*(E_{\nu} - E)}}{\pi^2 \hbar^3}, E \leq E_{\nu}$$

$$unit \equiv \left[\frac{Number\ of\ States}{m^3} \right]_{Joule} \Rightarrow \left[\frac{Number\ of\ States}{cm^3} \right]_{eV}$$





Density of States Concept

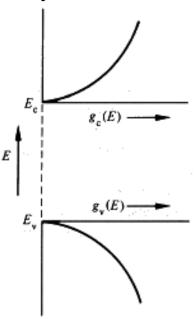


Figure 2.14 General energy dependence of $g_c(E)$ and $g_v(E)$ near the band edges. $g_c(E)$ and $g_v(E)$ are the density of states in the conduction and valence bands, respectively.

Thus, the number of states per cubic centimeter between energy E' and E'+dE is

$$g_c(E')dE$$
 if $E' \ge E_c$ and,
 $g_v(E')dE$ if $E' \le E_v$ and,
 0 otherwise



Probability of Occupation (Fermi Function) Concept

Now that we know the number of available states at each energy, how do the electrons occupy these states?

We need to know how the electrons are "distributed in energy".

Again, Quantum Mechanics tells us that the electrons follow the "Fermi-distribution function".

$$f(E) = \frac{1}{1 + e^{(E - E_F)/kT}}$$
 where $k = Boltzman$ constant, $T = Temperature$ in Kelvin

and $E_F \equiv$ Fermi energy (~ average energy of our electrons in the crystal)

f(E) is the probability that a state at energy E is *occupied*

1-f(E) is the probability that a state at energy E is *unoccupied*



An Aside about the fermi energy:

- •If the material had an imbalance of average electron energy from one side to another, electrons would flow from the region of high energy toward the region of low energy to balance out the average energy. Thus, gradients in the average energy would result in a current (flow of electrons) until the energy balance was satisfied. If an outside force maintained such an imbalance in energy from one region to another, a continuous current would result (see "quasi-fermi" energy concept introduced later).
- •When "assembling our crystal", if we brought extra electrons into the crystal such as in the case of donor doping, we have to do more work to introduce the extra valence electrons. Thus, the average electron energy would increase compared to the intrinsic doped case.
- •When "assembling our crystal", if we brought fewer electrons into the crystal such as in the case of acceptor doping, we have to do less work to introduce the fewer valence electrons. Thus, the average electron energy would decrease compared to the intrinsic doped case.



Probability of Occupation (Fermi Function) Concept

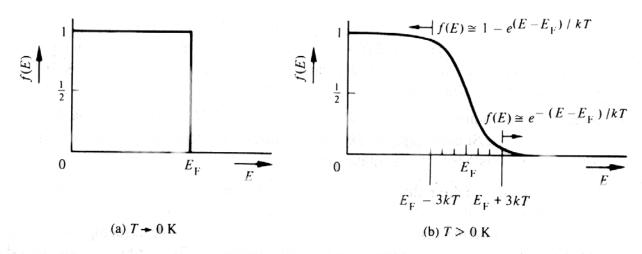


Figure 2.15 Energy dependence of the Fermi function. (a) $T \to 0$ K; (b) generalized T > 0 K plot with the energy coordinate expressed in kT units.

At T=0K, occupancy is "digital": No occupation of states above E_F and complete occupation of states below E_F

At T>0K, occupation probability is reduced with increasing energy.

 $f(E=E_F) = 1/2$ regardless of temperature.



Probability of Occupation (Fermi Function) Concept

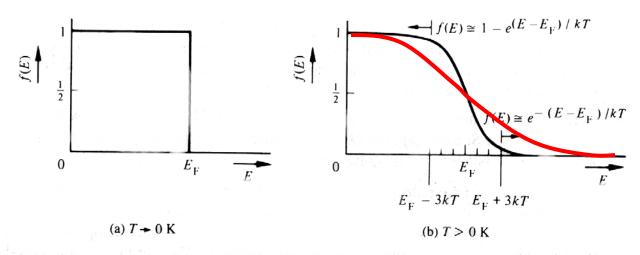


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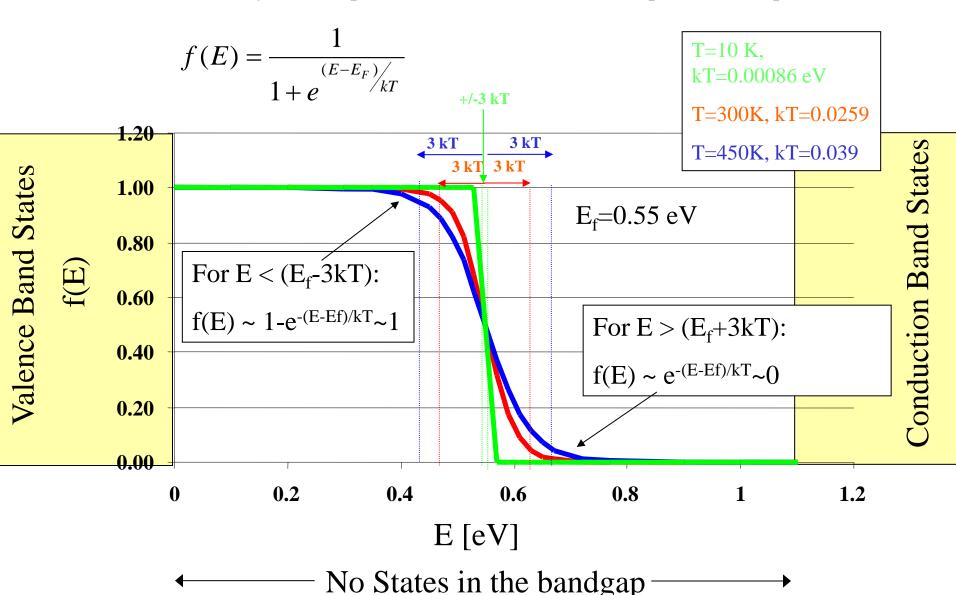
At higher temperatures, higher energy states can be occupied, leaving more lower energy states unoccupied (1-f(E)).

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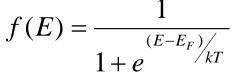
How do electrons and holes populate the bands?

Probability of Occupation (Fermi Function) Concept – Si Example

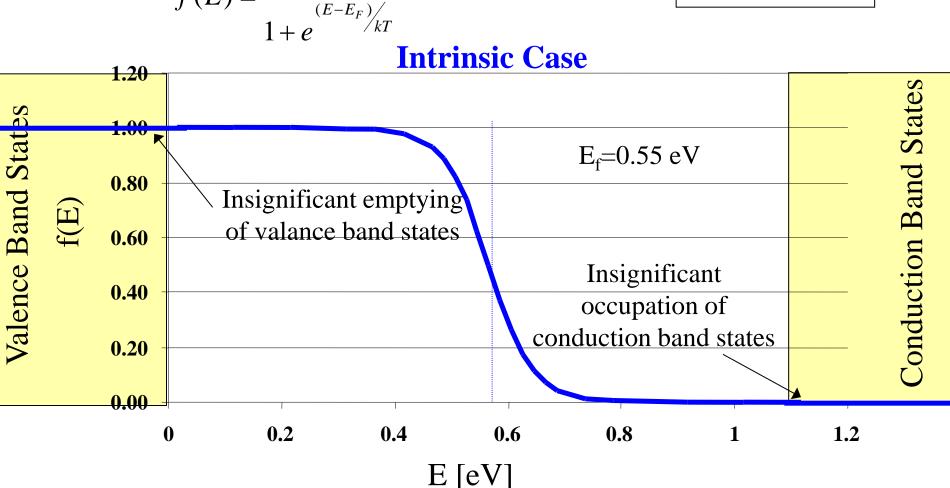




Probability of Occupation (Fermi Function) Concept – Si Example



T=450K, kT=0.039



No States in the bandgap

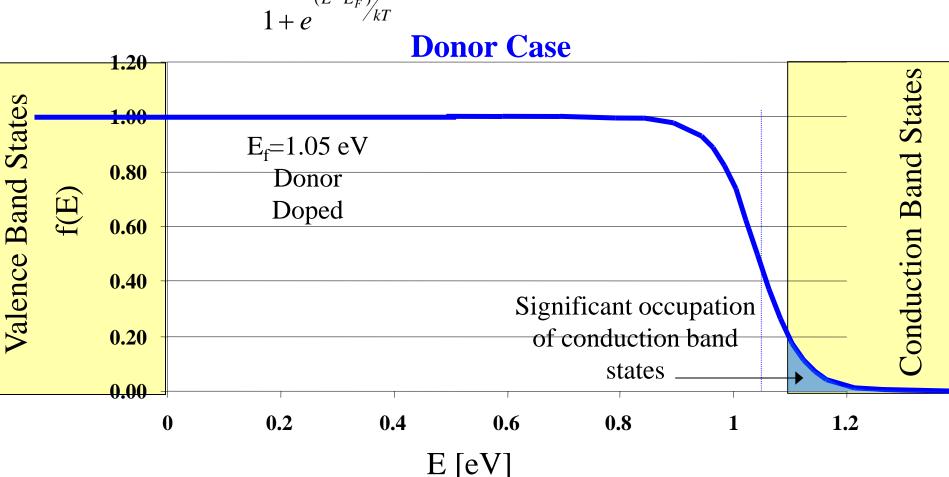
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Probability of Occupation (Fermi Function) Concept – Si Example

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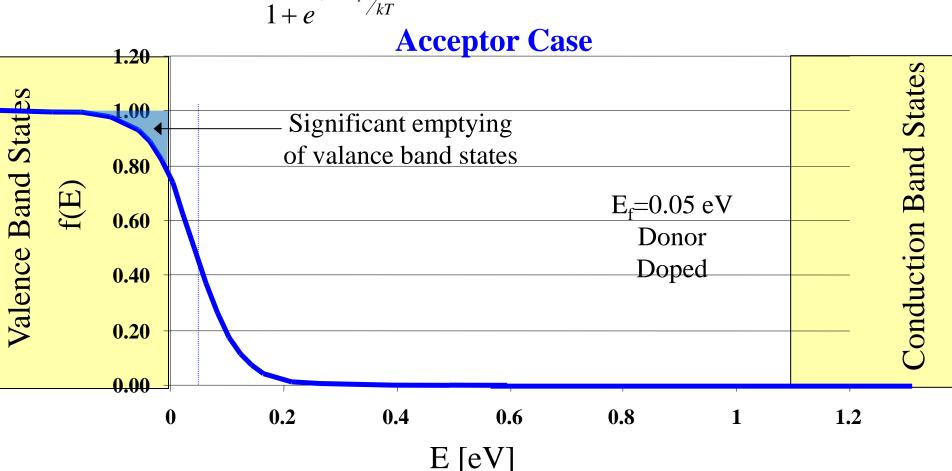
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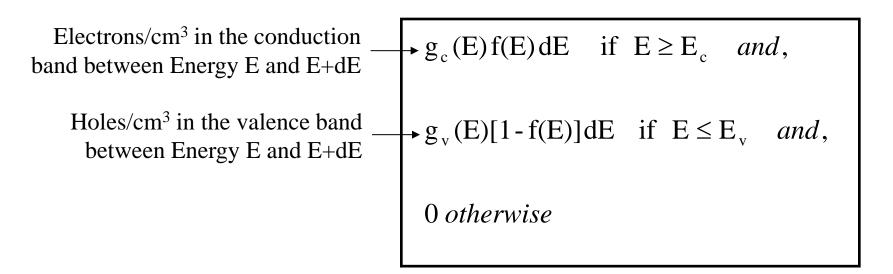
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Probability of Occupation

Thus, the density of electrons (or holes) occupying the states in energy between E and E+dE is:





Band Occupation

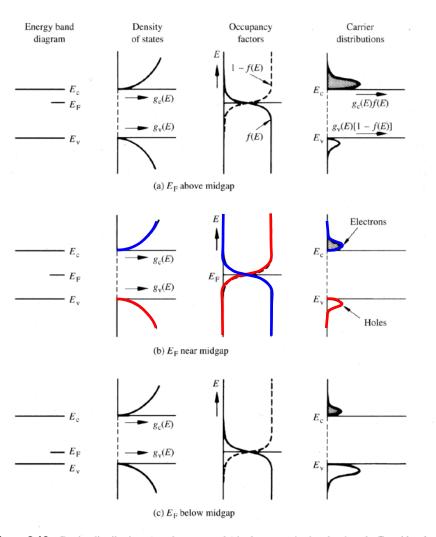


Figure 2.16 Carrier distributions (not drawn to scale) in the respective bands when the Fermi level is positioned (a) above midgap, (b) near midgap, and (c) below midgap. Also shown in each case are coordinated sketches of the energy band diagram, density of states, and the occupancy factors (the Fermi function and one minus the Fermi function).



Intrinsic Energy (or Intrinsic Level)

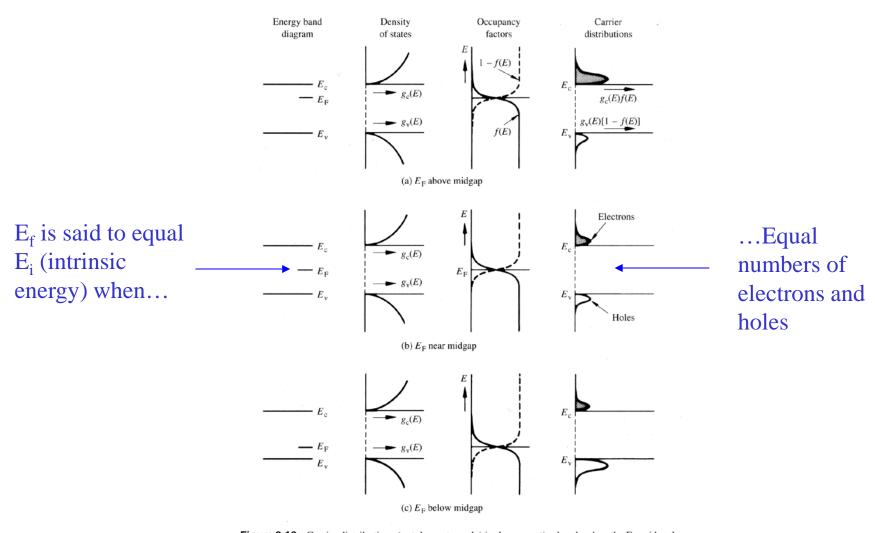


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Correction for the Presence of Additional Dopant States

