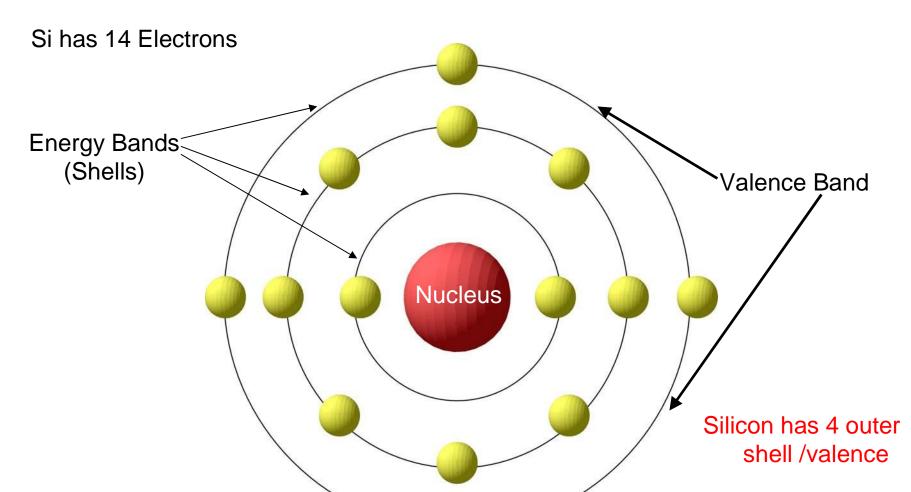




1

Silicon

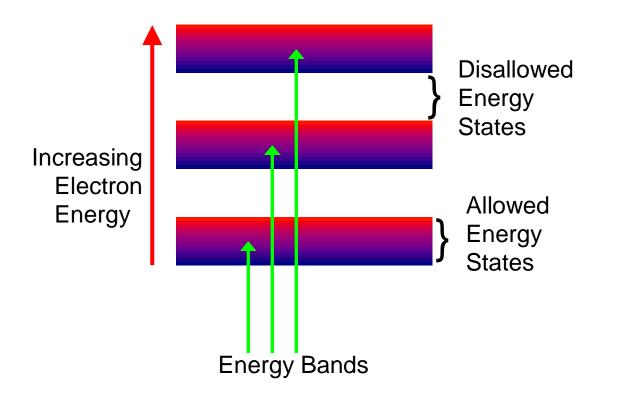
Silicon is the primary semiconductor used in VLSI systems



electrons

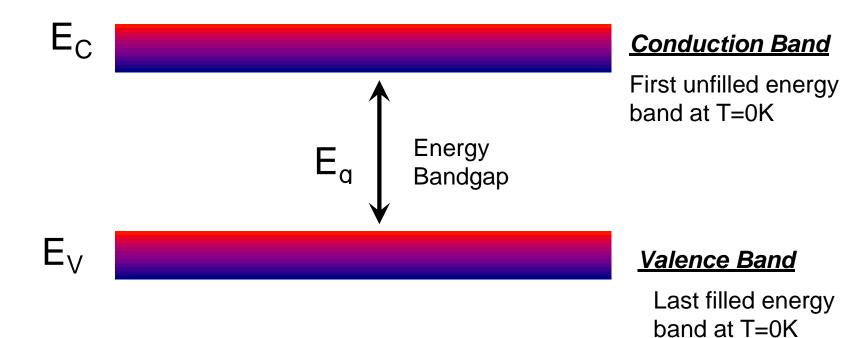
At T=0K, the highest energy band occupied by an electron is called the valence band.

Energy Bands

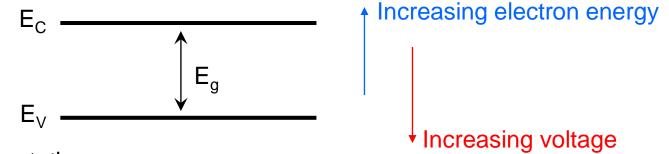


- Electrons try to occupy the lowest energy band possible
- Not every energy level is a legal state for an electron to occupy
- These legal states tend to arrange themselves in bands

Energy Bands



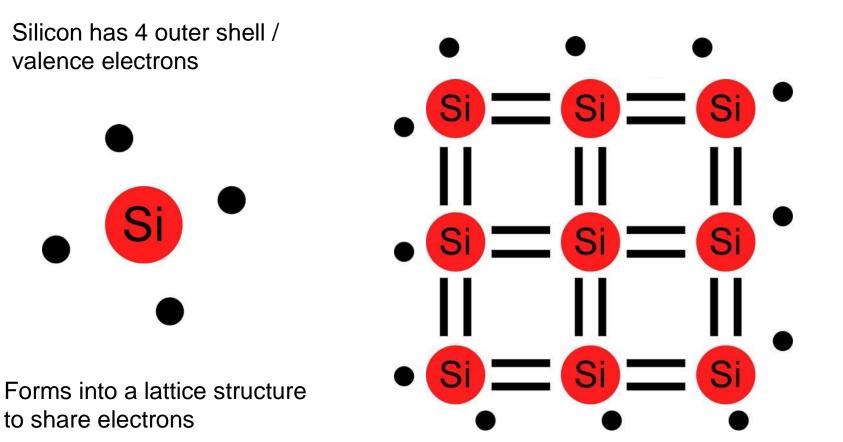
Band Diagrams



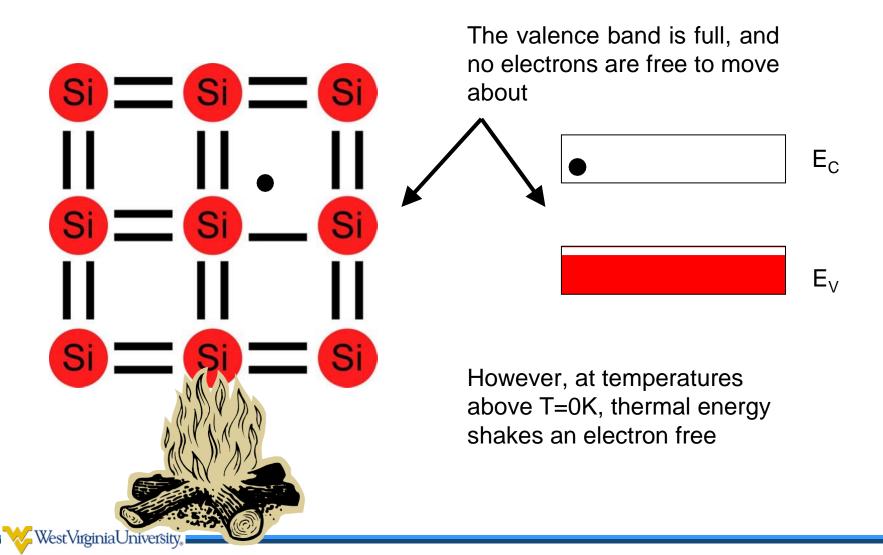
Band Diagram Representation Energy plotted as a function of position

- $E_{c} \rightarrow Conduction band$
 - \rightarrow Lowest energy state for a free electron
- $E_{V} \rightarrow Valence band$
 - \rightarrow Highest energy state for filled outer shells
- $E_G \rightarrow Band gap$
 - \rightarrow Difference in energy levels between E_C and E_V
 - \rightarrow No electrons (e⁻) in the bandgap (only above E_C or below E_V)
 - \rightarrow E_G = 1.12eV in Silicon

Intrinsic Semiconductor



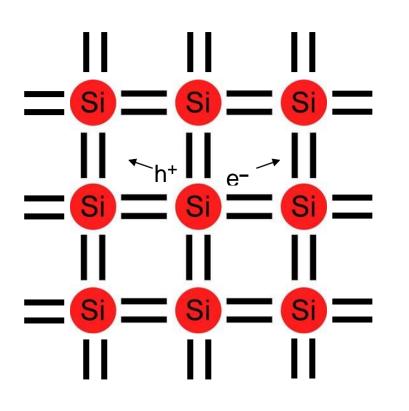
Intrinsic Silicon



Semiconductor Properties

<u>For T > 0K</u>

Electron shaken free and can cause current to flow

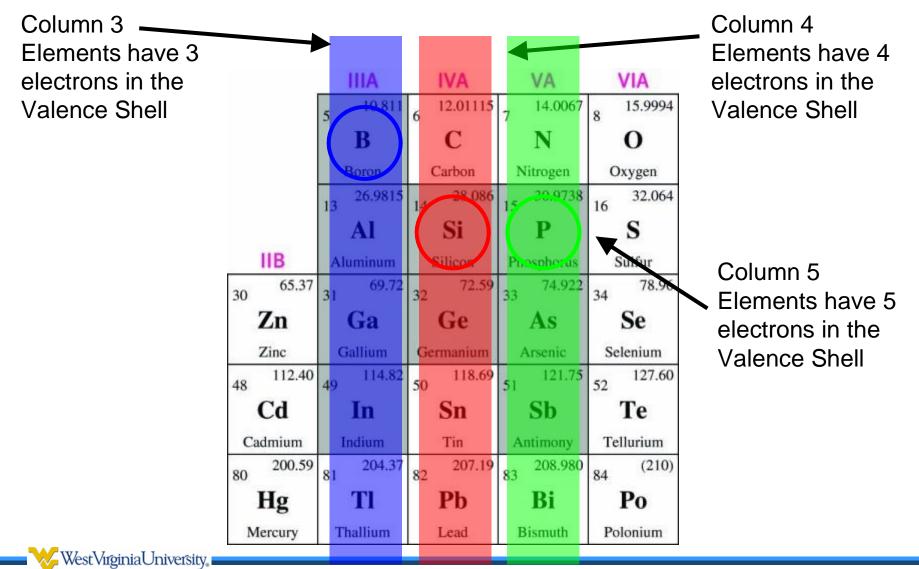


- <u>Generation</u> Creation of an electron (e⁻) and hole (h⁺) pair
- h⁺ is simply a missing electron, which leaves an excess positive charge (due to an extra proton)
- <u>Recombination</u> if an e⁻ and an h⁺ come in contact, they annihilate each other
- Electrons and holes are called "carriers" because they are charged particles – when they move, they carry current
- Therefore, semiconductors can conduct electricity for T > 0K ... but not much current (at room temperature (300K), pure silicon has only 1 free electron per 3 trillion atoms)

Doping

- <u>Doping</u> Adding impurities to the silicon crystal lattice to increase the number of carriers
- Add a small number of atoms to increase either the number of electrons or holes

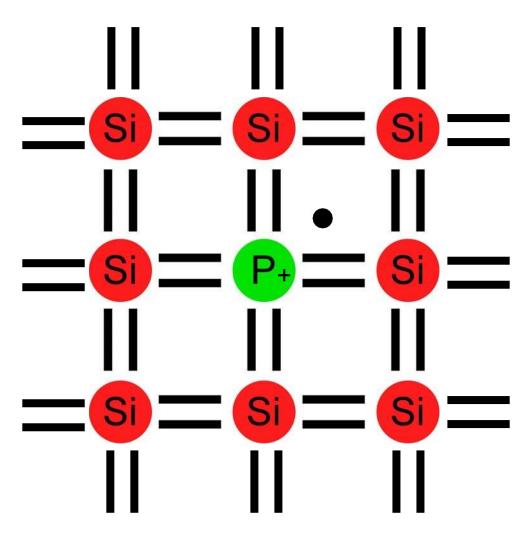
Periodic Table



Donors n-Type Material

<u>Donors</u>

- Add atoms with 5 valence-band electrons
- ex. Phosphorous (P)
- "Donates" an extra e⁻ that can freely travel around
- Leaves behind a positively charged nucleus (cannot move)
- Overall, the crystal is still electrically neutral
- Called "n-type" material (added <u>n</u>egative carriers)
- N_D = the concentration of donor atoms [atoms/cm³ or cm⁻³] ~10¹⁵-10²⁰cm⁻³
- e⁻ is free to move about the crystal (Mobility μ_n ≈1350cm²/V)

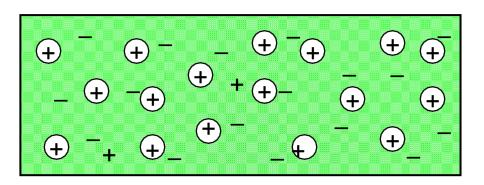


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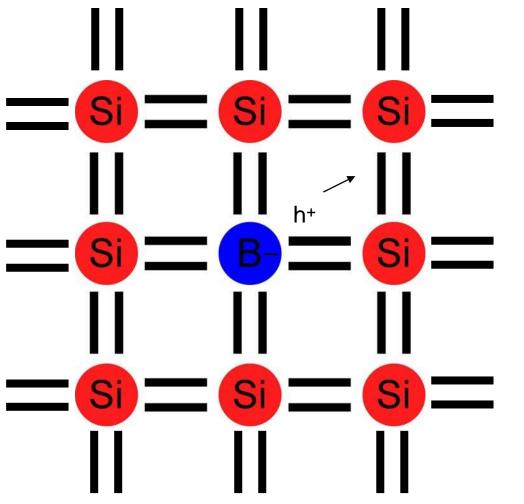
n-Type Material



Shorthand Notation

- Positively charged ion; immobile
- Negatively charged e-; mobile; Called "majority carrier"
- Positively charged h+; mobile; Called "minority carrier"

Acceptors Make p-Type Material

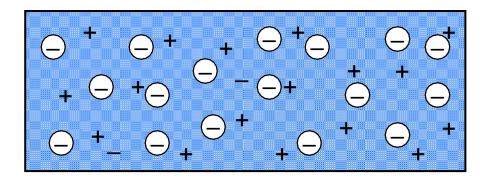


<u>Acceptors</u>

- Add atoms with only 3 valenceband electrons
- ex. Boron (B)
- "Accepts" e⁻ and provides extra h⁺ to freely travel around
- Leaves behind a negatively charged nucleus (cannot move)
- Overall, the crystal is still electrically neutral
- Called "p-type" silicon (added <u>p</u>ositive carriers)
- N_A = the concentration of acceptor atoms [atoms/cm³ or cm⁻³]
- Movement of the hole requires breaking of a bond! (This is hard, so mobility is low, µ_p ≈ 500cm²/V)

Acceptors Make p-Type Material

<u>p-Type Material</u>



Shorthand Notation O Negatively charged ion; immobile

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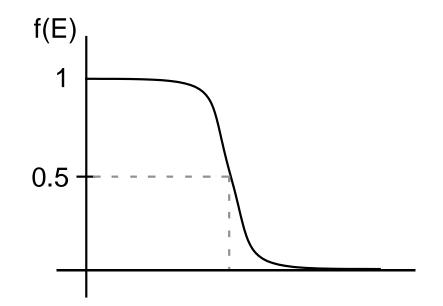
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The Fermi Function

<u>The Fermi Function</u>
Probability distribution function (PDF)
The probability that an available state at an energy E will be occupied by an e⁻

$$f(E) = \frac{1}{1 + e^{(E - E_f)/kT}}$$



The Fermi Function

- $E \rightarrow Energy level of interestE_{f} \rightarrow Fermi$ level
 - \rightarrow Halfway point
 - \rightarrow Where f(E) = 0.5
 - $k \rightarrow Boltzmann constant$
 - $= 1.38 \times 10^{-23} \text{ J/K}$
 - $= 8.617 \times 10^{-5} \, eV/K$
- $T \rightarrow$ Absolute temperature (in Kelvins)