

Fundamentals of Solid State Physics

Electronic Properties - The Tight-Binding Model

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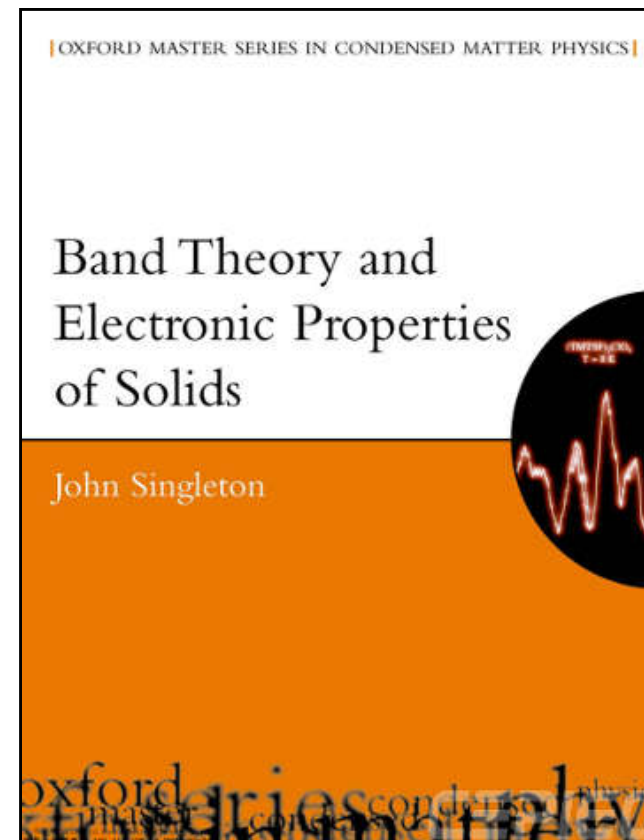
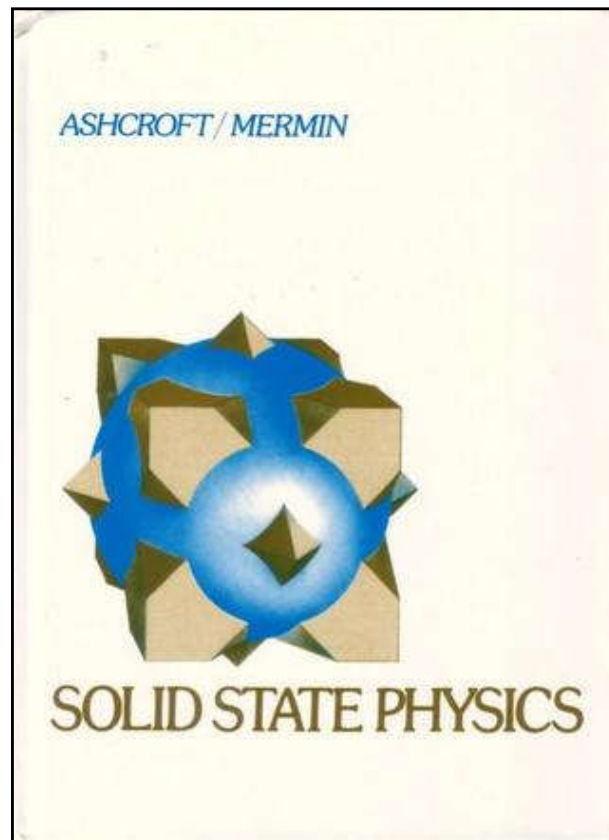


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Further Reading

- Ashcroft & Mermin, Chapter 10
- Singleton, Chapter 4



Real Electrons in Solids

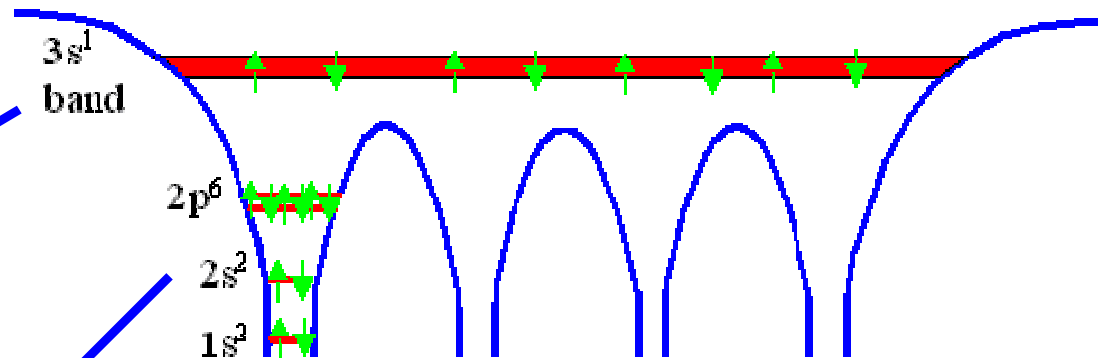
Electrons are in *periodic* potentials

→ **Bloch Wave**

$$\psi(\mathbf{r}) = \exp(i\mathbf{k} \cdot \mathbf{r}) \cdot u_{\mathbf{k}}(\mathbf{r})$$

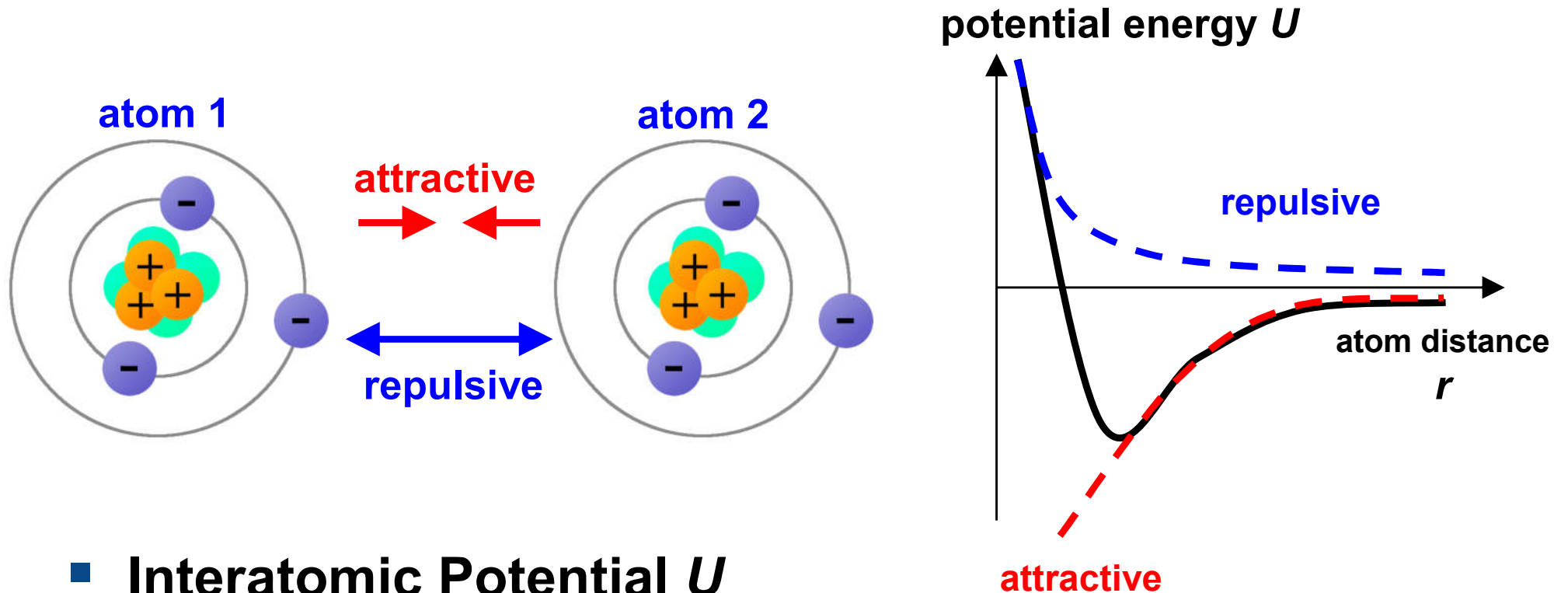
Nearly Free Electron Model
"近自由"近似

Tight Binding Model
"紧束缚"近似



Sodium (Na) [1s² 2s² 2p⁶] 3s¹

Atomic Bonding



■ Interatomic Potential U

- attraction: electrostatic (+ -)
- repulsion: electrostatic (+ + / - -)
and Pauli exclusion principle

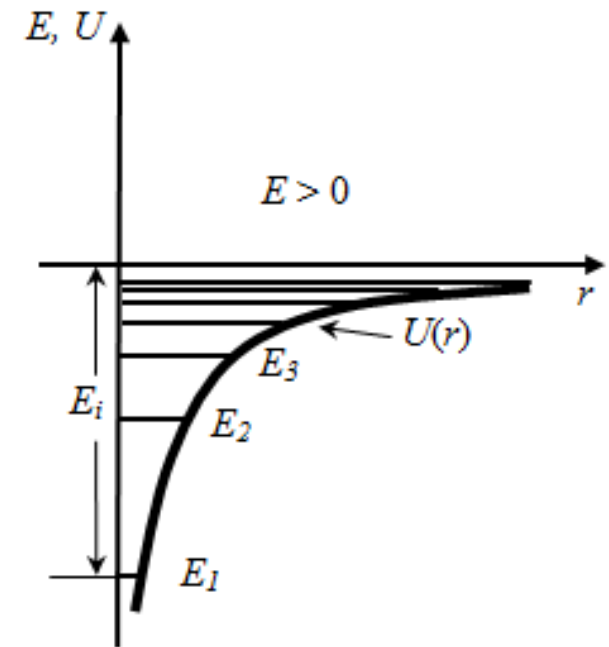
What are the quantum mechanic explanations?

Isolated Atoms

- Hydrogen atom

$$V(\mathbf{r}) = -\frac{e^2}{4\pi\epsilon_0} \frac{1}{r}$$

$$-\frac{\hbar^2}{2m} \nabla^2 \psi(\mathbf{r}) + V(\mathbf{r}) \cdot \psi(\mathbf{r}) = E \psi(\mathbf{r})$$



$$\psi(r, \theta, \varphi) = R_{nl}(r) \cdot Y_{lm}(\theta, \varphi)$$

$$E_n = -\frac{13.6 \text{ eV}}{n^2}$$

n, l, m - quantum numbers
 m_s - spin (+1/2, -1/2)

Isolated Atoms

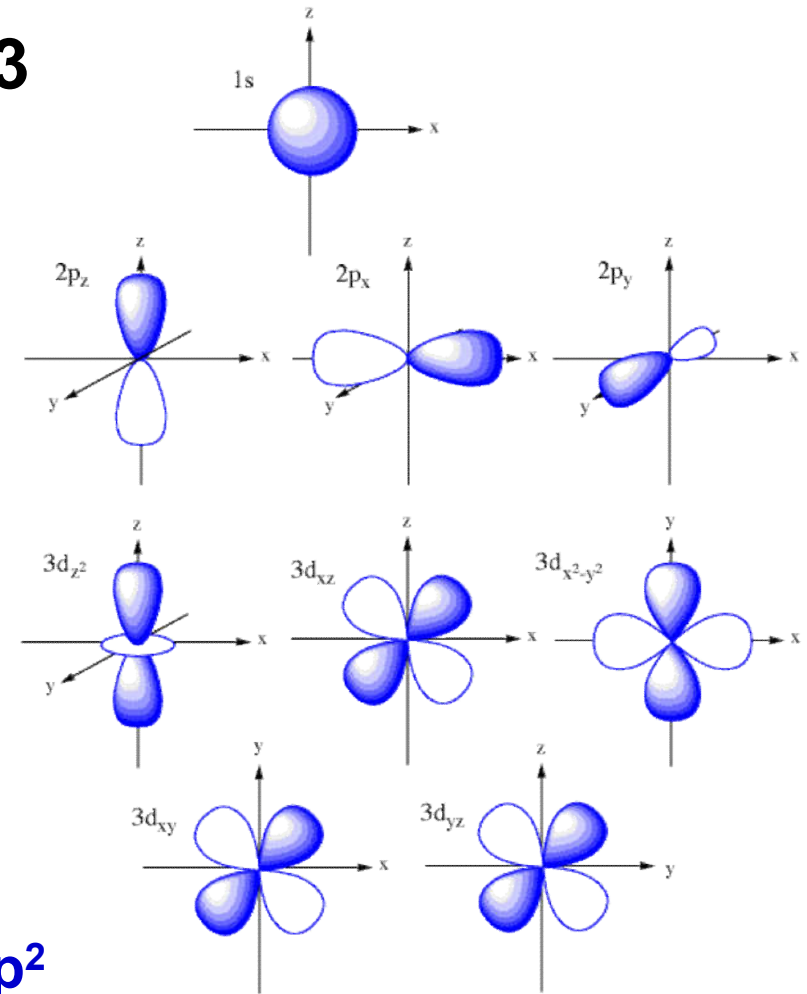
- Angular momentum: $l = 0, 1, 2, 3$
- Atomic orbitals: s p d f

- Examples

□ Hydrogen (H)	$1s^1$
□ Helium (He)	$1s^2$
□ Lithium (Li)	$[1s^2] 2s^1$
□ Carbon (C)	$[1s^2] 2s^2 2p^2$
□ Neon (Ne)	$[1s^2] 2s^2 2p^6$
□ Sodium (Na)	$[1s^2 2s^2 2p^6] 3s^1$
□ Silicon (Si)	$[1s^2 2s^2 2p^6] 3s^2 3p^2$

core electrons

valence electrons



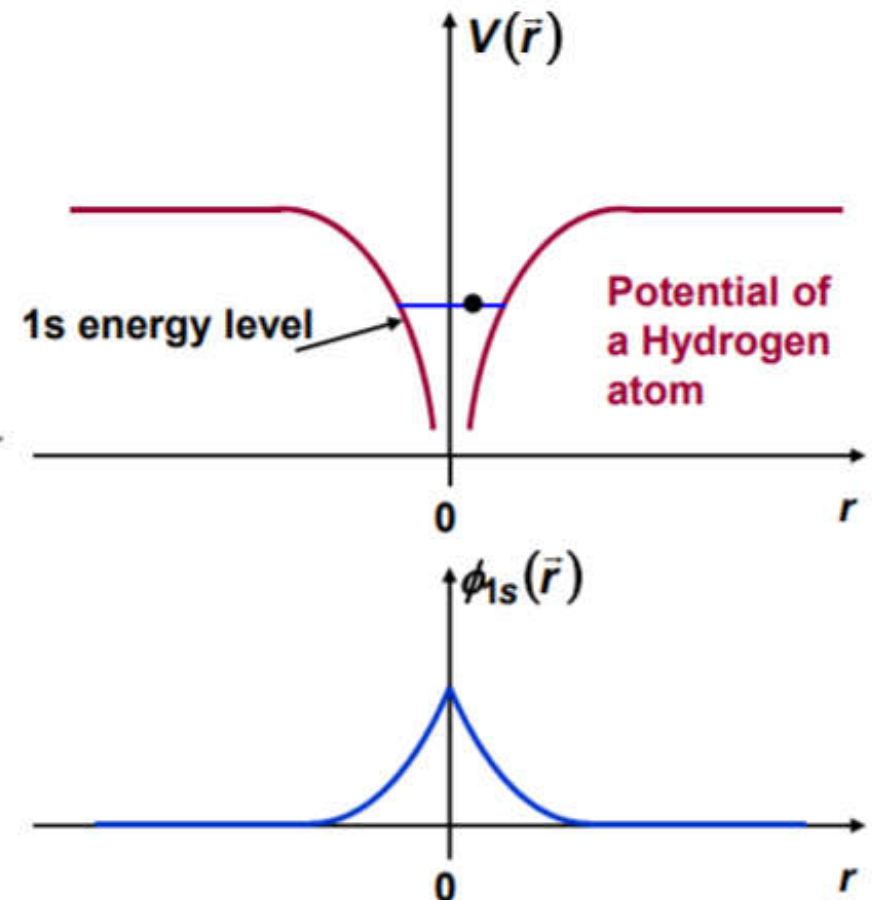
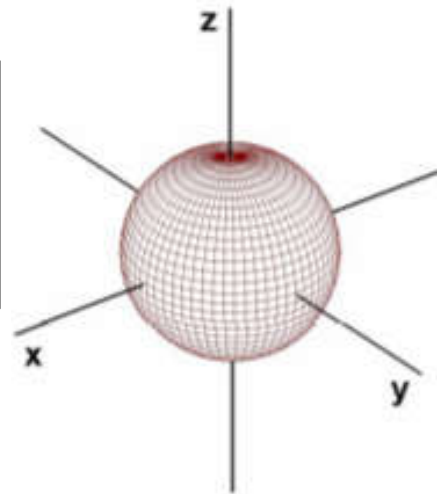
Hydrogen Atom

$$-\frac{\hbar^2}{2m} \nabla^2 \psi(\mathbf{r}) + V(\mathbf{r}) \cdot \psi(\mathbf{r}) = E \psi(\mathbf{r})$$

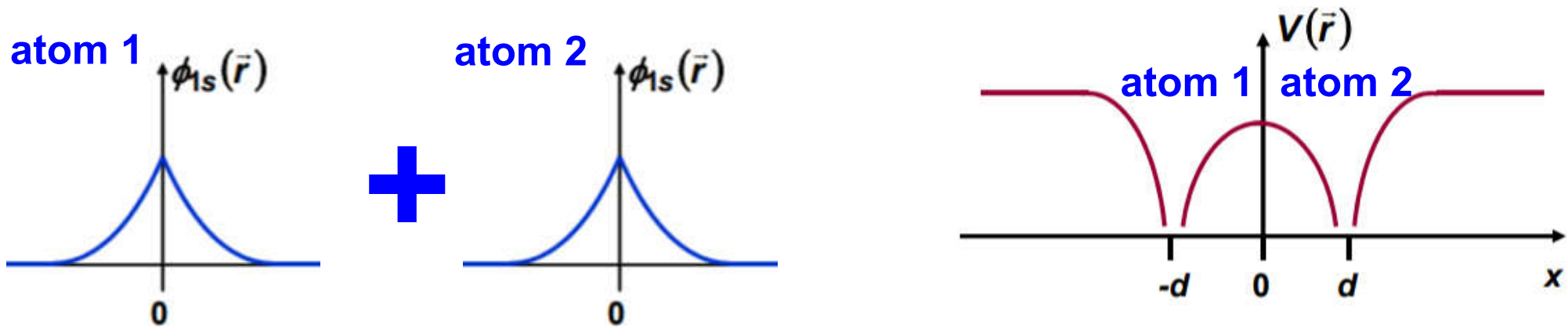
for 1s orbital →

$$\phi_{1s}(\mathbf{r}) = \frac{1}{\sqrt{\pi a_0^3}} e^{-r/a_0}$$

$$E_{1s} = -13.6 \text{ eV}$$



Hydrogen Molecule H-H



$$-\frac{\hbar^2}{2m} \nabla^2 \psi(\mathbf{r}) + [V_1(\mathbf{r}) + V_2(\mathbf{r})] \cdot \psi(\mathbf{r}) = E\psi(\mathbf{r})$$

Linear Combination of Atomic Orbitals (LCAO)

$$\psi(\mathbf{r}) = c_1 \phi_1(\mathbf{r}) + c_2 \phi_2(\mathbf{r})$$

Hydrogen Molecule H-H

Homework 5.5

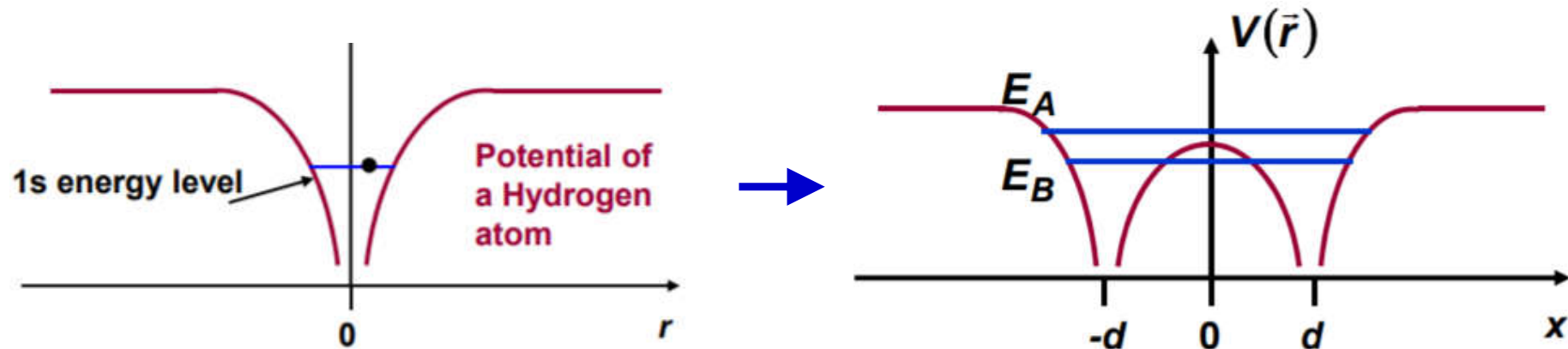
We have two solutions:

bonding molecular orbital (MO)

$$E_B = E_{1s} - V_{ss\sigma}$$

anti-bonding molecular orbital (MO)

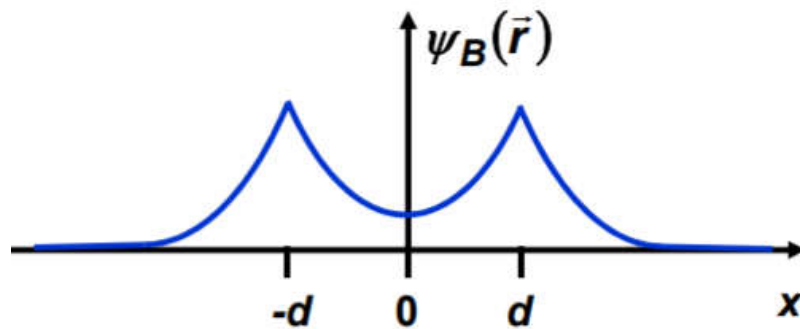
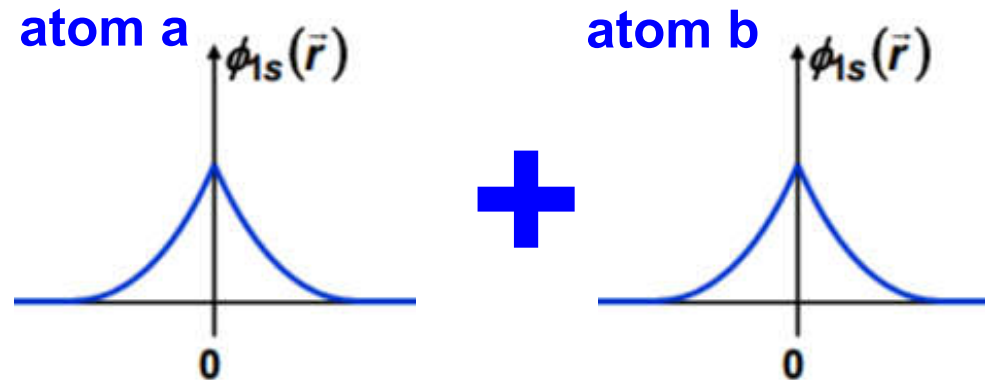
$$E_A = E_{1s} + V_{ss\sigma}$$



Pauli exclusion principle:

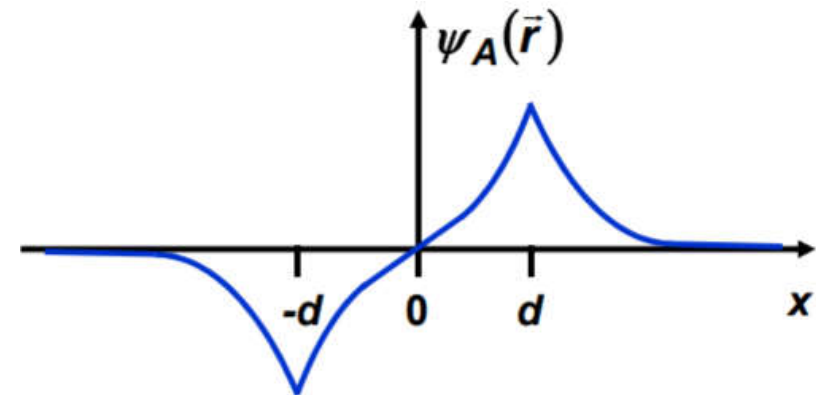
Two electrons cannot be in the same energy state

Hydrogen Molecule H-H



$$\psi_B(\mathbf{r}) \sim [\phi_1(\mathbf{r}) + \phi_2(\mathbf{r})]$$

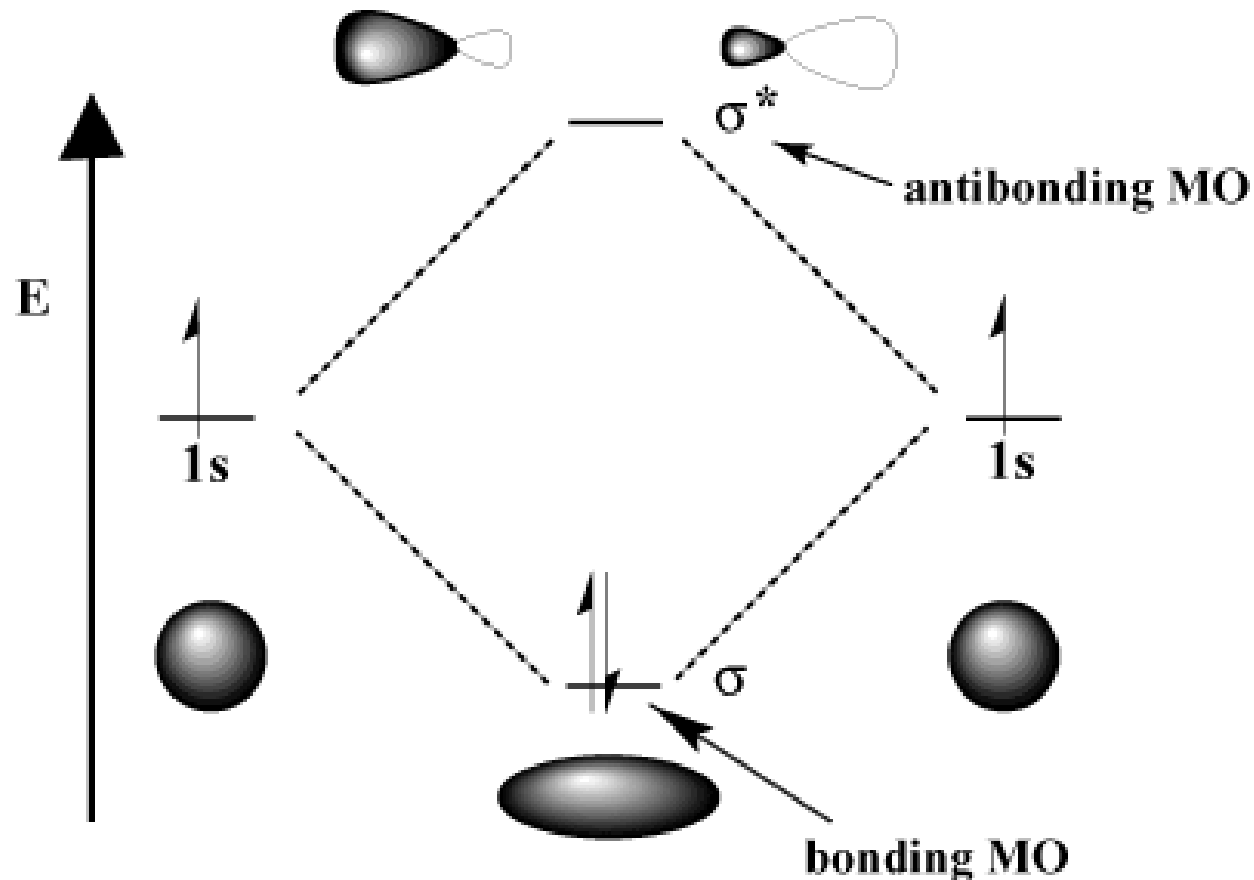
bonding orbital



$$\psi_A(\mathbf{r}) \sim [\phi_1(\mathbf{r}) - \phi_2(\mathbf{r})]$$

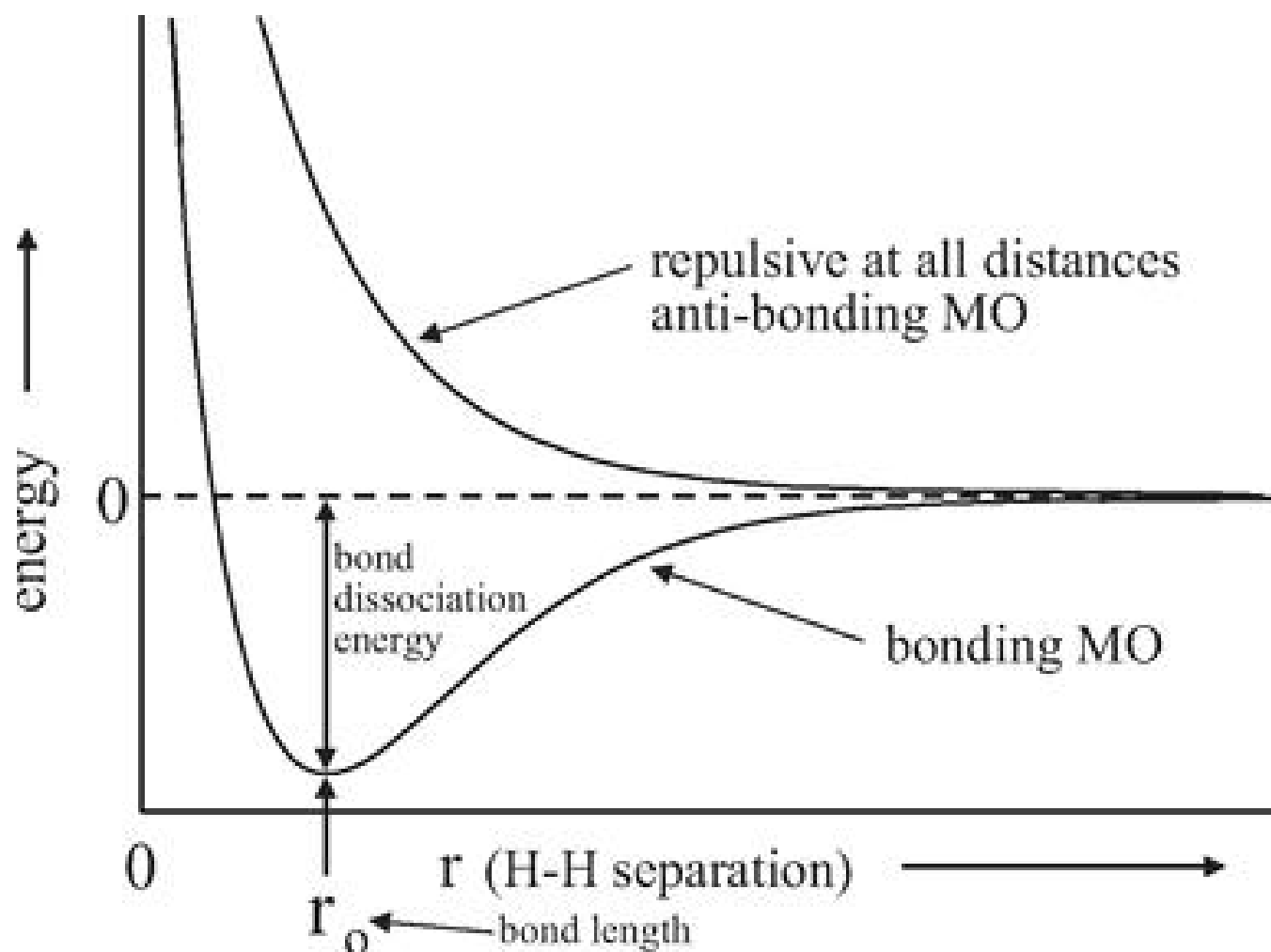
anti-bonding orbital

Hydrogen Molecule H-H



The two 1s orbitals on each Hydrogen atom combine to generate **two** molecular orbitals (MO): the **bonding MO** and the **anti-bonding MO**, with energy splitting

Hydrogen Molecule H-H



bonding energy vs. atom separation

A General Molecule A-B

Example: H-F, C-O, ...

Linear Combination of Atomic Orbitals (**LCAO**)

$$\psi(\mathbf{r}) = c_1\phi_1(\mathbf{r}) + c_2\phi_2(\mathbf{r})$$

For an electron, probabilities in A and B are different

$$P_A = \frac{c_1^2}{c_1^2 + c_2^2}$$

$$P_B = \frac{c_2^2}{c_1^2 + c_2^2}$$

A General Molecule A-B

Polarity (极性) f

$$f = \left| \frac{P_A - P_B}{P_A + P_B} \right| = \left| \frac{c_1^2 - c_2^2}{c_1^2 + c_2^2} \right|$$

For H-H, C-C, ...

$$|c_1| = |c_2| \longrightarrow f = 0 \quad \text{nonpolar bonding covalent (共价键)}$$

For H-F, C-H, Na-Cl, ...

$$|c_1| \neq |c_2| \longrightarrow 0 < f < 1 \quad \text{polar bonding covalent or ionic (离子键)}$$

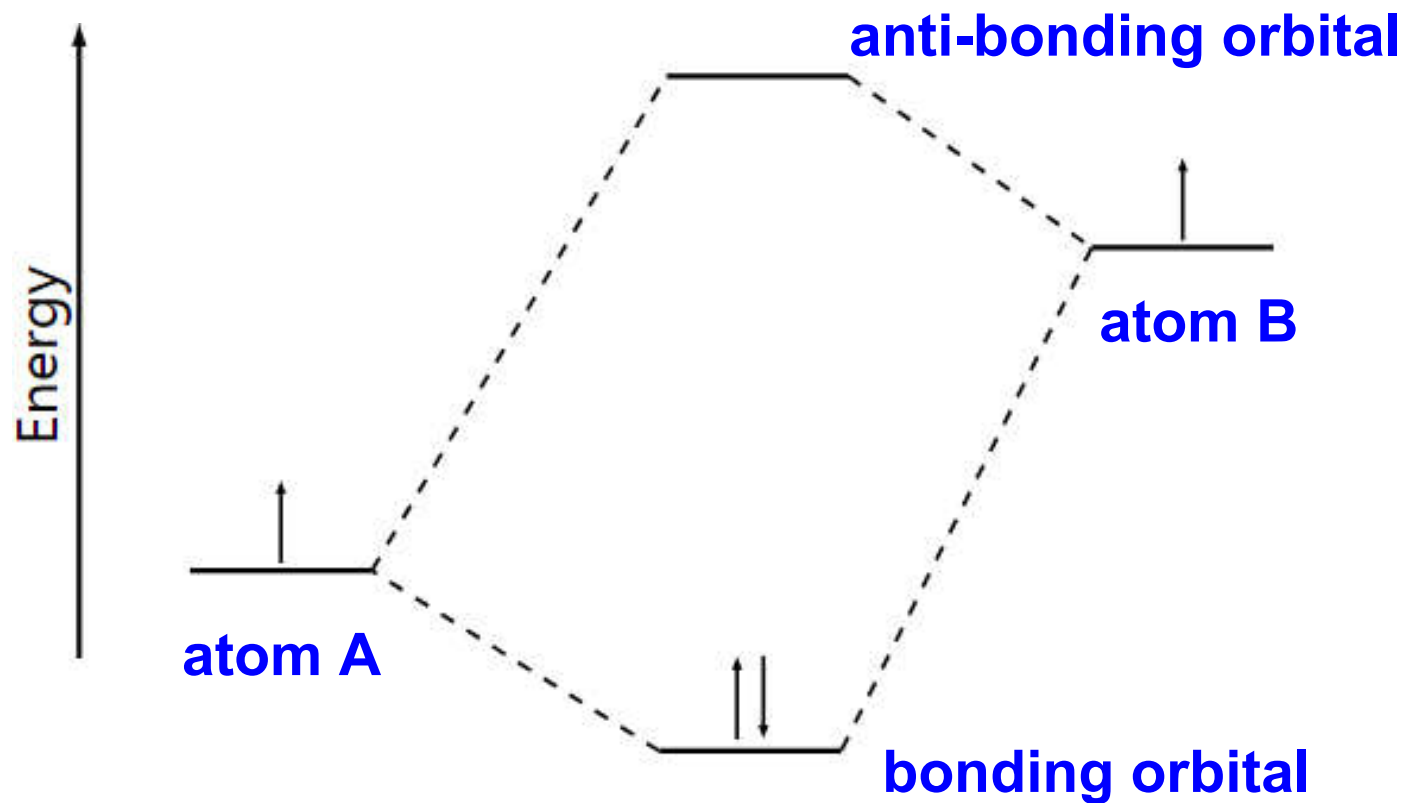
A General Molecule A-B

Polarity (极性) f

$$f = \left| \frac{P_A - P_B}{P_A + P_B} \right| = \left| \frac{c_1^2 - c_2^2}{c_1^2 + c_2^2} \right|$$

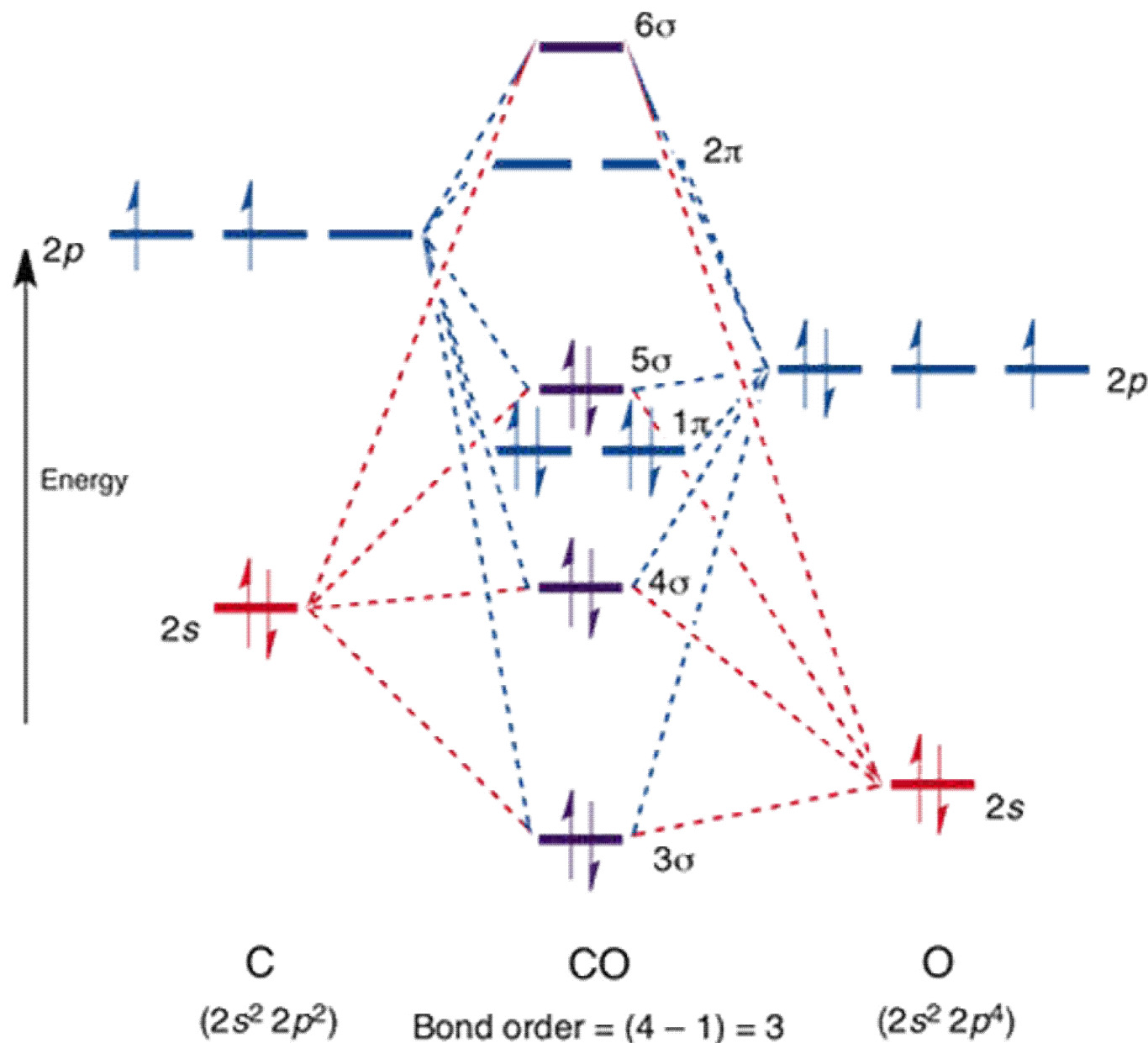
	C	Si	SiC	GaAs	ZnO
polarity f	0	0	0.177	0.310	0.616

A General Molecule A-B



Example: H-F, ...

Another Example: C-O bonding

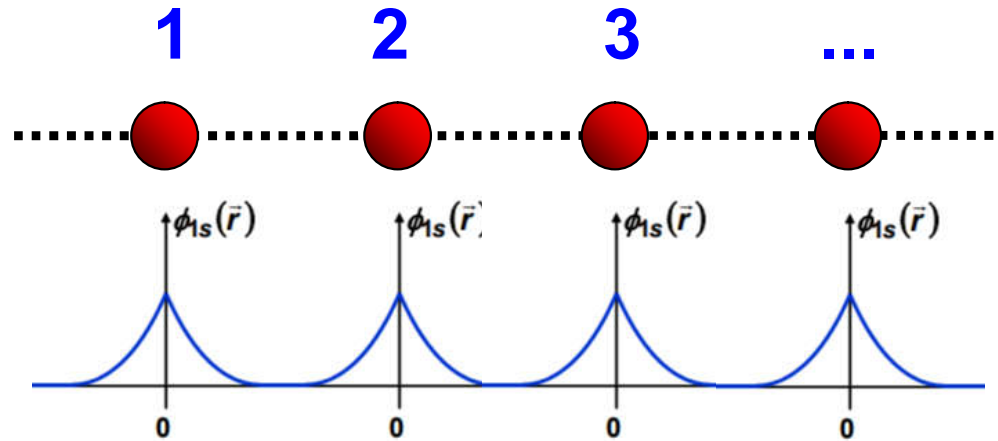


Chemical Bonding 化学键

- Metallic Bonding 金属键
- Ionic Bonding 离子键
- Covalent Bonding 共价键
- Van der Waals Bonding 范德华键
- Hydrogen Bonding 氢键
- ...

Chemical bonding originates from the electron wave functions distributed in multiple atoms

1D Chain of Atoms



Linear Combination of Atomic Orbitals (LCAO)

$$\psi(\mathbf{r}) = \sum_n c_n \phi_n(\mathbf{r})$$

use Bloch's Theorem

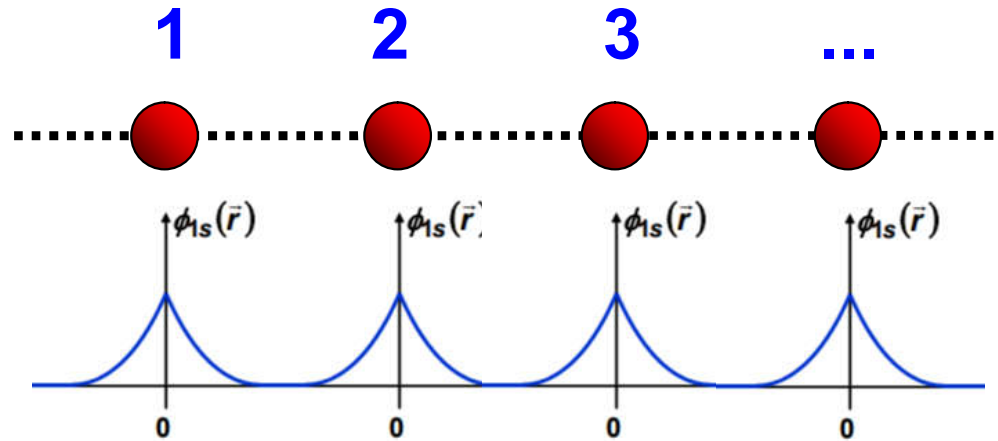


$$E(k) = E_{1s} - B - 2t \cos(ka)$$

$$k = \frac{2\pi}{a} \frac{n}{N}$$

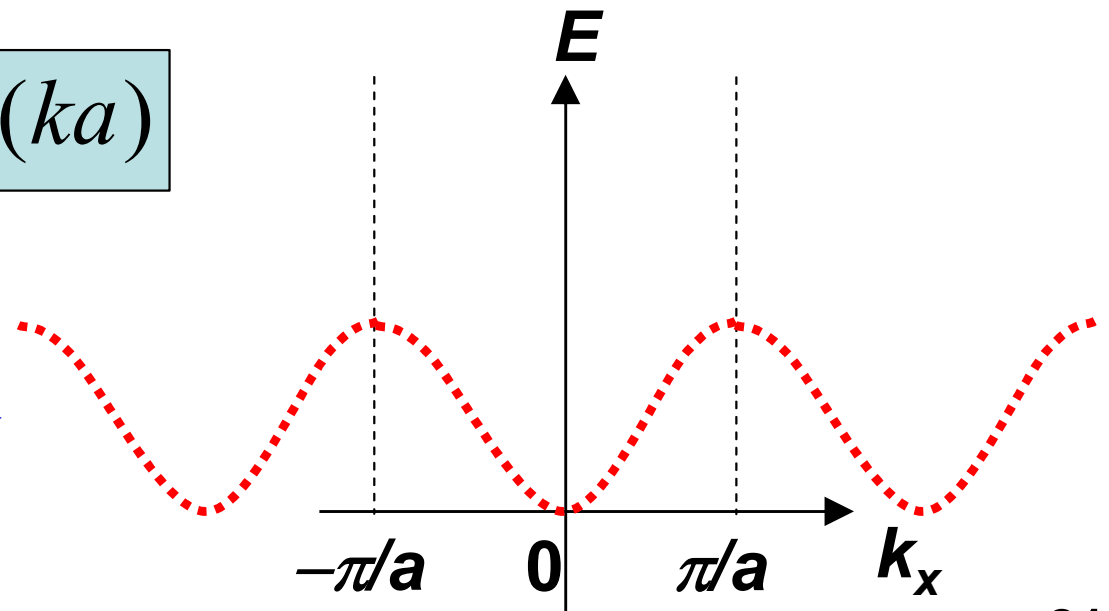
$$n = 0, \pm 1, \pm 2, \dots$$

1D Chain of Atoms



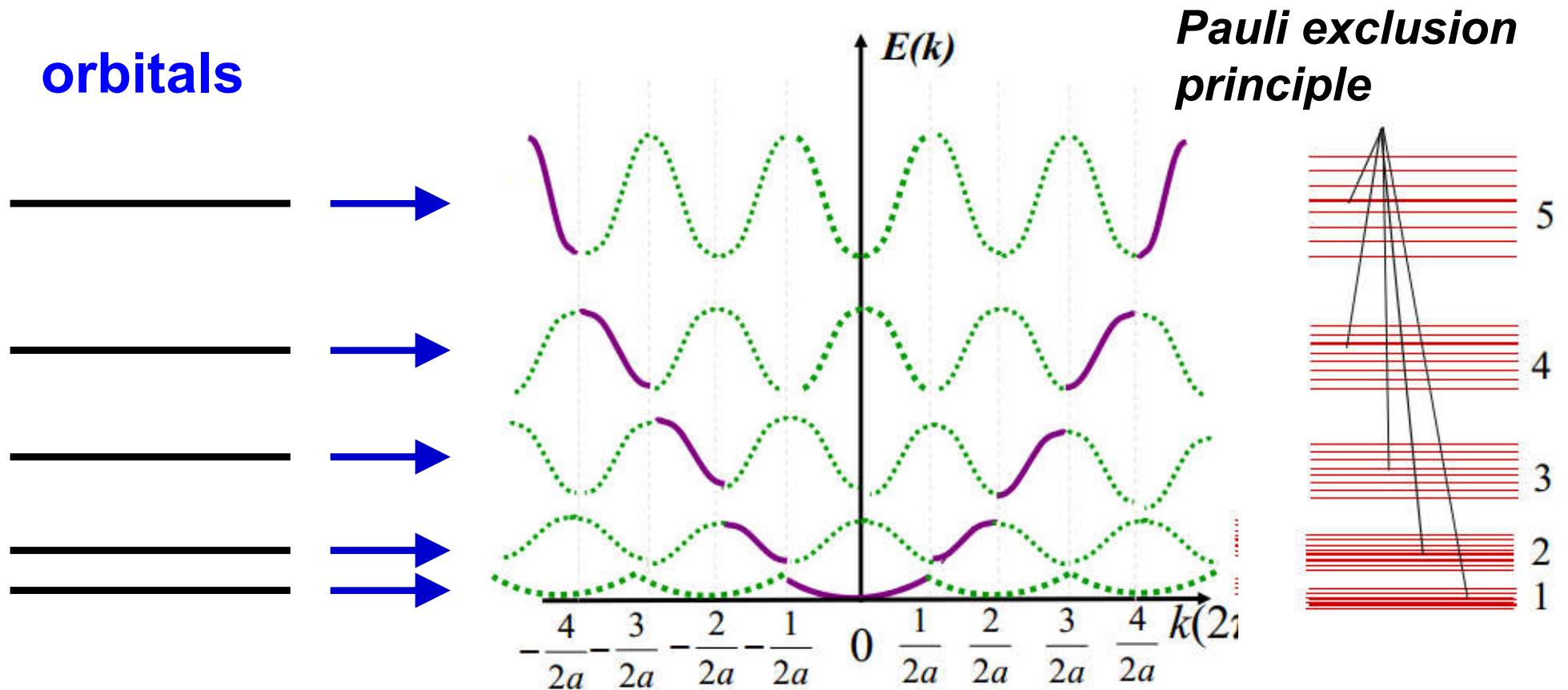
$$E(k) = E_{1s} - B - 2t \cos(ka)$$

1s orbital



Pauli exclusion principle

1D Chain of Atoms



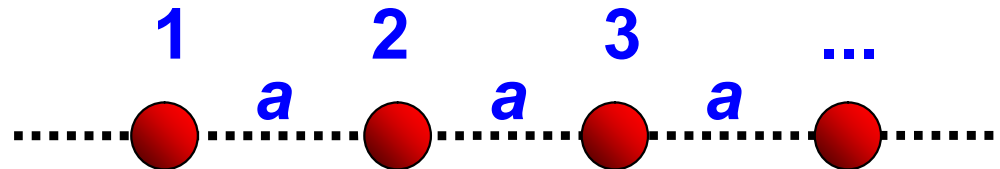
$$E(k) = E_{1s} - B - 2t \cos(ka)$$

$$k = \frac{2\pi}{a} \frac{n}{N}$$

$$n = 0, \pm 1, \pm 2, \dots$$

discrete orbitals become quasi-continuous bands

1D Chain of Atoms



$$E(k) = E_{1s} - B - 2t \cos(ka)$$

when ka is large

$$E(k) = \text{constant} \longrightarrow \text{discrete orbitals}$$

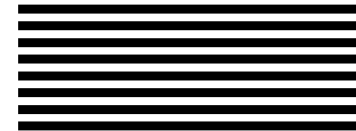
when ka is small

$$E(k) \approx E_{1s} - B - 2t + ta^2 k^2 \propto k^2$$

\longrightarrow nearly free electrons

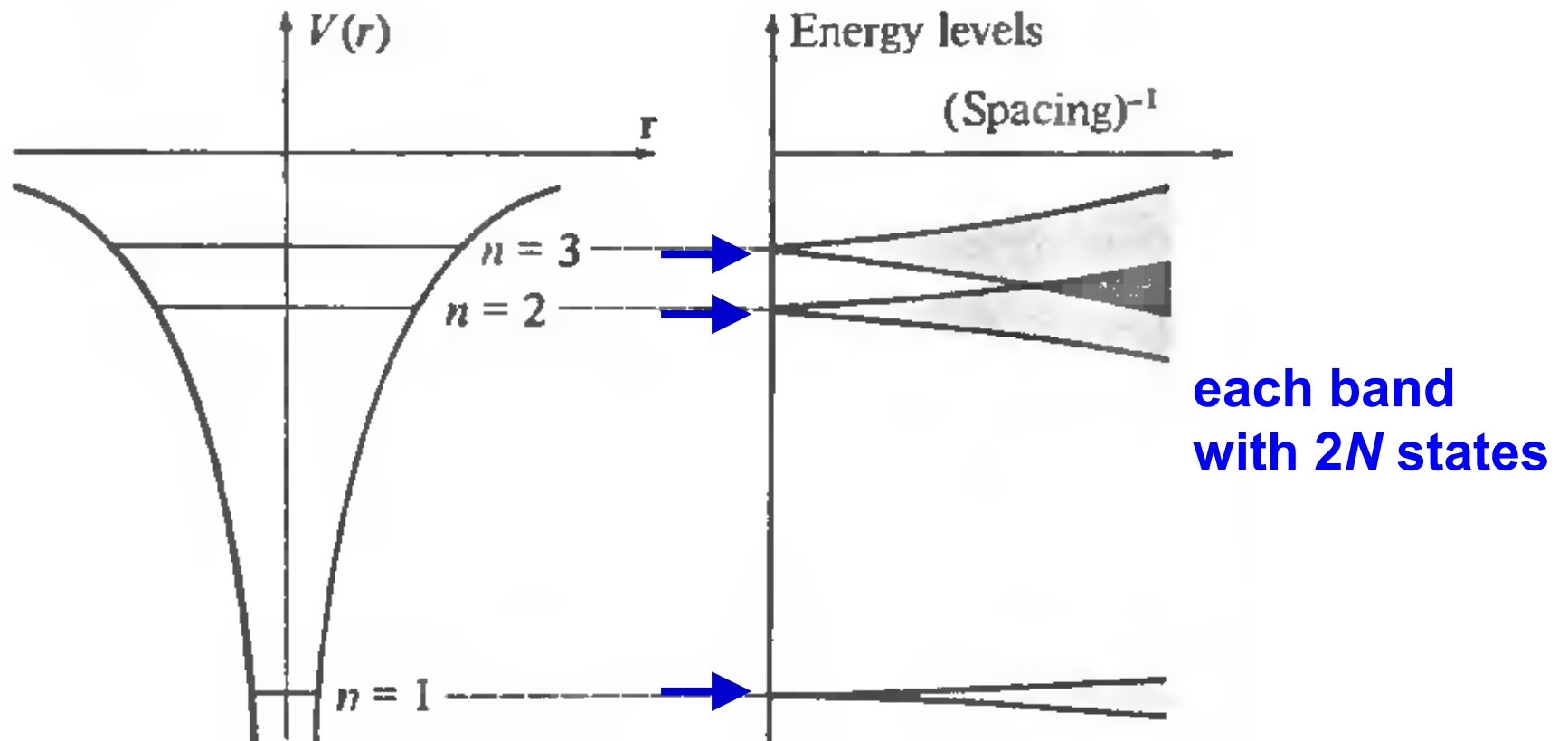
1D Chain of Atoms

1s orbital

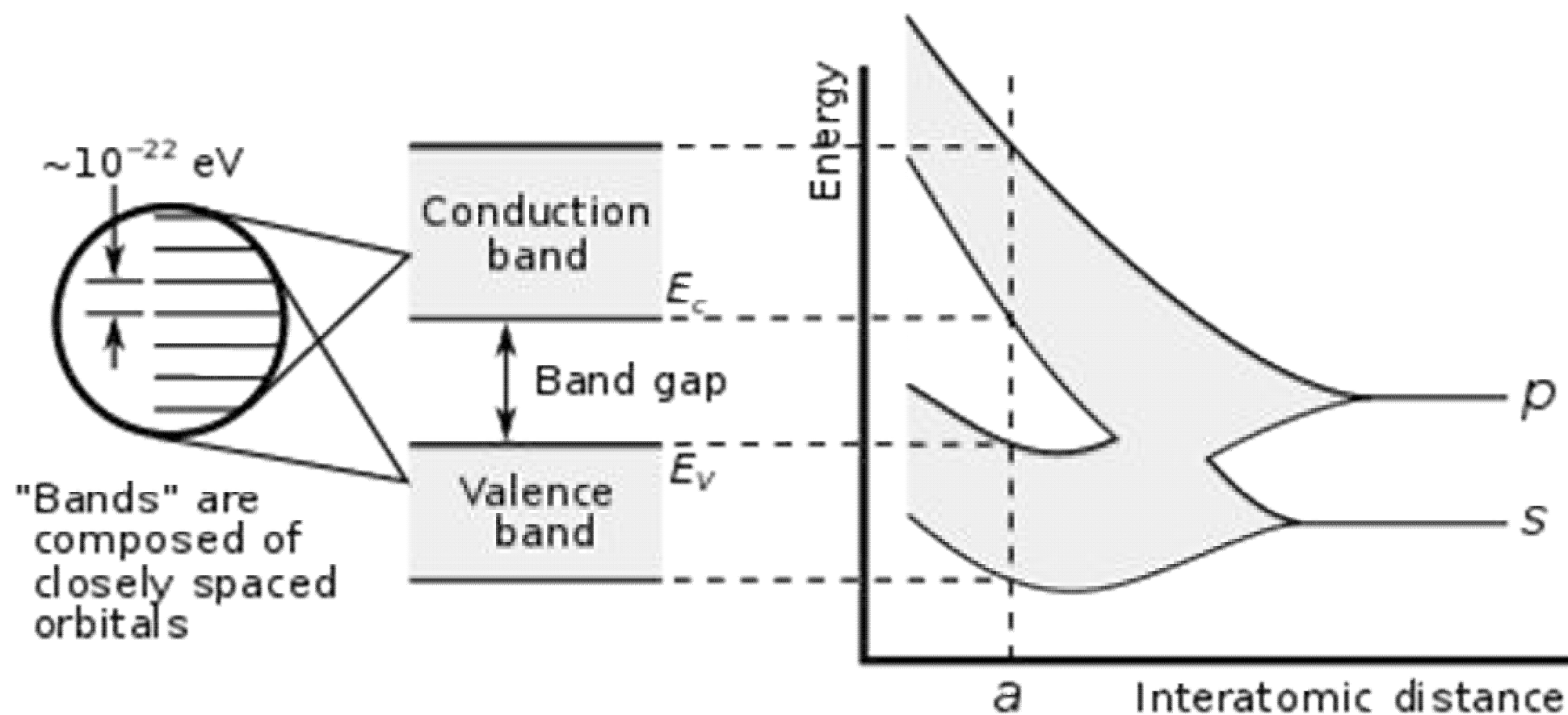


2N states

Pauli exclusion principle



Formation of bands and gaps



Thank you for your attention