

# 原子结构与键合

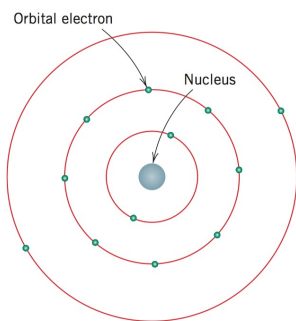
Dongsheng Wen

# 原子的结构

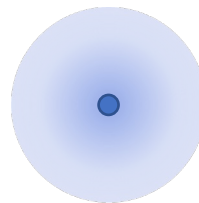
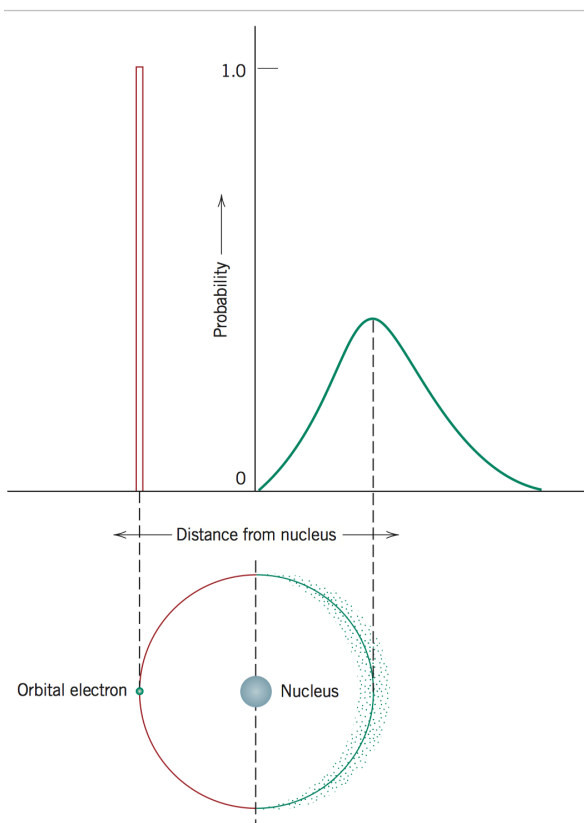
- 卢瑟福 (Rutherford) 的金箔实验 (1908-1913)
  - 三个人儿：Geiger, Marsden和Rutherford
  - 金箔
  - Alpha 粒子
  
- 结论

# 原子的结构

- 原子与核外电子（早期波尔模型-Bohr）

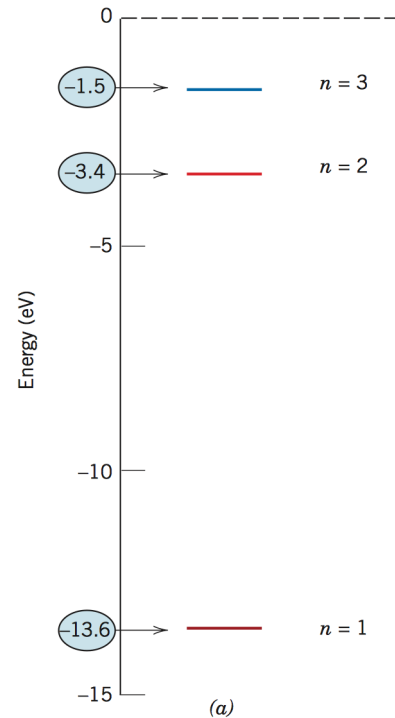
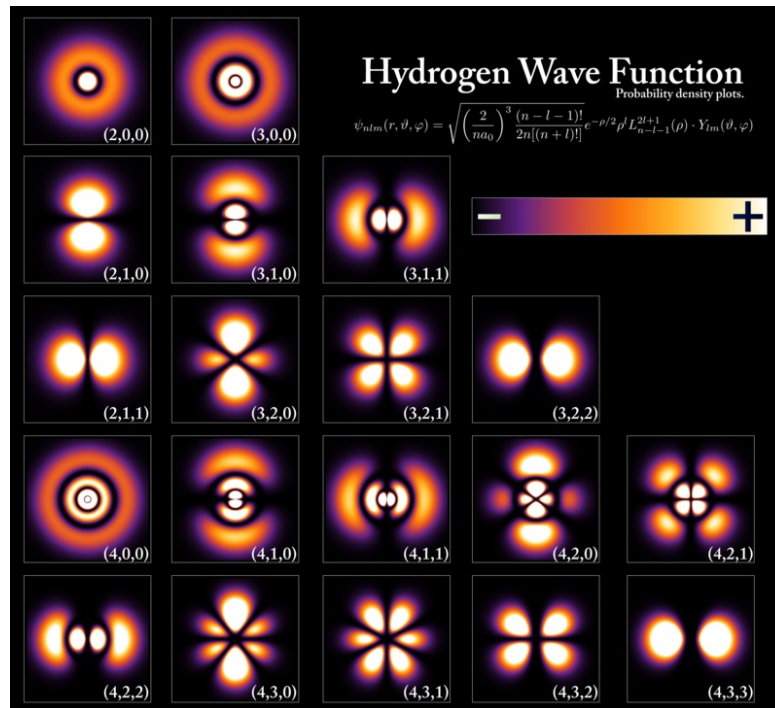
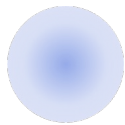


- 电子的波形态



# 原子的结构

- 电子轨道



[https://en.wikipedia.org/wiki/Hydrogen\\_atom](https://en.wikipedia.org/wiki/Hydrogen_atom)

# 原子的结构

- 电子轨道

IA		IIA												IIIA	IVA	VA	VIA	VIIA	0
1 H 1.0080		4 Be 9.0122												5 B 10.811	6 C 12.011	7 N 14.007	8 O 15.999	9 F 18.998	10 Ne 20.180
3 Li 6.941	12 Mg 24.305	IIIB		IVB	VB	VIB	VII	VIII		IB	IIB								
11 Na 22.990	20 Ca 40.08	21 Sc 44.956	22 Ti 47.87	23 V 50.942	24 Cr 51.996	25 Mn 54.938	26 Fe 55.845	27 Co 58.933	28 Ni 58.69	29 Cu 63.54	30 Zn 65.41	31 Ga 69.72	32 Ge 72.64	33 As 74.922	34 Se 78.96	35 Br 79.904	36 Kr 83.80		
19 K 39.098	20 Ca 40.08	21 Sc 44.956	22 Ti 47.87	23 V 50.942	24 Cr 51.996	25 Mn 54.938	26 Fe 55.845	27 Co 58.933	28 Ni 58.69	29 Cu 63.54	30 Zn 65.41	31 Ga 69.72	32 Ge 72.64	33 As 74.922	34 Se 78.96	35 Br 79.904	36 Kr 83.80		

Key

29	←	Atomic number
Cu	←	Symbol
63.54	←	Atomic weight

Metal

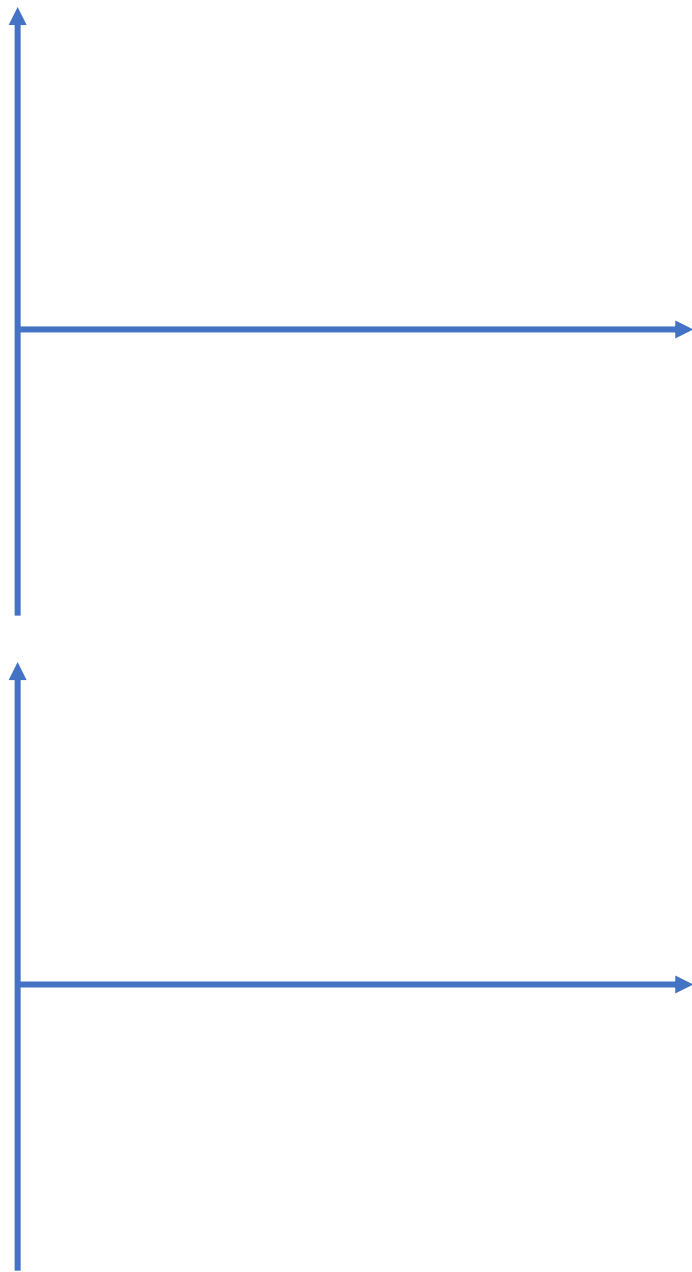
Nonmetal

Intermediate

# 原子间的力

- $F = F_A + F_R$

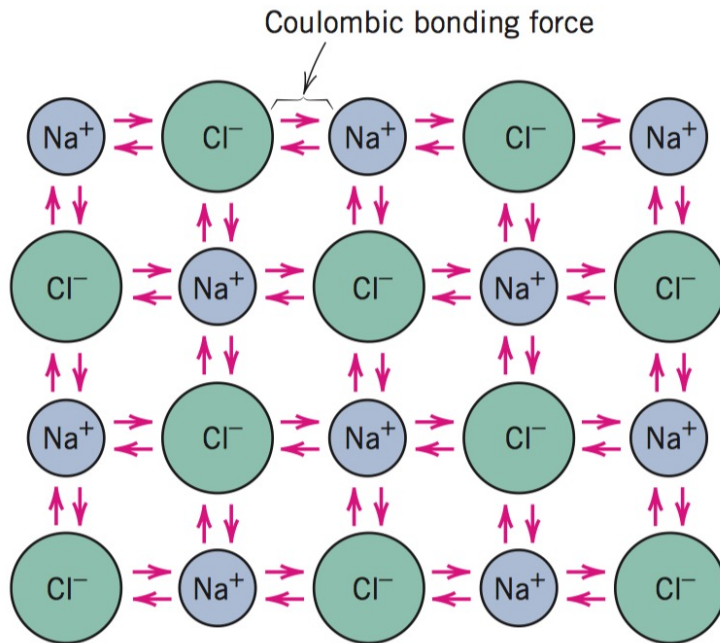
- $e = e_A + e_R$



# 化学键 (primary bonding)

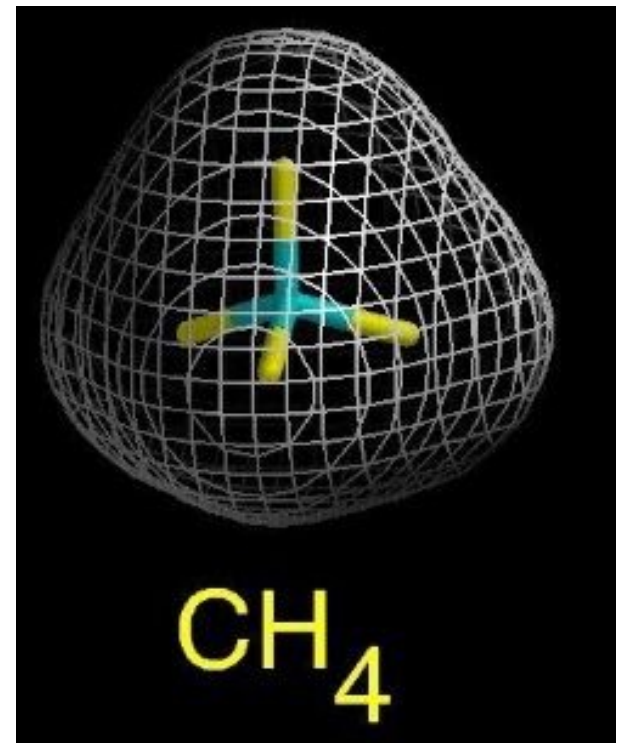
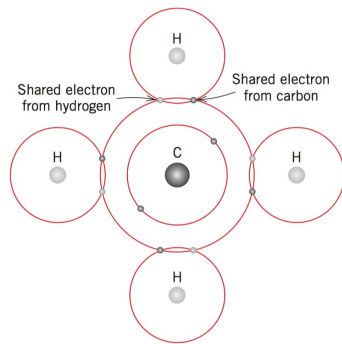
- 离子键 (ionic bonding)
- 共价键 (covalent bonding)
- 金属键 (metallic bonding)

# 离子键 (ionic bonding)

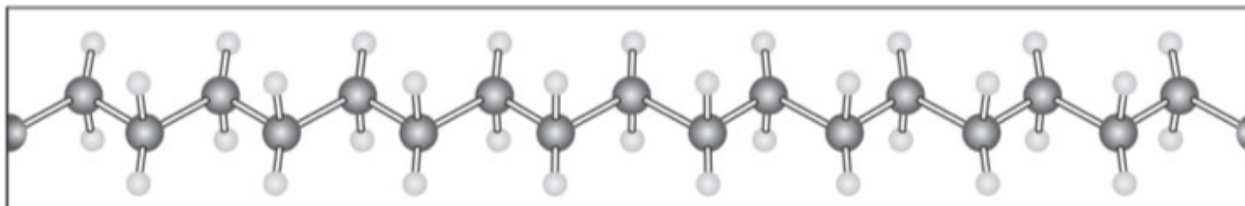
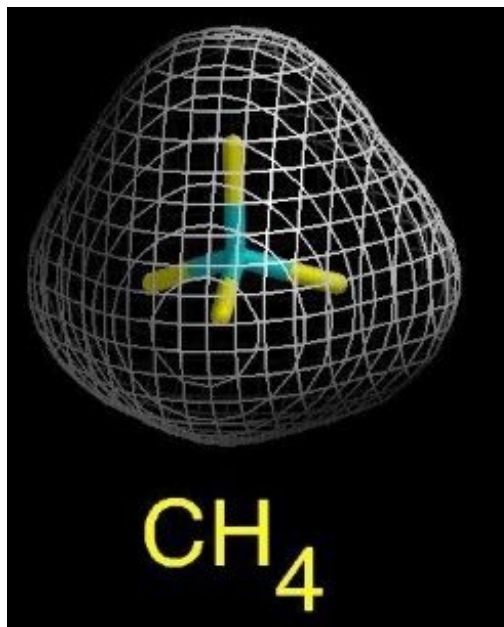
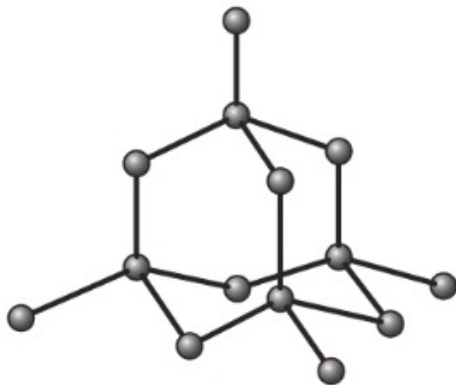




# 共价键 (covalent bonding)



# 共价键 (covalent bonding)



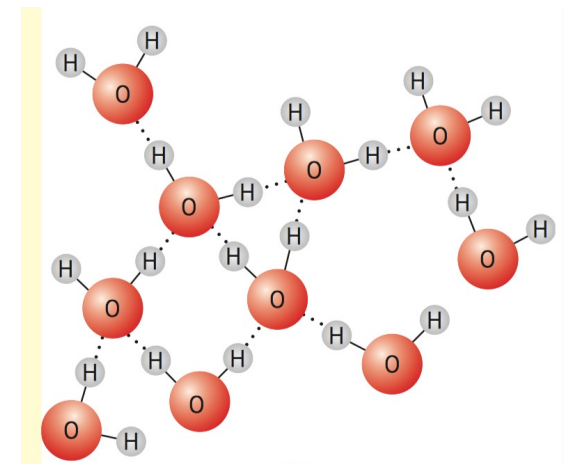
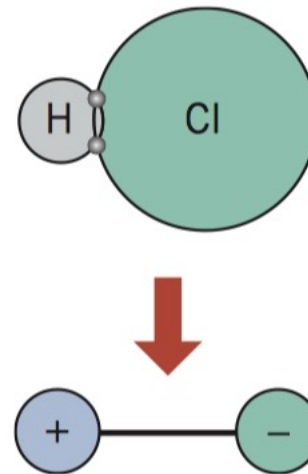
# 金属键 (metallic bonding)

# 物理键 (Secondary Bonding)

- 又称范德华键，比化学键要弱得多。



- 极化产生的范德华力



# 熔点与热膨胀

- $e = -\frac{A}{r^m} + \frac{B}{r^n}$



# 热膨胀

- $e = -\frac{A}{r^m} + \frac{B}{r^n}$



# 铝箔和塑料袋有啥不一样？



下一节课：晶向，晶面和点阵