

Ang. Mom. Addition → see review video

- rule: $j = |j_1 - j_2|, \dots, (j_1 + j_2)$
steps of 1
- reading CG tables: given $j_1 \times j_2 = |m_1, m_2\rangle \leftrightarrow |J, M\rangle$
individual composite
- q. nos vs e-values:

j	of $\hat{J}^2 = \hbar^2 j(j+1)$	$\neq M = -j, \dots, +j$ steps of 1
m	of $\hat{J}_z = \hbar m$	

Identical Particles

- Axiom 6: $\hat{P}_{ij} \psi(1, 2, \dots, i, \dots, j, \dots) = \begin{cases} -\psi_A(1, 2, \dots) & \text{if fermions} \\ +\psi_S(1, 2, \dots) & \text{if bosons} \end{cases}$
 exchange $i \leftrightarrow j$
 identical pcle

• $|\text{FULL QM STATE}\rangle = \text{space part} \times \text{spin part}$

$A =$ anti-symmetric
 $S =$ symmetric

↑
 this is what must obey Axiom 6

UNDER EXCHANGE
 of any pair ij
 of identical pcles

e.g. $|\text{STATE}\rangle_A$
 2 electrons →

$$\begin{cases} \psi_S(\vec{r}_1, \vec{r}_2) \cdot \chi_{A^{1,2}} \\ \psi_A(\vec{r}_1, \vec{r}_2) \cdot \chi_{S^{1,2}} \end{cases}$$

COROLLARY

▶ Pauli exclusion for identical fermions

▶ Exchange force: spatial w.f.

$\langle |\vec{r}_1 - \vec{r}_2| \rangle =$ larger for $\psi_A(\vec{r}_1, \vec{r}_2)$ = "repulsive"
 smaller for $\psi_S(\vec{r}_1, \vec{r}_2)$ = "attractive"

⊙ 2-spin 1/2 combos: (e.g. 2 electrons)

S=1 TRIPLET symmetric

$$\begin{cases} |1 +1\rangle_{SM} = |\uparrow\uparrow\rangle_{m_1 m_2} \\ |1 0\rangle_{SM} = \frac{1}{\sqrt{2}} [|\downarrow\uparrow\rangle + |\uparrow\downarrow\rangle]_{m_1 m_2} \\ |1 -1\rangle_{SM} = |\downarrow\downarrow\rangle_{m_1 m_2} \end{cases}$$

S=0 SINGLET antisymmetric

$$|0 0\rangle_{SM} = \frac{1}{\sqrt{2}} [|\downarrow\uparrow\rangle - |\uparrow\downarrow\rangle]_{m_1 m_2}$$

TECHNIQUE

⊙ Building ψ_A for > 2 particles: Slater determinant

⊙ 2-particle operators: use particle subscripts

e.g. $\hat{H}_{2 \text{ atomic electrons}} = -\frac{\hbar^2}{2m} \nabla_1^2 - \frac{\hbar^2}{2m} \nabla_2^2 + A \hat{S}_{1x} \hat{S}_{2x}$

acts on r_1, θ_1, ϕ_1 only

acts on x_2, y_2, z_2 only

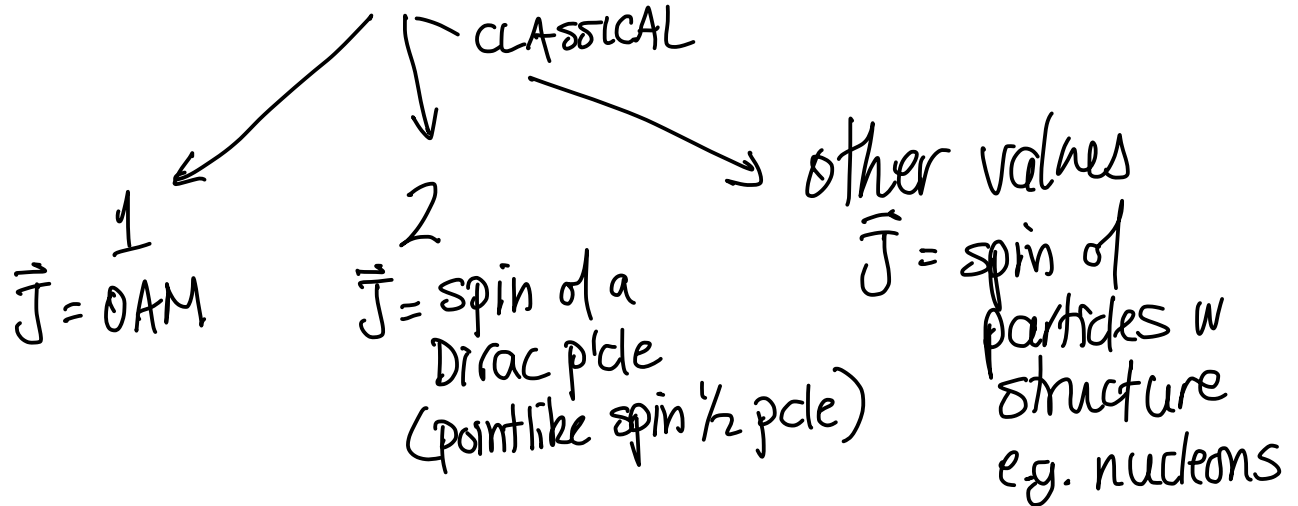
$$\begin{pmatrix} \hbar & 0 \\ \bar{z} & 0 \end{pmatrix} \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}$$

e.g. Disc 6 Q3

One- e^- Atoms: Bohr + spin-orbit

- Mag. Dipole Moment $\vec{\mu} = \gamma \vec{J}$

with $\gamma = g \cdot \left(\frac{e}{2m}\right)$



- Effect 1: $V_{\text{zeeman}} = -\vec{\mu} \cdot \vec{B}_{\text{EXTERNAL}}$

- Effect 2: $V_{\text{spin-orbit}} = -\vec{\mu} \cdot \vec{B}_{\text{INTERNAL}} \sim \hat{S} \cdot \hat{l}$
 (due to \vec{S} of atomic electron) (due to \vec{l} of electron around nucleus)

- good q. nos: when spin-orbit present,
 NOT (m_s, m_l) ... YES (j, m_j)

• atomic notation $\rightarrow l: s, p, d, f, \dots$

\rightarrow electronic config: $[\text{He}] 2s^2 2p^3$

\rightarrow term symbols: $2s+1 L_J$

• Hydrogenic energy corrections:
(in order)

Bohr

Fine Structure = spin-orbit +
Lamb shift [QED] relativistic KE

Hyperfine structure = $\vec{S} \cdot \vec{S}_{\text{NUC}}$

 Many-e Atoms : Bohr + V_{ee} + $V_{\text{spin-orbit}}$ +

• screening & central-force } approx to approximately
mean field } deal with V_{ee}

► filling order : $E_{nl} \sim (n+l)$

• closed shells : $S=0, L=0$

• good q. nos : L-S coupling scheme $\rightarrow S, L, J$
($V_{ee} \gg V_{s-o}$)

• Hund rules[#] (1, 2) \rightarrow minimize $V_{ee} \dots$ #3 \rightarrow min. V_{s-o}

e.g. energy-level diagram for Carbon: end of Lec 5B