

# Addition of Angular Momentum (aka Clebsch-Gordan Vector Coupling)

Remark: Recall from classical mechanics the connection between symmetry & conservation. If a coordinate  $q$  does not appear in Lagrangian ( $\frac{\partial L}{\partial q} = 0$ ) then the corresponding momentum is constant ( $\frac{\partial L}{\partial \dot{q}} = \text{const}$ ). In QM symmetry corresponds to  $[\hat{H}, Q] = 0 \Rightarrow$  degeneracy likely as  $\Psi$  eigenfunction  $\Rightarrow Q\Psi$  eigenfunction with same energy. Recall also the improved relation between  $\vec{L}$  & rotation - that  $\vec{L}$  "generates" rotations of the  $\vec{r}$  &  $\vec{p}$  that make it up.  $[\hat{H}, \vec{L}] = 0$  was start of program to show the "m degeneracy" -  $2\ell+1$  states with different orientation but the same  $L^2$

Note: the only thing we used in producing the  $|\ell m\rangle$  states and  $L_{\pm}$  operators was the commutation relations

(A)  $[\hat{L}_i, \hat{L}_j] = i\hbar \epsilon_{ijk} \hat{L}_k$ . If we discover new operators that follow that relations all the rest follows:  $2\ell+1$  eigenfunctions with  $L^2$  eigenvalue  $\hbar^2 \ell(\ell+1)$  &  $L_z$  eigenvalue  $m\hbar$

Consider orbital angular momentum operators  $\vec{L}$  & spin angular momentum operators  $\vec{S}$ . They operate on separate coordinates (CM location or Euler angle orientation) and so commute with each other. Separately they satisfy (A) so the sum  $\vec{J} = \vec{L} + \vec{S}$  also satisfies (A) and hence there must be a sequence of eigenfunctions of  $J^2$  with eigenvalue we call  $j$ :  $\hbar^2 j(j+1)$  &  $J_z$  eigenvalue  $\hbar m_j$ . For reasons I hope will be explained below - we want to find these eigenfunctions.

Consider two particles with coordinates  $\vec{r}_1$  &  $\vec{r}_2$ . There will be corresponding angular momentum operators,  $\vec{L}_1$  &  $\vec{L}_2$ . Since they operate on different coordinates they will commute; separately they satisfy (A), so the sum  $\vec{L} = \vec{L}_1 + \vec{L}_2$  also satisfies (A) and so there must exist a set of states with the same eigenvalue of  $L^2$  and different (stepwise) eigenvalues of  $L_z$

The previous examples aim to show that given two angular momentum operators we can form a sum angular momentum which must satisfy the rules/states generated by the raising/lowering operators. But why bother?

In the context of two orbiting particles (say in the atom) rotating, the coordinates of just one particle will generally not produce an equivalent problem - i.e.  $[H, \vec{L}_1] \neq 0$  but rotating the coordinates of every particle will leave an equivalent problem - i.e.  $[H, \vec{L}_1 + \vec{L}_2] = 0$  and in that case  $(L_1, L_2)_\pm$  will generate degenerate states whereas  $L_{1\pm}$  will not.

Remark: if we ignore the electrostatic interaction between the two electrons in the atom  $(\frac{e^2}{4\pi\epsilon_0 |\vec{r}_1 - \vec{r}_2|})$  then in fact  $[H, \vec{L}_1] = 0$

In the case of a single orbiting & spinning electron, if we rotate just  $\vec{L}$  (and leave  $\vec{S}$ ) the angle between  $\vec{L}$  &  $\vec{S}$  will change & it turns out the energy  $H$  does depend on the relative orientation of  $\vec{L}$  &  $\vec{S}$ :  $\vec{L} \cdot \vec{S}$  whereas if we rotate  $\vec{L}$  &  $\vec{S}$  by the same angle  $\vec{L} \cdot \vec{S}$  is invariant. This is a long way of saying  $[H, \vec{L}] \neq 0$  &  $[H, \vec{S}] \neq 0$  but  $[H, \vec{L} + \vec{S}] = 0$ , so the eigenfunctions of  $\vec{L} + \vec{S}$  will form a degenerate set.

In the case of a pair of isolated, spinning electrons we expect an interaction (i.e., one electron makes a  $\vec{B}$  the other then has  $PE = -\vec{\mu} \cdot \vec{B}$ ) But clearly that interaction depends only on the relative orientation of the spins i.e.  $\vec{S}_1 \cdot \vec{S}_2$ . Our dot product commutation rule then shows  $[\vec{S}_1 \cdot \vec{S}_2, \vec{S}_1 + \vec{S}_2] = 0$  so eigenfunctions of total spin will be degenerate.

Counter example: spinning electron in an external  $\vec{B}$ : rotating  $\vec{S}$  (but not  $\vec{B}$ ) changes energy &  $[H, \vec{S}] \neq 0$  degeneracy rules do not apply

Example: pair of isolated spin electrons with energy  $\vec{S}_1 \cdot \vec{S}_2$

Electrons are spin  $\frac{1}{2}$  so  $l=s=\frac{1}{2}$   $\therefore m_s = \pm\frac{1}{2}$  or  $-\frac{1}{2}$

notation  $|s m_s\rangle = |\frac{1}{2} \frac{1}{2}\rangle = \uparrow = \chi_+ \text{ text}$   
 $|\frac{1}{2} -\frac{1}{2}\rangle = \downarrow = \chi_- \text{ text}$

We have product wavefunctions where first arrow is always  $\uparrow$

Trick:  $\vec{S}_1 \cdot \vec{S}_2 = \underbrace{S_{1x}S_{2x} + S_{1y}S_{2y}}_{\text{compare to}} + S_{1z}S_{2z}$

$$\begin{aligned} S_{1+}S_{2-} + S_{1-}S_{2+} &= (S_{1x} + iS_{1y})(S_{2x} - iS_{2y}) + (S_{1x} - iS_{1y})(S_{2x} + iS_{2y}) \\ &= S_{1x}S_{2x} + S_{1y}S_{2y} + i(S_{1y}S_{2x} - S_{1x}S_{2y}) \\ &\quad + S_{1x}S_{2x} + S_{1y}S_{2y} - i(S_{1y}S_{2x} - S_{1x}S_{2y}) \\ &= 2(S_{1x}S_{2x} + S_{1y}S_{2y}) \end{aligned}$$

so  $\vec{S}_1 \cdot \vec{S}_2 = \frac{1}{2}(S_{1+}S_{2-} + S_{1-}S_{2+}) + S_{1z}S_{2z}$

general case:  $L_{\pm} |l m\rangle = \hbar \sqrt{l(l+1) - m(m\pm 1)} |l m\pm 1\rangle$

$S_+ \uparrow = 0$

$S_+ \downarrow = \hbar \sqrt{\frac{3}{4} - \frac{1}{2}(\frac{1}{2})} \uparrow = \hbar \uparrow$        $S_- \uparrow = \hbar \sqrt{\frac{3}{4} - \frac{1}{2}(-\frac{1}{2})} \downarrow = \hbar \downarrow$   
 $S_- \downarrow = 0$

We have 4 product states:  $\uparrow\uparrow, \uparrow\downarrow, \downarrow\uparrow, \downarrow\downarrow$

See how  $\vec{S}_1 \cdot \vec{S}_2$  affects each state

$[\frac{1}{2}(S_{1+}S_{2-} + S_{1-}S_{2+}) + S_{1z}S_{2z}] \uparrow\uparrow = \hbar^2 \frac{1}{4} \uparrow\uparrow \leftarrow \text{eigen function}$

$[\frac{1}{2}(S_{1+}S_{2-} + S_{1-}S_{2+}) + S_{1z}S_{2z}] \downarrow\downarrow = \hbar^2 \frac{1}{4} \downarrow\downarrow \leftarrow \text{degenerate eigen function}$

$[\frac{1}{2}(S_{1+}S_{2-} + S_{1-}S_{2+}) + S_{1z}S_{2z}] \uparrow\downarrow = \hbar^2 \left\{ \frac{1}{2} \downarrow\uparrow - \frac{1}{4} \uparrow\downarrow \right\}$

$[\frac{1}{2}(S_{1+}S_{2-} + S_{1-}S_{2+}) + S_{1z}S_{2z}] \downarrow\uparrow = \hbar^2 \left\{ -\frac{1}{4} \downarrow\uparrow + \frac{1}{2} \uparrow\downarrow \right\}$

see that  $[\ ] (\uparrow\downarrow + \downarrow\uparrow) = \hbar^2 \left\{ \frac{1}{4} \downarrow\uparrow + \frac{1}{4} \uparrow\downarrow \right\} \leftarrow \text{degenerate eigen function}$

see that  $[\ ] (\uparrow\downarrow - \downarrow\uparrow) = \hbar^2 \left\{ -\frac{3}{4} \uparrow\downarrow + \frac{3}{4} \downarrow\uparrow \right\}$

$\uparrow$   
eigenvalue =  $-\frac{3}{4} \hbar^2$

Faster: angular momentum addition:  $\frac{1}{2} + \frac{1}{2} = 0, 1$   
 $s = 0$   $s = 1$   $3x$  degenerate

$$\begin{aligned} \overline{S_1} \cdot \overline{S_2} &= \frac{1}{2} \left[ (\overline{S_1 + S_2})^2 - S_1^2 - S_2^2 \right] \\ &= \frac{\hbar^2}{2} \left[ s(s+1) - \frac{3}{4} - \frac{3}{4} \right] \end{aligned}$$

$s=1 \rightarrow +\frac{\hbar^2}{4}$   
 $s=0 \rightarrow -\frac{3}{4}\hbar^2$

Generalize: Given eigenstates of  $\vec{L}$   $|lm\rangle$  for given  $l$   
 Given eigenstates of  $\vec{S}$   $|sm_s\rangle$  for given  $s$   
 Construct product states  $|lm\rangle |sm_s\rangle$  and find combination of such states that is eigenfunction  $\vec{J} = \vec{L} + \vec{S}$

Note: since  $J_z |lm\rangle |sm_s\rangle = (L_z + S_z) |lm\rangle |sm_s\rangle$   
 $= \hbar (m + m_s) |lm\rangle |sm_s\rangle$   
 $\underbrace{m + m_s}_{m_j}$

we need only consider product state combinations that have  $m + m_s = \text{fixed}$

Classically we would expect total angular momentum to range from  $\vec{L} + \vec{S}$  to  $|\vec{L} - \vec{S}|$  - true in QM

Trick:  $J^2 = L^2 + S^2 + 2\vec{L} \cdot \vec{S} = L^2 + S^2 + 2L_z S_z + L_+ S_- + L_- S_+$   
 $\hbar^2 l(l+1)$   $\hbar^2 s(s+1)$   $m m_s \hbar^2$

Note: if we "start at the top"  $m=l$   $m_s=s$  discover it's already an eigenfunction of  $J^2$

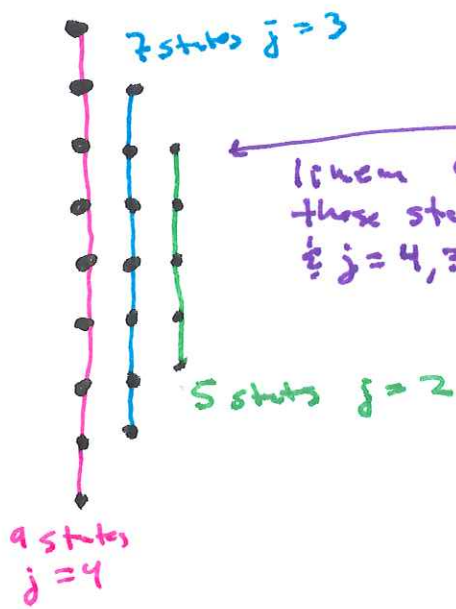
$$[L^2 + S^2 + 2L_z S_z + L_+ S_- + L_- S_+] |ll\rangle |ss\rangle = \hbar^2 \left\{ l(l+1) + s(s+1) + 2ls \right\} |ll\rangle |ss\rangle$$

$\leftarrow 0$   $\leftarrow 0$   $\leftarrow l(l+s+1) + s(l+s+1)$

we can then lower this state with  $(L_- + S_-)$  to get the linear combo  $|ll-1\rangle |ss\rangle$  &  $|ll\rangle |ss-1\rangle$  that has  $m_j = j-1$

the orthogonal combination of these two must have  $j = l+s-1$   
 $m_j = l+s-1$   
 etc.

Example:  $l=3$   $s=1$ . We label states just with  $m_l m_s$ . Make rows of states with same  $m_j = m_l + m_s$



- (3,1)
- (2,1) (3,0)
- (1,1) (2,0) (3,-1)
- (0,1) (1,0) (2,-1)
- (-1,1) (0,0) (1,-1)
- (-2,1) (-1,0) (0,-1)
- (-3,1) (-2,0) (-1,-1)
- (-3,0) (-2,-1)
- (-3,-1)

Note: the dimension of a vector space does not depend on basis:  $(2l+1)(2s+1)$  product states so must have same number of  $|j m_j\rangle$  states:

$$\sum_{j=l-s}^{l+s} 2j+1 = 2 \left( \frac{2s+1}{2} \right) (2l) + (2s+1) = (2s+1)(2l+1) \checkmark$$

← Gauss Trick

Notation:

$$|l s j m_j\rangle = \sum_{\substack{m_l + m_s = m_j \\ \text{fixed}}} \underbrace{|l m_l\rangle |s m_s\rangle}_{\text{product state}} \underbrace{\langle l s m_l m_s | l s j m_j \rangle}_{\text{Clebsch-Gordan}}$$

state with total angular momentum  $j$  is  $z$  component  $m_j$  made up of states with  $L \rightarrow l, S \rightarrow s$

Restatement: only angular momentum that is conserved will generate degeneracy. Often only total momentum is conserved.