## **Lecture 13 Highlights**

The eigenfunctions of  $J^2$  can be expressed as linear combinations of states with different values of  $m_\ell$  and  $m_s$  using the world-famous Clebsch-Gordan coefficients  $(C_{m_\ell - m_s - m_s}^{\ell - s - j})$  as:

$$\left|j m_{j}\right\rangle = \sum_{m_{\ell}+m_{s}=m_{j}} C_{m_{\ell} m_{s} m_{s} m_{j}}^{\ell s j} \left|\ell m_{\ell}\right\rangle \left|s m_{s}\right\rangle$$

Where the ket  $|\ell| m_{\ell}$  represents the spherical harmonics  $Y_{\ell}^{m_{\ell}}$ . The C-G coefficient values are given in Table 4.8 on page 188 of Griffiths. Remember that the all of the coefficients should appear under a square root, with the minus sign (if any) out front. Also note that we have dropped the radial part of the wavefunction  $(R_{n\ell})$  because it plays no role in combining angular momenta. Don't forget to put it back later.

Where do these coefficients come from? Consider starting with a product wavefunction at the top of the  $m_j$  ladder (it is a product of the wavefunctions with maximum values of  $m_\ell$  and  $m_s$ ). Now apply the  $J_-$ lowering operator, and construct orthonormal states on lower rungs of the ladder. The coefficients on the terms of those wavefunctions are the C-G coefficients.

We did a specific example of a hydrogen atom with  $\ell=1$  and spin s=1/2. In this case the angular momentum vector and spin vector can either be "parallel" or "antiparallel." Consider the two cases:

- 1) "Parallel"  $\vec{L}$  and  $\vec{S}$ : The maximum value of  $m_\ell$  is +1, while the value of  $m_s$  is +1/2 for the "parallel" case. This means that  $m_j = m_\ell + m_s = 3/2$ . This is the state at the top of the ladder. There must also be states with  $m_j = +1/2, -1/2, -3/2$ . This is a set of 4 states on the ladder of j = 3/2. Thus the eigenvalues of  $J^2$  for this ladder must be  $\frac{3}{2}(\frac{3}{2}+1)\hbar^2 = \frac{15}{4}\hbar^2$ . Note that  $\vec{L} \cdot \vec{S} > 0$  is this case, giving a positive spin-orbit Hamiltonian perturbation.
- 2) "Anti-Parallel"  $\vec{L}$  and  $\vec{S}$ : The maximum value of  $m_\ell$  is +1, while the value of  $m_s$  is -1/2 for the "anti-parallel" case. This means that  $m_j = m_\ell + m_s = 1/2$ . This is the state at the top of the ladder. There must also be a state with  $m_j = -1/2$ . This is a set of 2 states on the ladder of j = 1/2. Thus the eigenvalues of  $J^2$  for this ladder must be  $\frac{1}{2}(\frac{1}{2}+1)\hbar^2 = \frac{3}{4}\hbar^2$ . Note that  $\vec{L} \cdot \vec{S} < 0$  is this case, giving a negative spin-orbit Hamiltonian perturbation.

There are a total of 6 states possible by simply combining the orbital angular momentum with  $\ell = 1$  and spin angular momentum with s = 1/2! Just imagine what happens when you combine 3 or more angular momentum vectors!

Now for an example of how to construct states that are simultaneous eigenfunctions of  $L^2$ ,  $S^2$ ,  $J^2$  and  $J_z$ . Take the case again of hydrogen with  $\ell=1$  and spin s=1/2. How do we find the state with j=3/2 and  $m_j=-1/2$  in terms of the  $Y_\ell^{m_\ell}$  and spinors? Look at the  $1\times\frac{1}{2}$  Table on page 188. We are led to this table because we are combining an angular momentum vector with  $\ell=1$  and spin vector with  $\ell=1/2$ . Now look under the column labeled " $\frac{3/2}{-1/2}$ ". It says:

$$\left| \frac{3}{2} - \frac{1}{2} \right\rangle = \sum_{m_{\ell} + m_{s} = -1/2} C_{m_{\ell} - m_{s} - 1/2}^{1 - 1/2} \left| 1 m_{\ell} \right\rangle \left| \frac{1}{2} m_{s} \right\rangle$$

$$\left| \frac{3}{2} - \frac{1}{2} \right\rangle = \sqrt{\frac{2}{3}} \left| 1 0 \right\rangle \left| \frac{1}{2} - \frac{1}{2} \right\rangle + \sqrt{\frac{1}{3}} \left| 1 - 1 \right\rangle \left| \frac{1}{2} \frac{1}{2} \right\rangle$$

This can be written in a more familiar way in terms of spherical harmonics and spinors as:

$$\left|\frac{3}{2} - \frac{1}{2}\right\rangle = \sqrt{\frac{2}{3}}Y_1^0 \chi_- + \sqrt{\frac{1}{3}}Y_1^{-1} \chi_+$$

One can move back and forth between the coupled and un-coupled representations using the Clebsch-Gordan table on page 188. Here is the schematic layout for the CG table for combining two spins (called  $\vec{S}_1$ ,  $\vec{S}_2$ ) to form a total spin  $\vec{S} = \vec{S}_1 + \vec{S}_2$  ( $S^2$  has eigenvalue  $s(s+1)\hbar^2$ ):

General Schematic of the C-G Table

$$\vec{S} = \vec{S}_1 + \vec{S}_2$$
 Coupled Representation  $m_s$ 

$$m_{s_1}$$
  $m_{s_2}$   $CG\#$ 

Un-Coupled Representation

We considered what happens when two spin-1/2 spins are combined. The lecture followed Griffiths pages 184-188, although not in that order. We considered the 4 naïve product states of the two spins:

$$|\uparrow\rangle|\uparrow\rangle, \ |\uparrow\rangle|\downarrow\rangle, \ |\downarrow\rangle|\uparrow\rangle, \ |\downarrow\rangle|\downarrow\rangle$$
 where 
$$|\uparrow\rangle|\uparrow\rangle \text{ represents the product ket } |\frac{1}{2} + \frac{1}{2}\rangle_1|\frac{1}{2} + \frac{1}{2}\rangle_2, \text{ where the first }$$

number in each ket represents  $s_1$  and  $s_2$ , respectively, and the second number represents  $m_{s1}$  and  $m_{s2}$ . We found that the eigen-kets of the  $S^2$  operator (where  $\vec{S} = \vec{S}_1 + \vec{S}_2$ ) are in two ladders of states:

$$\begin{vmatrix} 1 & 1 \rangle = |\uparrow\rangle|\uparrow\rangle \\ |1 & 0 \rangle = \frac{1}{\sqrt{2}} (|\downarrow\rangle|\uparrow\rangle + |\uparrow\rangle|\downarrow\rangle)$$
 This is the s = 1 ladder of 3 states. **TRIPLET** 
$$|1 & -1\rangle = |\downarrow\rangle|\downarrow\rangle$$

and

$$|0 \quad 0\rangle = \frac{1}{\sqrt{2}} (|\downarrow\rangle|\uparrow\rangle - |\uparrow\rangle|\downarrow\rangle)$$
 This is the s = 0 ladder of 1 state. **SINGLET**

The kets on the left are written in the "coupled representation" while those on the right are in the "un-coupled representation." These are 4 orthonormal states that span the Hilbert space for the two spins.

Note that we started with spins that individually live on half-integer ladders, but the combined spin lives on integer ladders. This will have important ramifications for the physics of multi-particle systems, such as multi-electron atoms and the theory of superconductivity, among other things.