

Lecture 33: Quantum Mechanical Spin

Phy851 Fall 2009



Intrinsic Spin

- Empirically, we have found that most particles have an additional *internal* degree of freedom, called 'spin'
- The Stern-Gerlach experiment (1922):



- Each type of particle has a discrete number of internal states:
 - 2 states --> spin _
 - 3 states --> spin 1
 - Etc....



Interpretation

- It is best to think of spin as just an additional quantum number needed to specify the state of a particle.
 - Within the Dirac formalism, this is relatively simple and requires no new physical concepts
- The physical meaning of spin is not wellunderstood
- Fro Dirac eq. we find that for QM to be Lorentz invariant requires particles to have both anti-particles and spin.
- The 'spin' of a particle is a form of angular momentum



Spin Operators

• Spin is described by a vector operator:

$$\vec{S} = S_x \vec{e}_x + S_y \vec{e}_y + S_z \vec{e}_z$$

 The components satisfy angular momentum commutation relations:

$$[S_x, S_y] = i\hbar S_z$$
$$[S_y, S_z] = i\hbar S_x$$
$$[S_z, S_x] = i\hbar S_y$$

 This means simultaneous eigenstates of S² and S_z exist:

$$S^{2} = S_{x}^{2} + S_{y}^{2} + S_{z}^{2}$$
$$S^{2} | s, m_{s} \rangle = \hbar^{2} s (s+1) | s, m_{s} \rangle$$
$$S_{z} | s, m_{s} \rangle = \hbar m | s, m_{s} \rangle$$



Allowed quantum numbers

• For any set of 3 operators satisfying the angular momentum algebra, the allowed values of the quantum numbers are:

$$j \in \{0, \frac{1}{2}, 1, \frac{3}{2}, \dots\}$$
$$m_{j} \in \{-j, -j+1, \dots, j\}$$

- For orbital angular momentum, the allowed values were further restricted to only integer values by the requirement that the wavefunction be single-valued
- For spin, the quantum number, s, can only take on one value
 - The value depends on the type of particle
 - **S=0:** Higgs
 - *s*=1/2: Electrons, positrons, protons, neutrons, muons, neutrinos, quarks,...
 - **s=1:** Photons, W, Z, Gluon

- **s=2:** graviton

$$m_s \in \{-s, -s+1, \dots, s\}$$

Complete single particle basis

 A set of 5 commuting operators which describe the independent observables of a single particle are:

$$\vec{R}, S^2, S_z$$

- Or equivalently:

$$R, L^2, L_z, S^2, S_z$$

• Some possible basis choices:

$$\begin{cases} \vec{r}, s, m_s \\ \vec{p}, s, m_s \\ \end{cases} \\ \begin{cases} \vec{p}, s, m_s \\ \end{cases} \\ \begin{cases} r, \ell, m_\ell, s, m_s \\ \end{cases} \\ \end{cases} \\ \end{cases}$$

 When dealing with a single-particle, it is permissible to drop the s quantum number

Intrinsic Magnetic Dipole Moment

• Due to spin, an electron has an *intrinsic magnetic dipole moment*:

$$\vec{\mu}_e = -\frac{g_e|e|}{2m_e}\vec{S}$$

- g_e is the electron g-factor
- For an electron, we have:

$$g_e = 2.0023193043622 \pm 0.00000000015$$

- The is the most precisely measured physical quantity
- For most purposes, we can take g_e≈ 2, so that

$$\vec{\mu}_e = -\frac{\left|e\right|}{m_e}\vec{S}$$

• For any charged particle we have:

$$\vec{\mu} = \frac{g \ q}{2M} \vec{S}$$
Each particle
has a different
g-factor



Hamiltonian for an electron in a magnetic field

- Because the electron is a point-particle, the dipole-approximation is always valid for the spin degree of freedom
- Any `kinetic' energy associated with S² is absorbed into the rest mass
- To obtain the full Hamiltonian of an electron, we must add a single term: $\sqrt{2} = -\frac{1}{\mu} \cdot \frac{B}{B}$

$$H \rightarrow H + \frac{|e|}{m_e} \vec{S} \cdot \vec{B}(\vec{R})$$

$$H = \frac{1}{2m_e} \left[\vec{P} + |e|\vec{A}(\vec{R}) \right] - |e|\vec{O}(\vec{R}) + \frac{|e|}{m_e} \vec{S} \cdot \vec{B}(\vec{R})$$



Uniform Weak Magnetic Field as a perturbation

• For a weak uniform field, we find

$$H = \frac{P^2}{2m_e} + \frac{|e|B_0}{2m_e} (L_z + 2S_z)$$

• With the addition of a spherically symmetric potential, this gives:

$$H = \frac{P^2}{2m_e} + V(R) + \frac{|e|B_0}{2m_e} (L_z + 2S_z)$$

- If the zero-field eigenstates are known $H_0 | n, \ell, m_\ell \rangle = E_{0,n} | n, \ell, m_\ell \rangle$

- The weak-uniform-field eigenstates are: $\begin{cases}
n, \ell, m_{\ell}, m_{s} \rangle \\
H | n, \ell, m_{\ell}, m_{s} \rangle = E_{n, m_{\ell}, m_{s}} | n, \ell, m_{\ell}, m_{s} \rangle \\
E_{n, m_{\ell}, m_{s}} = E_{0, n} + \mu_{B} B_{0} \left(m_{\ell} + 2m_{s} \right) \\
\mu_{B} = \frac{|e|\hbar}{2m} \quad \text{'Bohr Magneton'}
\end{cases}$

Wavefunctions

- In Dirac notation, all spin does is add two extra quantum numbers
- The separate concept of a 'spinor' is unnecessary
- Coordinate basis: $\left\{ \vec{r}, S, m_s \right\}$

– Eigenstate of
$$ec{R},S^2,S_z$$

• Projector:

$$I = \sum_{m_s = -s}^{s} \int_{V} d^3 r |\vec{r}, s, m_s\rangle \langle \vec{r}, s, m_s |$$

• Wavefunction:

$$\psi_{m_s}\left(\vec{r}\right) := \left\langle \vec{r}, s, m_s \left| \psi \right\rangle \right.$$



Spinor Notation: $\psi_{m_s}(\vec{r}) \coloneqq \langle \vec{r}, s, m_s | \psi \rangle$

 We think of them as components of a length 2s+1 vector, where each component is a wavefunction

• Example:
$$s=1/2$$

 $\psi_{\uparrow}(\vec{r}) := \langle \vec{r}, s, \frac{1}{2} | \psi \rangle$
 $\psi_{\downarrow}(\vec{r}) := \langle \vec{r}, s, -\frac{1}{2} | \psi \rangle$

- Spinor wavefunction definition: $\begin{bmatrix} \psi \ \vec{r} \end{bmatrix} := \begin{pmatrix} \psi_{\uparrow} (\vec{r}) \\ \psi_{\downarrow} (\vec{r}) \end{pmatrix} = \begin{pmatrix} \psi_{\frac{1}{2}} (\vec{r}) \\ \psi_{-\frac{1}{2}} (\vec{r}) \end{pmatrix}$
 - If external and internal motions are not entangled, we can factorize the spinor wavefunction:

$$\left[\psi \right] \vec{r} := \begin{pmatrix} c_{\uparrow} \\ c_{\downarrow} \end{pmatrix} \psi \left(\vec{r} \right)$$

$$\begin{pmatrix} C_{\uparrow} \\ C_{\downarrow} \end{pmatrix} \begin{matrix} I \\ F \end{matrix}$$

Is then a *pure spinor*

Schrödinger's Equation

• We start from:

$$i\hbar \frac{d}{dt} |\psi\rangle = H |\psi\rangle$$

• Hit from left with with $\left< \vec{r}, m_s \right|$

$$i\hbar \frac{d}{dt} \langle \vec{r}, m_s | \psi \rangle = \langle \vec{r}, m_s | H | \psi \rangle$$

• Insert the projector

$$i\hbar \frac{d}{dt} \langle \vec{r}, m_s | \psi \rangle = \sum_{m'_s = -s}^{s} \int_{V} d^3 r' \langle \vec{r}, m_s | H | \vec{r}', m'_s \rangle \langle \vec{r}', m' | \psi \rangle$$

• Let:
$$H = \frac{P^2}{2M} \llbracket I \rrbracket + \llbracket V \rrbracket \vec{R}$$

- For s=1/2:

$$\begin{bmatrix}I\\\end{bmatrix} = \begin{pmatrix} 1 & 0\\ 0 & 1 \end{pmatrix} \quad \begin{bmatrix}V\\\end{bmatrix} \vec{r} = \begin{pmatrix} V_{\uparrow\uparrow}(\vec{r}) & V_{\uparrow\downarrow}(\vec{r})\\ V_{\downarrow\uparrow}(\vec{r}) & V_{\downarrow\downarrow}(\vec{r}) \end{pmatrix}$$

$$i\hbar\frac{d}{dt}\left[\psi\right]\vec{r} = -\frac{\hbar^2}{2M}\nabla^2\left[\psi\right]\vec{r} + \left[V\right]\vec{r}\psi\vec{r}$$

Example: Electron in a Uniform
Field

$$i\hbar \frac{d}{dt} [\psi](\vec{r}) = -\frac{\hbar^2}{2M} \nabla^2 [\psi](\vec{r}) + [[V]](\vec{r})(\psi](\vec{r}))$$

$$i\hbar\frac{d}{dt}\begin{pmatrix}\psi_{\uparrow}(\vec{r})\\\psi_{\downarrow}(\vec{r})\end{pmatrix} = \begin{bmatrix}-\frac{\hbar^{2}}{2m_{e}}\nabla^{2} - i\mu_{B}B_{0}\frac{\partial}{\partial\phi}\end{bmatrix}\begin{pmatrix}\psi_{\uparrow}(\vec{r})\\\psi_{\downarrow}(\vec{r})\end{pmatrix} + \mu_{B}B_{0}\begin{pmatrix}1&0\\0&-1\end{pmatrix}\begin{pmatrix}\psi_{\uparrow}(\vec{r})\\\psi_{\downarrow}(\vec{r})\end{pmatrix}$$

• This is just a representation of two separate equations:

$$i\hbar \frac{d}{dt} \psi_{\uparrow}(\vec{r}) = \left[-\frac{\hbar^2}{2m_e} \nabla^2 - i\mu_B B_0 \frac{\partial}{\partial\phi} \right] \psi_{\uparrow}(\vec{r}) + \mu_B B_0 \psi_{\uparrow}(\vec{r})$$
$$i\hbar \frac{d}{dt} \psi_{\downarrow}(\vec{r}) = \left[-\frac{\hbar^2}{2m_e} \nabla^2 - i\mu_B B_0 \frac{\partial}{\partial\phi} \right] \psi_{\downarrow}(\vec{r}) - \mu_B B_0 \psi_{\downarrow}(\vec{r})$$

 We would have arrived at these same equations using Dirac notation, without ever mentioning 'Spinors'



$$\frac{\text{Pauli Matrices}}{i\hbar \frac{d}{dt} \begin{pmatrix} \psi_{\uparrow}(\vec{r}) \\ \psi_{\downarrow}(\vec{r}) \end{pmatrix}} = \left[-\frac{\hbar^2}{2m_e} \nabla^2 - i\mu_B B_0 \frac{\partial}{\partial \phi} \right] \begin{pmatrix} \psi_{\uparrow}(\vec{r}) \\ \psi_{\downarrow}(\vec{r}) \end{pmatrix} + \mu_B B_0 \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} \begin{pmatrix} \psi_{\uparrow}(\vec{r}) \\ \psi_{\downarrow}(\vec{r}) \end{pmatrix}$$

 Here we see that we have recovered one of the Pauli Matrices:

$$\sigma_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$$

• The other Pauli matrices are:

$$\sigma_x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \qquad \sigma_y = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}$$

• Then in the basis of eigenstates of S_z we have: $\neg \hbar$

$$\vec{S} = \frac{n}{2}\vec{\sigma}$$

$$i\hbar\frac{d}{dt}\left[\psi\right]\vec{r}\left[\vec{r}\right] = \left(-\frac{\hbar^2}{2M}\nabla^2 + \mu_B B_0\left(-i\frac{\partial}{\partial\phi} + \sigma_z\right)\right)\left[\psi\right]\vec{r}\right)$$

• If we only care about spin dynamics: $i \frac{d}{dt} [c] = \frac{|e|}{2m_e} \vec{\sigma} \cdot \vec{B} [c]$

Two particles with spin

 How do we treat a system of two particles with masses M₁ and M₂, charges q₁ and q₂, and spins s₁ and s₂?

- Basis:

$$\left| \vec{r}_{1}, s_{1}, m_{s1}; \vec{r}_{2}, s_{2}, m_{s2} \right\rangle$$

- Wavefunction:

$$\psi_{s_1,m_{s_1},s_2,m_{s_2}}(\vec{r}_1,\vec{r}_2) := \langle \vec{r}_1,s_1,m_{s_1};\vec{r}_2,s_2,m_{s_2} | \psi \rangle$$

 Hamiltonian w/out motional degrees of freedom:

$$H = -\frac{q_1}{M_1} \vec{S}_1 \cdot \vec{B}(\vec{R}_1) - \frac{q_2}{M_2} \vec{S}_2 \cdot \vec{B}(\vec{R}_2)$$

 Hamiltonian w/ motional degrees of freedom:

$$H = \frac{1}{2M_1} \left(\vec{P}_1 - q_1 \vec{A}(\vec{R}_1) \right) + q_1 \Phi(\vec{R}_1) - \frac{q_1}{M_1} \vec{S}_1 \cdot \vec{B}(\vec{R}_1) = \frac{1}{2M_2} \left(\vec{P}_2 - q_2 \vec{A}(\vec{R}_2) \right) + q_2 \Phi(\vec{R}_2) - \frac{q_2}{M_2} \vec{S}_2 \cdot \vec{B}(\vec{R}_2)$$

Example #1

 A spin _ particle is in the ↑ state with respect to the z-axis. What is the probability of finding it in the ↓-state with respect to the x-axis?

• Let:
$$|\psi\rangle = |\uparrow_z\rangle$$

• In the basis, $\{\uparrow_z\rangle, |\downarrow_z\rangle\}$ the operator for the x-component of spin is:

$$\sigma_x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}$$

- By symmetry, $\sigma_{\rm X}$ must have eigenvalues +1 and -1
- The eigenvector corresponding to -1 is defined by:

$$\sigma_{x} |\downarrow_{x}\rangle = - |\downarrow_{x}\rangle$$



Example #1 continued:

$$\sigma_{x} = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \qquad \sigma_{x} | \downarrow_{x} \rangle = - | \downarrow_{x} \rangle$$

$$| \downarrow_{x} \rangle = -\sigma_{x} | \downarrow_{x} \rangle$$

$$\langle \uparrow_{z} | \downarrow_{x} \rangle = - \langle \uparrow_{z} | \sigma_{x} | \downarrow_{x} \rangle$$

$$= - \langle \downarrow_{z} | \downarrow_{x} \rangle$$

• This implies that:

$$\left|\downarrow_{x}\right\rangle = \frac{1}{\sqrt{2}} \left(\uparrow_{z}\right) - \left|\downarrow_{z}\right\rangle\right)$$

$$P = \left| \left\langle \downarrow_x \right| \uparrow_z \right\rangle \right|^2 = \frac{1}{2}$$



Example #2

 Two identical spin-1/2 particles are placed in a uniform magnetic field. Ignoring motional degrees of freedom, what are the energy-levels and degeneracies of the system?

• States:
$$\{\uparrow\uparrow\rangle, |\uparrow\downarrow\rangle, |\downarrow\uparrow\rangle, |\downarrow\downarrow\rangle\}$$

- Z-axis chosen along B-field

• Hamiltonian:
$$H = -\frac{gqB_0}{2M} \left(S_{1z} + S_{2z}\right)$$

• Basis states are already eigenstates:

$$H|\uparrow\uparrow\rangle = -\frac{\hbar g q B_0}{2M}|\uparrow\uparrow\rangle \qquad E_1 = -\frac{\hbar g q B_0}{2M}; \quad d_1 = 1$$
$$H|\uparrow\downarrow\rangle = H|\downarrow\uparrow\rangle = 0 \qquad E_2 = 0; \quad d_2 = 2$$
$$H|\downarrow\downarrow\rangle = \frac{\hbar g q B_0}{2M}|\downarrow\downarrow\rangle \qquad E_3 = \frac{\hbar g q B_0}{2M}; \quad d_3 = 1$$