• Fine structure

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Homework Assignment #04: Chapter 7:8,12,13,17,33,36 due Tuesday, February 11, 2020

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Homework Assignment #04: Chapter 7:8,12,13,17,33,36 due Tuesday, February 11, 2020

Homework Assignment #05: Chapter 7:20,21,24,28,29,37 due Tuesday, February 18, 2020

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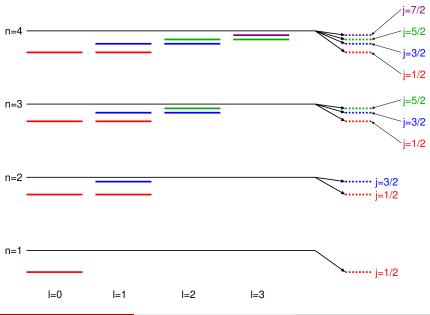
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Fine structure of hydrogen



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$$B_{\rm ext} \ll B_{\rm int}$$
 weak-field

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 $B_{\text{ext}} \ll B_{\text{int}}$ weak-field $B_{\text{ext}} pprox B_{\text{int}}$ intermediate-field

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> $B_{ext} \approx B_{int}$

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depending on the regime, we can use different kinds of perturbation theory

Consider the following relationship between angular momentum and magnetic field:

$$\vec{B} = \frac{1}{4\pi\epsilon_0} \frac{e}{mc^2 r^3} \vec{L}$$

use it to estimate the internal field in hydrogen, and characterize quantitatively a "strong" and "weak" Zeeman field.

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The internal field that gives rise to the spin-orbit interactions is given by

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use it to estimate the internal field in hydrogen, and characterize quantitatively a "strong" and "weak" Zeeman field.

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- a "strong" Zeeman field is $B_{\rm ext}\gg 10{\rm T}$
- a "weak" Zeeman field is $B_{ext} \ll 10 \mathrm{T}$

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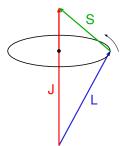
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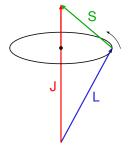
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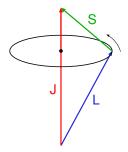
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Weak-field Zeeman effect

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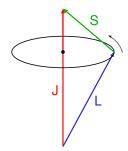
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 $\vec{S}_{ave} = \frac{\vec{S} \cdot \vec{J}}{2} \vec{J}$

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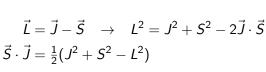
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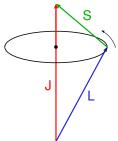
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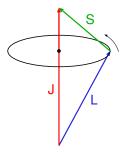
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$$ec{L} = ec{J} - ec{S} \quad o \quad L^2 = J^2 + S^2 - 2 ec{J} \cdot ec{S}$$
 $ec{S} \cdot ec{J} = rac{1}{2}(J^2 + S^2 - L^2)$

$$= rac{\hbar^2}{2}[j(j+1) + s(s+1) - l(l+1)]$$



$$\left\langle \vec{L} + 2\vec{S} \right\rangle = \left\langle \vec{J} + \vec{S} \right\rangle$$

$$\left\langle \vec{L} + 2\vec{S} \right\rangle = \left\langle \vec{J} + \vec{S} \right\rangle = \left\langle \left(1 + \frac{\vec{S} \cdot \vec{J}}{J^2} \right) \vec{J} \right\rangle$$

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$$E_Z^{(1)}=rac{e}{2m}B_{\mathrm{ext}}g_J\hbar m_j \hspace{1.5cm} \mu_B\equivrac{e\hbar}{2m}=5.788 imes10^{-5}\mathrm{eV/T}$$

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where g_J is the so-called Landé g-factor and the full energy correction becomes

$$E_Z^{(1)} = \frac{e}{2m} B_{\text{ext}} g_J \hbar m_j = \mu_B g_J B_{\text{ext}} m_j$$
 $\mu_B \equiv \frac{e\hbar}{2m} = 5.788 \times 10^{-5} \text{eV/T}$

the total energy includes both the spin-orbit and Zeeman corrections and the 2j+1 states then have unique energies

When $B_{\rm ext}\gg B_{\rm int}$, the spin-orbit coupling must be treated as the perturbation and the solutions must be eigenfunctions of the unperturbed wave functions with good quantum numbers: n, l, m_l, s, m_s .

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$$E_{fs}^{1} = \left\langle n I m_{I} m_{s} \left| (H_{r}^{\prime} + H_{so}^{\prime}) \right| n I m_{I} m_{s} \right\rangle$$

$$\langle \vec{S} \cdot \vec{L} \rangle = \langle S_{x} \rangle \langle L_{x} \rangle + \langle S_{y} \rangle \langle L_{y} \rangle + \langle S_{z} \rangle \langle L_{z} \rangle$$

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Intermediate-field Zeeman effect

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for l=0 there is only one value, $j=\frac{1}{2}$, possible with two possible values of m_j directly related to the value of m_s

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for l=1, there are many more mixing possibilities since $m_l=-1,0,1$

Intermediate-field Zeeman effect

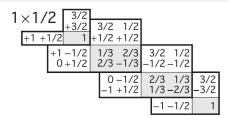
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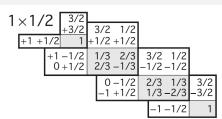
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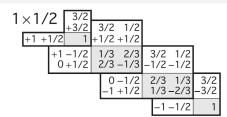
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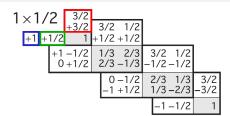
for l=1, there are many more mixing possibilities since $m_l=-1,0,1$ it is convenient to use the $1\times\frac{1}{2}$ Clebsch-Gordan table to generate all the possible states of $|j|m_i\rangle$







$$l=1$$



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Complete n = 2, l = 0, 1 degenerate set

The complete set of 8 n=2, l=0,1 degenerate states using quantum numbers l, s, j, and m_i can now be listed

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now build the W matrix for the $H' = H'_Z + H'_{fs}$ perturbation

$$\gamma \equiv (\alpha/8)^2 \, 13.6 \, \text{eV}, \qquad \beta \equiv \mu_B B_{\text{ext}}$$

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$$W = -$$

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$$\begin{array}{c}
5\gamma - \beta \\
5\gamma + \beta
\end{array}$$

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$$\gamma \equiv (lpha/8)^2 \, 13.6 \, \mathrm{eV}, \qquad eta \equiv \mu_B B_{\mathrm{ext}}$$
 $S\gamma - eta \qquad \qquad \gamma - 2eta \qquad \qquad \gamma + 2eta \qquad \qquad \gamma - rac{2}{3}eta \qquad \qquad \qquad S\gamma - rac{1}{3}eta \qquad \qquad S\gamma - rac{1}{3}\eta \qquad \qquad S\gamma - rac{1}{$

$$\gamma \equiv (lpha/8)^2 \, 13.6 \, \mathrm{eV}, \qquad eta \equiv \mu_B B_{\mathrm{ext}}$$
 $V = -egin{pmatrix} 5\gamma - eta & & & & \\ & 5\gamma + eta & & & \\ & & \gamma - 2eta & & \\ & & & \gamma + 2eta & & \\ & & & \gamma - rac{2}{3}eta & & \\ & & & 5\gamma - rac{1}{3}eta & \\ & & & \gamma + rac{2}{3}eta & & \\ & & & & \gamma + rac{2}{3}eta & & \\ & & & & \gamma + rac{2}{3}eta & & \\ & & & & \gamma + rac{2}{3}eta & & \\ & & & & \gamma + rac{2}{3}eta & & \\ & & & & \gamma + rac{2}{3}eta & & \\ & & & & \gamma + rac{2}{3}eta & & \\ & & & & \gamma + rac{2}{3}eta & & \\ & & & & \gamma + rac{2}{3}eta & & \\ & & & \gamma + rac{2}{3}eta & & \\ & & & \gamma + rac{2}{3}eta & & \\ & & & \gamma + rac{2}{3}eta & & \\ & & & \gamma + rac{2}{3}eta & & \\ & & \gamma + rac{2}{3}eta & & \\ & & \gamma + rac{2}{3}\eta & & \\ & & \gamma + rac{2}{3}\eta & & \\ & \gamma +$

$$\gamma \equiv (\alpha/8)^2 \, 13.6 \, \text{eV}, \qquad \beta \equiv \mu_B B_{\text{ext}}$$

$$W = -\begin{pmatrix}
5\gamma - \beta & & & & \\
& 5\gamma + \beta & & & \\
& & \gamma - 2\beta & & & \\
& & & \gamma + 2\beta & & & \\
& & & \gamma - \frac{2}{3}\beta & & & \\
& & & & 5\gamma - \frac{1}{3}\beta & & \\
& & & & \gamma + \frac{2}{3}\beta & & \\
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\end{pmatrix}$$

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$$W = -\begin{pmatrix} 5\gamma - \beta & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 5\gamma + \beta & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & \gamma - 2\beta & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & \gamma + 2\beta & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & \gamma - \frac{2}{3}\beta & \frac{\sqrt{2}}{3}\beta & 0 & 0 \\ 0 & 0 & 0 & 0 & \frac{\sqrt{2}}{3}\beta & 5\gamma - \frac{1}{3}\beta & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & \gamma + \frac{2}{3}\beta & \frac{\sqrt{2}}{3}\beta \\ 0 & 0 & 0 & 0 & 0 & 0 & \frac{\sqrt{2}}{3}\beta & 5\gamma + \frac{1}{3}\beta \end{pmatrix}$$

All the matrix elements in this basis set will be calculated by you when you do problem 7.29. Defining the fine structure and Zeeman terms with

 ψ_1 , ψ_2 , ψ_3 , and ψ_4 are clearly already eigenfunctions of the full perturbation.

All the matrix elements in this basis set will be calculated by you when you do problem 7.29. Defining the fine structure and Zeeman terms with

 ψ_1 , ψ_2 , ψ_3 , and ψ_4 are clearly already eigenfunctions of the full perturbation. The other 4 eigenfunctions can be solved by solving two 2×2 blocks and then negating the solutions

$$0 = \det \left| \begin{array}{cc} \gamma - \frac{2}{3}\beta - \lambda & \frac{\sqrt{2}}{3}\beta \\ \frac{\sqrt{2}}{3}\beta & 5\gamma - \frac{1}{3}\beta - \lambda \end{array} \right|$$

$$0=det\left|egin{array}{ccc} \gamma-rac{2}{3}eta-\lambda & rac{\sqrt{2}}{3}eta \ rac{\sqrt{2}}{3}eta & 5\gamma-rac{1}{3}eta-\lambda \end{array}
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The first block is solved by diagonalizing the 2×2 submatrix

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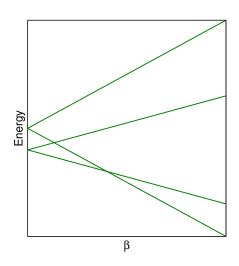
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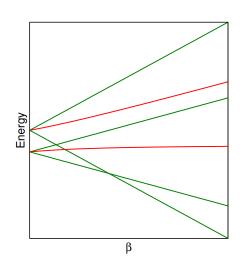
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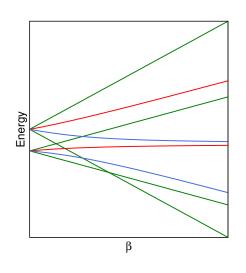
The four solutions which were already eigenfunctions start at two energies and vary linearly with $\boldsymbol{\beta}$



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The two solutions from the first block appear linear at small values of β but show curvature as β increases

The two solutions of the second block mirror those of the first

$$\vec{\mu}_p = \frac{g_p e}{2m_p} \vec{S}_p$$

$$\vec{\mu}_p = rac{g_p e}{2m_p} \vec{S}_p, \quad \vec{\mu}_e = -rac{g_e e}{2m_e} \vec{S}_e$$

where
$$g_p = 5.59$$

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Like the electron, the proton has an associated magnetic dipole where $g_p=5.59$

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$$E_{hf}^{(1)} = \frac{\mu_0 g_p e^2}{8\pi m_p m_e} \left\langle \frac{3(\vec{S}_p \cdot \hat{r})(\vec{S}_e \cdot \hat{r}) - \vec{S}_p \cdot \vec{S}_e}{r^3} \right\rangle$$

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$$\begin{split} H'_{hf} &= \frac{\mu_0 g_p e^2}{8\pi m_p m_e} \frac{[3(\vec{S}_p \cdot \hat{r})(\vec{S}_e \cdot \hat{r}) - \vec{S}_p \cdot \vec{S}_e]}{r^3} + \frac{\mu_0 g_p e^2}{3m_p m_e} (\vec{S}_p \cdot \vec{S}_e) \delta^3(\vec{r}) \\ E^{(1)}_{hf} &= \frac{\mu_0 g_p e^2}{8\pi m_p m_e} \left\langle \frac{3(\vec{S}_p \cdot \hat{r})(\vec{S}_e \cdot \hat{r}) - \vec{S}_p \cdot \vec{S}_e}{r^3} \right\rangle + \frac{\mu_0 g_p e^2}{3m_p m_e} \langle \vec{S}_p \cdot \vec{S}_e \rangle |\psi(0)|^2 \end{split}$$

Problem 7.31

Let \vec{a} and \vec{b} be two constant vectors. Show that

$$\int (\vec{a} \cdot \hat{r})(\vec{b} \cdot \hat{r}) \sin \theta \, d\theta \, d\phi = \frac{4\pi}{3} (\vec{a} \cdot \vec{b})$$

Use this result to demonstrate that

$$\left\langle \frac{3(\vec{S}_p \cdot \hat{r})(\vec{S}_e \cdot \hat{r}) - \vec{S}_p \cdot \vec{S}_e}{r^3} \right\rangle = 0$$

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the angular integrals A and B can be done first, ignoring the specifics of the radial function

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$$B = -\vec{S_p} \cdot \vec{S_e} \int_0^{\pi} \int_0^{2\pi} \sin\theta \, d\theta \, d\phi = -\vec{S_p} \cdot \vec{S_e} 4\pi$$

A and B cancel exactly, giving the desired result

$$E_{hf}^{(1)} = \frac{\mu_0 g_p e^2}{8\pi m_p m_e} \left\langle \frac{3(\vec{S}_p \cdot \hat{r})(\vec{S}_e \cdot \hat{r}) - \vec{S}_p \cdot \vec{S}_e}{r^3} \right\rangle + \frac{\mu_0 g_p e^2}{3m_p m_e} \langle \vec{S}_p \cdot \vec{S}_e \rangle |\psi(0)|^2$$

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for states with I=0 the first term vanishes and we are left with an energy correction which depends on a coupling between the spin of the proton and the spin of the electron

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for the ground state,

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