

Zeeman effect

Prob1: A spectrometer can just resolve a wavelength of 0.5 \AA at 5000 \AA . What should be the strength of external magnetic field to observe the normal Zeeman effect?

Sol: The Zeeman shift is $\Delta \nu = \frac{eB}{4\pi m}$.

We have, $\nu = \frac{c}{\lambda} \Rightarrow \Delta \nu = -\frac{c}{\lambda^2} \Delta \lambda$

$\therefore \Delta \lambda = \frac{\lambda^2}{c} \frac{eB}{4\pi m} \quad \text{or} \quad B = \frac{\Delta \lambda}{\lambda^2} c \frac{4\pi m}{e}$

We have $\Delta \lambda = 0.5 \text{ \AA}$ and $\lambda = 5000 \text{ \AA}$, $B \sim 4.29 \text{ wb/m}^2$.

Prob2: One of the most prominent spectral lines of Calcium is one with wavelength $\lambda = 4226.73 \text{ \AA}$. Calcium atoms exhibit normal Zeeman patterns when placed in a uniform magnetic field of 4 wb/m^2 . Calculate the wavelength of the three components of normal Zeeman pattern & the separation between them.

Soln: The Zeeman separation is $\Delta \nu = \frac{eB}{4\pi m}$

$\& \Delta \lambda = \frac{\lambda^2}{c} \frac{eB}{4\pi m} \sim 0.33 \text{ \AA}$

\therefore The wavelengths of the normal Zeeman components are $(4226.73 \pm 0.33) \text{ \AA}$.

Q Find the equatorial velocity v of an electron under the assumption that it is a uniform sphere of radius $r = 5.00 \times 10^{-17} \text{ m}$ that is rotating about an axis through its centre.

Soln: The angular momentum is $L = I\omega = \frac{2}{5} m r^2 \omega$ & $\omega = \frac{v}{r}$ is

its angular velocity. The total spin angular momentum is

$$S = \sqrt{3(3+1)} \hbar = \frac{\sqrt{3}}{2} \hbar = \frac{2}{5} m r^2 \frac{v}{r} = \frac{2}{5} m v r$$

$$\Rightarrow v = \frac{5\sqrt{3}}{4} \frac{\hbar}{m r} \sim 5.01 \times 10^{12} \text{ ms}^{-1} \sim 10^4 c \sim \text{impossible}$$

Prob: Calculate the possible orientation of spin vector \vec{S} with respect to magnetic field direction.

$$\text{Sol: } \cos \theta = \frac{S_z}{S} = \frac{m_s \hbar}{\sqrt{s(s+1)} \hbar} = \frac{m_s}{\sqrt{\frac{1}{2}(\frac{1}{2}+1)}} = \frac{2}{\sqrt{3}} m_s$$

$$\text{now, } m_s = \pm \frac{1}{2} \quad \text{so } \theta = \cos^{-1}\left(\frac{1}{\sqrt{3}}\right) \quad \text{and } \theta = \cos^{-1}\left(-\frac{1}{\sqrt{3}}\right)$$

$$= 54.7^\circ \quad \text{and } 125.3^\circ$$

Prob: A beam of electrons enters a uniform magnetic field of flux density 1.2 Wb/m^2 . Find the energy difference in eV between electrons whose spins are parallel and antiparallel to the field.

$$\text{Sol: } \vec{\mu}_s = -\frac{e}{m} \vec{S} \hbar, \quad \text{Along } z \text{ axis, } \mu_{sz} = -\frac{e}{m} S_z = \mp \frac{eh}{4\pi m}$$

$$\therefore U_m = -\vec{\mu}_s \cdot \vec{B} = -\mu_{sz} B = \mp \frac{ehB}{4\pi m}$$

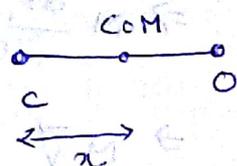
$$\therefore \text{difference} = \frac{ehB}{4\pi m} - \left(-\frac{ehB}{4\pi m}\right) = \frac{ehB}{2\pi m}$$

Centre of mass problems.

→ The distance between the centre of Carbon and oxygen atoms in the carbon monoxide (CO) gas molecule is $1.130 \times 10^{-10} \text{ m}$. Locate the centre of mass of the molecule relative to the carbon atom.

Sol: The CO gas molecule is a linear molecule.

So the COM is likely to remain in the line



Joining the carbon and oxygen atom. So

if x is the distance of COM from C, then we can always write

$$\Rightarrow 12x = 16(1.130 \times 10^{-10} - x) \quad \Rightarrow 28x = 16 \times 1.130 \times 10^{-10}$$

$$\Rightarrow x = 6.487 \times 10^{-11} \text{ m}$$

Thus, the splitting is shown below, with the selection rules

$$\Delta M_J = 0, \pm 1 \quad (M_J = 0 \not\leftrightarrow \mp M_J = 0 \text{ if } \Delta J = 0).$$

→ The Paschen-Back Effect :-

In anomalous Zeeman effect, remember that the applied external magnetic field was weak. This means that the internal mag. field produced due to the motion of \vec{e} is ~~less~~ strong enough as compared to the external B . This means that for such an atom, \vec{L} and \vec{S} keeps on precessing around \vec{J} .

However, when one increases the mag. field \vec{B} (external), the separation between the anomalous Zeeman components keep on increasing until they exceed the separation between multiple fine-structure components. (For eg- in case of D_1 & D_2 lines in a weak mag. field (i.e. anomalous Zeeman effect) if you keep on increasing the external magnetic field \vec{B} , then at one point of time, the separation between the levels will keep on increasing in such a way that at one point of time there will be overlapping between the levels (see the various M_J levels in anomalous Zeeman effect in page 208, Raj Kumar). As a result at some point of time instead of 4 & 6 lines from D_1 and D_2 , we shall see only three lines exactly similar to the normal Zeeman effect). This is the Paschen Back effect.

Explanation :- As mentioned previously, when a mag. field \vec{B} is applied & becomes stronger than the internal fields, \vec{J} starts precessing around \vec{B} very fast. Remember that in a weak mag. field, the precession of \vec{L} & \vec{S} is fast around \vec{J} as compared to \vec{J} around \vec{B} . Thus, the coupling between \vec{L} & \vec{S} remains intact. On applying a strong B field, the exactly

reverse scenario happens. \vec{S} starts rotating around \vec{B} very fast as compared to \vec{L} & \vec{S} around \vec{J} . As the field is increased more, \vec{L} & \vec{S} starts independently precessing around \vec{B} . While precessing independently, the component \vec{L} & \vec{S} around \vec{B} gets quantised so that

$$L_z = M_L \frac{h}{2\pi} \quad \& \quad S_z = M_S \frac{h}{2\pi}$$

where $M_L = L, L-1, \dots, -L$ &

$M_S = S, S-1, \dots, -S$.

The velocity of these independent precessions are

$$\omega_L = \frac{e}{2m} B \quad \& \quad \omega_S = 2 \frac{e}{2m} B$$

& the interacting energy of these precession are

$$\Delta E_L = \omega_L L_z = \frac{e}{2m} B M_L \hbar \quad (\text{Refer previous class})$$

$$\Delta E_S = \omega_S S_z = 2 \frac{e}{2m} B M_S \hbar$$

∴ the total interaction energy is

$$\Delta E = (M_L + 2M_S) \frac{e\hbar}{4\pi m} B$$

& wave number shift is,

$$-\Delta T = \frac{\Delta E}{hc} = (M_L + 2M_S) \frac{eB}{4\pi mc}$$

In Lorentz unit,

$$-\Delta T = (M_L + 2M_S) L'$$

This is the expression for strong-field interaction energy. Now M_L has $(2L+1)$ values & M_S has $(2S+1)$ components. So each of the field free level is split into $(2L+1)(2S+1)$ levels.

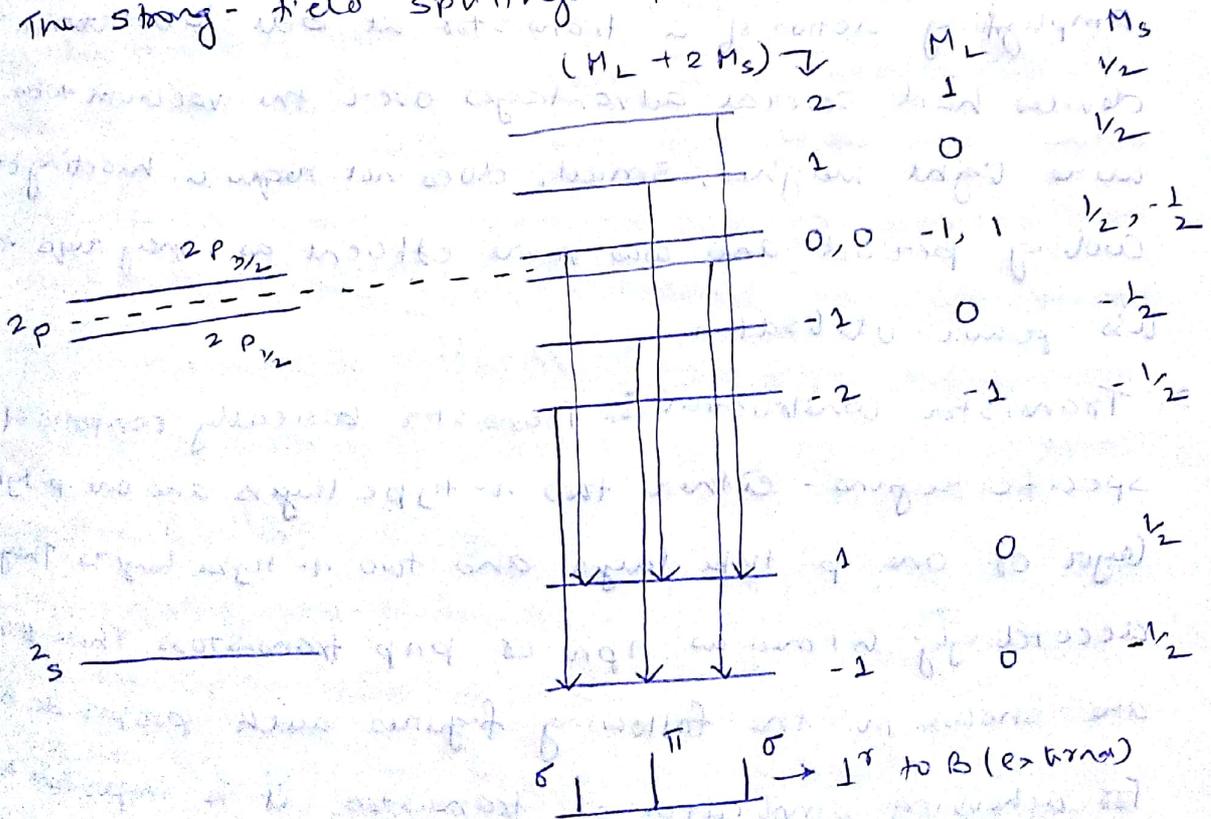
Let us consider the following examples

$$2p - 2s$$

This is the transition responsible for D_1 and D_2 . Now,

Term	No. of Strong-field levels ($2L+1)(2S+1)$	M_L	M_S	Shift in L unit ($M_L + 2M_S$)
		1	$\frac{1}{2}, -\frac{1}{2}$	2, 0
		0	$\frac{1}{2}, -\frac{1}{2}$	1, -1
$2p$ $L=1, S=\frac{1}{2}$	6	-1	$\frac{1}{2}, -\frac{1}{2}$	0, -2
		0	$\frac{1}{2}, -\frac{1}{2}$	1, -1
$2s$ $L=0, S=\frac{1}{2}$	2	0	$\frac{1}{2}, -\frac{1}{2}$	1, -1

The strong-field splittings of the terms 2^1p & 2^2s have



$$\Delta M_L = 0 \text{ (} \parallel \text{ to field)}$$

$$\Delta M_L = \pm 1 \text{ (} \perp \text{ to field)}$$

$$\Delta M_S = 0$$