

Inner Time and The Inner ear

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Abstract

This document will contain the relevant equations I need to numerically calculate the Paschen-Back for arbitrary I and J .

I. THE WHOLE PAPER

The hyperfine splitting up to magnetic octopole contributions is given by:

$$\begin{aligned}
H_{\text{hfs}} = & A_{\text{hfs}} \frac{\mathbf{I} \cdot \mathbf{J}}{\hbar^2} \\
& + B_{\text{hfs}} \frac{\frac{3}{\hbar^2} (\mathbf{I} \cdot \mathbf{J})^2 + \frac{3}{2\hbar} \mathbf{I} \cdot \mathbf{J} - I(I+1)J(J+1)}{2I(2I-1)J(2J-1)} \\
& + C_{\text{hfs}} \frac{\frac{10}{\hbar^3} (\mathbf{I} \cdot \mathbf{J})^3 + \frac{20}{\hbar^2} (\mathbf{I} \cdot \mathbf{J})^2 + \frac{2}{\hbar} \mathbf{I} \cdot \mathbf{J} [I(I+1) + J(J+1) + 3] - 3I(I+1)J(J+1) - 5I(I+1)J(J+1)}{I(I-1)(2I-1)J(J-1)(2J-1)}
\end{aligned} \tag{1}$$

For Sodium 23, we have:

Magnetic Dipole Constant, $3^2S_{1/2}$	$A_{3^2S_{1/2}}$	$h \cdot 885.813$ MHz	(2)
Magnetic Dipole Constant, $3^2P_{1/2}$	$A_{3^2P_{1/2}}$	$h \cdot 94.44$ MHz	
Magnetic Dipole Constant, $3^2P_{3/2}$	$A_{3^2P_{3/2}}$	$h \cdot 18.534$ MHz	
Electric Quadrupole Constant, $3^2P_{3/2}$	$B_{3^2P_{3/2}}$	$h \cdot 2.724$ MHz	

For Lithium 6:

Magnetic Dipole Constant, $2^2S_{1/2}$	$A_{2^2S_{1/2}}$	$h \cdot 152.137$ MHz	(3)
Magnetic Dipole Constant, $2^2P_{1/2}$	$A_{2^2P_{1/2}}$	$h \cdot 17.386$ MHz	
Magnetic Dipole Constant, $2^2P_{3/2}$	$A_{2^2P_{3/2}}$	$h \cdot -1.155$ MHz	
Electric Quadrupole Constant, $2^2P_{3/2}$	$B_{2^2P_{3/2}}$	$h \cdot -0.10$ MHz	

For the DC Zeeman shift, we want to compute the eigenstates of

$$H_B^{(\text{fs})} + H_B^{(\text{hfs})} \tag{4}$$

where for a B -field in the z direction, we have

$$H_B^{(\text{fs})} = -\mu_S \cdot \mathbf{B} - \mu_L \cdot \mathbf{B} = \frac{\mu_B}{\hbar} (g_S S_z + g_L L_z) B \tag{5}$$

and

$$H_B^{(\text{hfs})} = -\mu_I \cdot \mathbf{B} = \frac{\mu_B}{\hbar} g_I I_z B. \tag{6}$$

For now, we will lump together the effect of S and L into J via the Lande g_J factor, and treat only the Paschen-Back effect due to the hyperfine interaction (since this occurs

at much lower energy B -fields than the separation of S and L eigenstates). Thus we must diagonalize

$$H_B^{(\text{fs})} + H_B^{(\text{hfs})} = \frac{\mu_B}{\hbar} (g_J J_z + g_I I_z) B \quad (7)$$

where

$$\begin{aligned} g_J &\equiv g_L + (g_S - g_L) \frac{J(J+1) + S(S+1) - L(L+1)}{2J(J+1)} \\ g_L &\equiv \frac{\text{Reduced Mass}}{m_e} = \frac{1}{1 + m_e/m_n} \\ g_S &= 2.002319. \end{aligned} \quad (8)$$

In order to compute matrix elements of $\mathbf{I} \cdot \mathbf{J}$ in $H_B^{(\text{hfs})}$, we will need to use:

$$\begin{aligned} \mathbf{I} \cdot \mathbf{J} &= I_z J_z + \frac{I_+ J_- + I_- J_+}{2} \\ (\mathbf{I} \cdot \mathbf{J})^2 &= (I_z J_z)^2 + \frac{1}{2} \{I_z J_z, I_+ J_- + I_- J_+\} + \frac{(I_+ J_-)^2 + (I_- J_+)^2}{4} + \frac{I_+ I_- J_- J_+ + I_- I_+ J_+ J_-}{4} \end{aligned} \quad (9)$$

We will calculate the energy splittings ignoring the octopole, C_{hfs} term in eq. (1). For the case of $I = 1/2$ or $J = 1/2$, the B_{hfs} term vanishes (but does it really? The denominator is certainly zero—what about the numerator?), and the result would be the Breit-Rabi formula. Thus, for a first non-trivial go at numerical calculations, let's try the case of the Lithium 6