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Physics 472

Exam 2

Spring 2019

Non-Degenerate Perturbation Theory

If a level n state with energy $E_n^{(0)}$ and eigenfunction $\psi_n^{(0)}$ is non-degenerate, then the perturbed energy and eigenfunction are given by

$$E_n = E_n^{(0)} + \langle \psi_n^{(0)} | H' | \psi_n^{(0)} \rangle + \sum_{k \neq n} \frac{|\langle \psi_k^{(0)} | H' | \psi_n^{(0)} \rangle|^2}{E_n^{(0)} - E_k^{(0)}} + \dots$$

$$\psi_n = \psi_n^{(0)} + \sum_{k \neq n} \frac{\psi_k^{(0)} \langle \psi_k^{(0)} | H' | \psi_n^{(0)} \rangle}{E_n^{(0)} - E_k^{(0)}} + \dots$$

where H' is the perturbation.

Degenerate Perturbation Theory

If an energy level is n -fold degenerate with its n orthogonal eigenfunctions labelled as $\psi_1, \psi_2, \dots, \psi_n$ and $H'_{jk} = \langle \psi_j | H' | \psi_k \rangle$, where H' is a perturbation, then the first order energy shifts to this level are determined by

$$\begin{vmatrix} H'_{11} - E & H'_{12} & H'_{13} & \dots & H'_{1n} \\ H'_{21} & H'_{22} - E & H'_{23} & \dots & H'_{2n} \\ H'_{31} & H'_{32} & H'_{33} - E & \dots & H'_{3n} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ H'_{n1} & H'_{n2} & H'_{n3} & \dots & H'_{nn} - E \end{vmatrix} = 0$$

Variational Method

If H is a Hamiltonian with ground state energy E_0 , then for any normalized state $|\psi\rangle$,

$$E_0 \leq \langle \psi | H | \psi \rangle.$$

Symmetries

For hydrogen eigenstates $|n\ell m\rangle$, rotational invariance implies that $(m' - m)\langle n\ell' m' | z | n\ell m \rangle = 0$, $\ell' = \ell \pm 1$ and parity invariance implies $\langle n\ell' m' | z | n\ell m \rangle = (-1)^{\ell' + \ell + 1} \langle n\ell' m' | z | n\ell m \rangle$.

Zeeman Effect

The perturbative Hamiltonian for the Zeeman effect gives, in first order

$$\langle H'_Z \rangle = \frac{\mu_B}{\hbar} B \langle \ell \frac{1}{2} j m_j | J_z + S_z | \ell \frac{1}{2} j m_j \rangle = \mu_B B g_L m_j,$$

where B is the magnitude of the external magnetic field.

Miscellaneous

<u>Harmonic Oscillator</u>	$\left(\frac{p^2}{2m} + \frac{1}{2} m \omega^2 x^2 \right) \psi_n = E_n \psi_n$	$E_n = \hbar \omega (n + \frac{1}{2}),$
<u>Hydrogen</u>	$\left(\frac{\vec{p}^2}{2m} - \frac{\hbar c \alpha}{r} \right) \psi_{n\ell m} = E_n \psi_{n\ell m}$	$E_n = -\frac{m c^2 \alpha^2}{2n^2} = -\frac{13.6 \text{eV}}{n^2}.$

1. The ground state wave function and energy for the hydrogen atom are

$$\psi_0(\vec{r}) = \left(\frac{1}{\pi a^3} \right)^{\frac{1}{2}} e^{-r/a}, \quad E_0 = -\frac{1}{2} m c^2 \alpha^2 = -13.6 \text{ eV},$$

where $a = \hbar/mc\alpha$ is the Bohr radius. The atom is perturbed by a potential linear in r resulting in the Hamiltonian

$$H = \frac{\vec{p}^2}{2m} - \frac{\hbar c \alpha}{r} + \frac{2}{3} m c^2 \alpha^2 \lambda \frac{r}{a},$$

where $\lambda > 0$ and the term proportional to λ is the perturbation H' .

- (a) (3 pt) Calculate the ground state energy using first order perturbation theory. (*I.e.*, calculate $E^{(0)} + E^{(1)}$ for the ground state.)
- (b) (3 pt) To obtain an improved estimate of the ground state energy, use the trial wave function

$$\psi(\vec{r}) = \left(\frac{Z^3}{\pi a^3} \right)^{\frac{1}{2}} e^{-Zr/a}$$

and calculate $\langle \psi | H | \psi \rangle$. Note: With this trial wave function we have

$$\langle \psi | \vec{p}^2 | \psi \rangle = m^2 c^2 \alpha^2 Z^2, \quad \langle \psi | \frac{1}{r} | \psi \rangle = \frac{m c \alpha Z}{\hbar}, \quad \langle \psi | r | \psi \rangle = \frac{3a}{2Z}.$$

- (c) (2 pt) Derive the equation satisfied by Z which determines the minimum of $\langle \psi | H | \psi \rangle$. (You do not have to solve this equation.)

2. A hydrogen atom is placed in a static electric field $\vec{\mathcal{E}}$ of magnitude $2 \times 10^4 \text{ V/m}$. The atom experiences a perturbation given by the Stark interaction $H' = e\vec{r} \cdot \vec{\mathcal{E}}$.
- (a) (1 pt) Is it possible for this perturbation to shift the energy of the ground ($n = 1$) state? If so, in what order of perturbation theory does this first occur?
 - (b) (1 pt) Is it possible for this perturbation to shift the energy of the first excited ($n = 2$) state? If so, in what order of perturbation theory does this first occur?
 - (c) (2 pt) List the $n = 2$ states (ignoring spin) and state which matrix elements of H' between these states can be nonzero.
 - (d) (2 pt) Calculate the leading order energy shifts to the $n = 2$ levels. You needn't evaluate any non-vanishing matrix elements that occur, but clearly define your notation.
 - (e) (1 pt) Given that the Bohr radius $a = 0.53 \times 10^{-10} \text{ m}$, estimate the size of a typical Stark energy shift in electron volts for the given electric field strength $\mathcal{E} = 2 \times 10^4 \text{ V/m}$.

3. Give a brief but reasoned answer to each of the following:

- (a) (2 pt) When describing the weak field Zeeman effect in hydrogen, the expectation value of the magnetic interaction is expressed in terms of the Landé g -factor g_L as

$$E_{\text{Zee}} = \frac{\mu_B B}{\hbar} \langle n\ell \frac{1}{2} j m_j | (J_z + S_z) | n\ell \frac{1}{2} j m_j \rangle = \mu_B B g_L m_j ,$$

where $\mu_B = e\hbar/2m_e$ is the Bohr magneton and the eigenvalue of J_z is $\hbar m_j$. What is the value of g_L for an electron in the state $\psi_{32j}^{m_j}(\vec{r}) = R_{32}(r)Y_2^2(\theta, \phi)\chi^{1/2}$?

- (b) (1 pt) Given that $v/c \approx \alpha = 1/137$ in the hydrogen atom and that the relativistic kinetic energy correction is $H'_{KE} = -(\vec{p}^2)^2/8m^3c^2 = -(\vec{p}^2/2m)(\frac{\vec{p}^2}{4m^2c^2})$, estimate the size of the fine-structure corrections in eV. (An order of magnitude estimate is all that is wanted here.)
- (c) (2 pt) The state vectors $|n j m_j \ell \frac{1}{2}\rangle$ are eigenvectors of the operator $\vec{L} \cdot \vec{S}$. What is the eigenvalue for an electron in the $4D_{\frac{3}{2}}$ state of hydrogen? (Note: here we use the notation $n\ell_j$ for neutral hydrogen, since $s = \frac{1}{2}$ is always assumed.)

