

Physics 472 – Spring 2009

Homework #9, due Friday, March 27

(Point values are in parentheses.)

1. [7] The isotropic 2-dimensional harmonic oscillator is easily solved by writing the Hamiltonian as a sum of x and y Hamiltonians:

$$\hat{H}^0 = \frac{\hat{P}_x^2 + \hat{P}_y^2}{2m} + \frac{1}{2}m\omega^2(x^2 + y^2) = \hat{H}_x^0 + \hat{H}_y^0 \quad \text{with} \quad [\hat{H}_x^0, \hat{H}_y^0] = 0.$$

Simultaneous eigenstates of \hat{H}_x^0 and \hat{H}_y^0 obey $\hat{H}^0 |n_x, n_y\rangle = (n_x + n_y + 1)\hbar\omega |n_x, n_y\rangle$.

- a) Consider the perturbation $\lambda\hat{H}' = \lambda m\omega^2 xy$. Calculate the first and second order energy shifts of the ground state. In class we used $\hat{X} = \sqrt{\frac{\hbar}{2m\omega}}(\hat{a}_x^+ + \hat{a}_x)$ to evaluate the matrix elements:

$$\langle n_x' | x | n_x \rangle = \sqrt{\frac{\hbar}{2m\omega}} (\sqrt{n_x} \delta_{n_x', n_x-1} + \sqrt{n_x+1} \delta_{n_x', n_x+1}). \quad \text{The same holds for } \hat{Y} = \sqrt{\frac{\hbar}{2m\omega}}(\hat{a}_y^+ + \hat{a}_y).$$

To evaluate the matrix elements of \hat{H}' in the $|n_x n_y\rangle$ tensor product states, use:

$$\langle n_x' n_y' | xy | n_x n_y \rangle = \langle n_x' | x | n_x \rangle \langle n_y' | y | n_y \rangle$$

- b) Use degenerate P.T. to calculate the first-order energy shifts of the first excited states, as well as the “correct” linear combinations of those two states that diagonalize \hat{H}' .

- c) The full Hamiltonian, $\hat{H} = \hat{H}^0 + \lambda\hat{H}'$, is exactly solvable if you make the coordinate transformation $u = (x + y)/\sqrt{2}$, $v = (x - y)/\sqrt{2}$. Express \hat{H} in terms of u , v , and their conjugate momenta P_u , and P_v . You should find that the harmonic oscillator in the “ u ” direction has a higher frequency than before, while in the “ v ” direction the frequency is lower. Calculate the exact energies of the new basis states $|n_u, n_v\rangle$. For the ground state, expand the energy to second order in λ . For the next two higher states, expand the energies to first order in λ . Compare your results with those you obtained in parts (a) and (b).

2. [7] Consider an electron in a 3-dimensional isotropic harmonic oscillator potential, in the presence of a uniform magnetic field $\vec{B} = B_{ext} \hat{k}$. The full Hamiltonian for the system is:

$$\hat{H} = \frac{\hat{P}_x^2 + \hat{P}_y^2 + \hat{P}_z^2}{2m} + \frac{1}{2}m\omega^2(x^2 + y^2 + z^2) + \frac{eB_{ext}}{2m}(L_z + 2S_z)$$

This problem is exactly solvable, but we’ll use the machinery of perturbation theory to get organized. Treat the first two terms of H as H^0 , and the Zeeman term as $\lambda H'$. The eigenstates of H^0 satisfy: $\hat{H}^0 |n_x, n_y, n_z, m_s\rangle = (n_x + n_y + n_z + \frac{3}{2})\hbar\omega |n_x, n_y, n_z, m_s\rangle$, where $m_s = \pm \frac{1}{2}$, and $S_z |n_x, n_y, n_z, m_s\rangle = \hbar m_s |n_x, n_y, n_z, m_s\rangle$. (I am putting the space and spin quantum numbers together inside the same ket to avoid using the cumbersome tensor product notation.)

a) Express \hat{L}_z in terms of the harmonic oscillator raising and lowering operators. Hint: you should get $\hat{L}_z = i\hbar(a_x a_y^+ - a_x^+ a_y)$.

b) The ground state of H^0 is 2-fold degenerate, due to spin. But since the two $|0,0,0, m_s\rangle$ states are already eigenstates of H' , you can use standard first-order perturbation theory to calculate the energy shifts due to the magnetic field. Express your answers in terms of $\mu_B B_{ext}$ and m_s .

c) The first excited state of H^0 is six-fold degenerate (3 spatial states \times 2 spin states). Calculate the linear combinations of states that diagonalize H' . To help you keep track of what you are doing, here are some suggestions. First, since all your states are eigenstates of S_z , leave spin out of the problem until the end; then you only have to diagonalize a 3×3 matrix rather than a 6×6 matrix. Since the original basis states $|n_x, n_y, n_z, m_s\rangle$ are not eigenstates of \hat{L}_z , you need to find linear combinations of them that are. Label the new states this way: $|n, l, m_l, m_s\rangle$, where $n = n_x + n_y + n_z$. You don't need to know l to do this problem – you just need m_l . But you can probably guess what l is once you know what m_l is for the three states. Finally, when you have found the states that diagonalize H' , calculate the Zeeman energy shifts of those states. How many distinct energies are there? Make a plot of energy vs. $\mu_B B_{ext}$ for all the states.

d) The second excited state of H^0 is twelve-fold degenerate (6 spatial states \times 2 spin states). Forget about spin altogether so you don't get lost. Construct the 6×6 matrix representation of \hat{L}_z . If you choose the order of your 6 states judiciously, your 6×6 matrix should break up into a 2×2 block, a 3×3 block, and a 1×1 block. Calculate the eigenvalues of \hat{L}_z and their degeneracies. Guess what the values of l are for this six-dimensional subspace. Don't bother to calculate the 12 Zeeman energies – I know you could do it if you had to!

3. [6] Griffiths problem 6.37. Follow the same strategy you used to solve Griffiths problem 6.36. Use symmetries to figure out which matrix elements of the form $\langle n, l', m_l' | z | n, l, m_l \rangle$ are zero.

Rotational symmetry, $[\hat{z}, \hat{L}_z] = 0$, implies $(m_l' - m_l) \langle n, l', m_l' | z | n, l, m_l \rangle = 0$. The parity transformation, $\hat{\Pi} \hat{z} \hat{\Pi} = -\hat{z}$, implies $(-1)^{l'+l} \langle n, l', m_l' | z | n, l, m_l \rangle = -\langle n, l', m_l' | z | n, l, m_l \rangle$.

Use these same symmetries to do Griffiths problem 6.37. First, show which elements of the 9×9 matrix are zero. Then calculate the first non-zero matrix element, $\langle 3, 0, 0 | z | 3, 1, 0 \rangle$, using Tables 4.3 and 4.7 in Griffiths. Use Mathematica to do the radial integration. You can take the values of the other nonzero matrix elements from Griffiths. Construct the 9×9 matrix representation of \hat{z} . If you choose the order of the 9 states carefully, then the matrix should break into a 3×3 block, two 2×2 blocks, and two trivial 1×1 blocks. Calculate the eigenvalues and their degeneracies. Don't forget to multiply the eigenvalues by eE_{ext} to get the energies.