CAS PY 452 — Quantum Physics II

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Perturbing two harmonic oscillators

Problem statement: Consider a particle in a two-dimensional isotropic quantum harmonic oscillator; that is, consider the following Hamiltonian:

$$\hat{H}_0 = \frac{\hat{p}_x^2}{2m} + \frac{\hat{p}_y^2}{2m} + \frac{1}{2}m\omega^2\hat{x}^2 + \frac{1}{2}m\omega^2\hat{y}^2.$$
 (1)

Now perturb the Hamiltonian $\hat{H} = \hat{H}_0 + \hat{H}'$ with,

$$\hat{H}' = \lambda \hat{x}\hat{y} \tag{2}$$

where $\lambda \ll m\omega^2/2$. Find the first and second-order corrections to the energy of the first excited state (which is doubly-degenerate), and find the first-order corrections to the first excited state.

Solution We will use ladder operators to compute inner products; if any of this is unclear, you should look at the discussion notes from Oct. 5. Here are some identities for reference:

$$\hat{a} = \sqrt{\frac{m\omega}{2\hbar}}(\hat{x} + \frac{i}{m\omega}\hat{p}), \qquad \hat{a}^{\dagger} = \sqrt{\frac{m\omega}{2\hbar}}(\hat{x} - \frac{i}{m\omega}\hat{p}),$$
 (3)

$$\hat{x} = \sqrt{\frac{\hbar}{2m\omega}} (\hat{a}^{\dagger} + \hat{a}), \qquad \hat{p} = i\sqrt{\frac{\hbar m\omega}{2}} (\hat{a}^{\dagger} - \hat{a}). \tag{4}$$

Now that we are dealing with two quantum harmonic oscillators, there are two sets of ladder operators: $\hat{a}_x, \hat{a}_x^{\dagger}$ and $\hat{a}_y, \hat{a}_y^{\dagger}$. The first few states are,

$$|\psi_{00}\rangle = |\psi_0(x)\rangle |\psi_0(y)\rangle \tag{5}$$

$$|\psi_{10}\rangle = |\psi_1(x)\rangle |\psi_0(y)\rangle \tag{6}$$

$$|\psi_{01}\rangle = |\psi_0(x)\rangle |\psi_1(y)\rangle \tag{7}$$

$$|\psi_{11}\rangle = |\psi_1(x)\rangle |\psi_1(y)\rangle \tag{8}$$

$$|\psi_{20}\rangle = |\psi_2(x)\rangle |\psi_0(y)\rangle \tag{9}$$

$$|\psi_{02}\rangle = |\psi_0(x)\rangle |\psi_2(y)\rangle \tag{10}$$

$$\vdots (11)$$

We can write the un-perturbed Hamiltonian $\hat{H} = \hbar\omega(\hat{a}_x^{\dagger}\hat{a}_x + \hat{a}_y^{\dagger}\hat{a}_y + 10) = \hbar\omega(\hat{N}_x + \hat{N}_y + 1)$ as a matrix in the basis $|\psi_{00}\rangle, |\psi_{10}\rangle, |\psi_{01}\rangle, |\psi_{11}\rangle, |\psi_{20}\rangle, |\psi_{02}\rangle, |\psi_{12}\rangle, \dots$:

$$\hat{H}_{0} = \hbar\omega \begin{pmatrix} 1 & \cdot \\ \cdot & 2 & \cdot \\ \cdot & \cdot & 2 & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & 3 & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \cdot & 3 & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \cdot & 3 & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \cdot & \cdot & 4 & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \cdot & \cdot & 4 & \cdot & \cdot & \cdot \\ \vdots & \ddots \end{pmatrix}$$

$$(12)$$

and the perturbation as a matrix:

We can see that the first excited states, $|\psi_{10}\rangle$, $|\psi_{01}\rangle$, have off-diagonal matrix elements within their block of the matrix \hat{H}' . This means that when we compute the sum,

$$\sum_{m_x, m_y \neq 0, 1} \frac{\langle \psi_{m_x, m_y} | \hat{H}' | \psi_{n_x, n_y} \rangle}{E_n^{(0)} - E_m^{(0)}} | \psi_{m_x, m_y} \rangle, \tag{15}$$

there will be a zero in the denominator for the $m_x, m_y = 1, 0$ term without a zero in the numerator. However, if we diagonalize the block $\begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \rightarrow \begin{pmatrix} -1 & 0 \\ 0 & 1 \end{pmatrix}$ by choosing new basis states,

$$|\psi_{-}\rangle = \frac{1}{\sqrt{2}} (|\psi_{01}\rangle - |\psi_{10}\rangle), \qquad |\psi_{+}\rangle = \frac{1}{\sqrt{2}} (|\psi_{01}\rangle + |\psi_{10}\rangle),$$
 (16)

we can compute,

$$\begin{split} \hat{H}_{\pm,m_xm_y}' &= \langle \psi_{m_xm_y} | \hat{H}' | \psi_\pm \rangle = \frac{\hbar \lambda}{2\sqrt{2}m\omega} \left\langle \psi_{m_xm_y} | \left(\hat{a}_x^\dagger + \hat{a}_x \right) \left(\hat{a}_y^\dagger + \hat{a}_y \right) \left(|\psi_{01}\rangle \pm |\psi_{10}\rangle \right) \\ &= \frac{\hbar \lambda}{2\sqrt{2}m\omega} \left\langle \psi_{m_xm_y} | \left(\hat{a}_x^\dagger + \hat{a}_x \right) \left(\sqrt{2} \left| \psi_{02} \right\rangle + \left| \psi_{00} \right\rangle \pm \left(\left| \psi_{11} \right\rangle + 0 \right) \right) \\ &= \frac{\hbar \lambda}{2\sqrt{2}m\omega} \left\langle \psi_{m_xm_y} | \left(\sqrt{2} (\left| \psi_{12} \right\rangle + 0) + \left(\left| \psi_{10} \right\rangle + 0 \right) \pm \left(\sqrt{2} \left| \psi_{21} \right\rangle + \left| \psi_{01} \right\rangle \right) \right) \\ &= \frac{\hbar \lambda}{2m\omega} \left\langle \psi_{m_xm_y} | \left(\left| \psi_{12} \right\rangle \pm \left| \psi_{21} \right\rangle \mp \left| \psi_\pm \right\rangle \right) = \frac{\hbar \lambda}{2m\omega} \left(\delta_{12,m_xm_y} \pm \delta_{21,m_xm_y} \mp \delta_{\pm,m_xm_y} \right). \end{split}$$

These new states are still energy eigenstates because a linear combination of eigenvectors with a common eigenvalue is also an eigenvector with the same eigenvalue; for a visualization of these new states in the position representation,

see figure 1. In the new basis $|\psi_{00}\rangle$, $|\psi_{-}\rangle$, $|\psi_{+}\rangle$, $|\psi_{11}\rangle$, $|\psi_{20}\rangle$, $|\psi_{02}\rangle$, $|\psi_{21}\rangle$, $|\psi_{12}\rangle$, ..., the perturbed Hamiltonian is,

$$\hat{H}' = \frac{\hbar\lambda}{2m\omega} \begin{pmatrix} \cdot & \cdot & \cdot & 1 & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & -1 & 0 & \cdot & \cdot & \cdot & -1 & 1 & \cdot \cdot \\ \cdot & 0 & 1 & \cdot & \cdot & \cdot & 1 & 1 & \cdot \cdot \\ 1 & \cdot & \cdot & \cdot & \sqrt{2} & \sqrt{2} & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \sqrt{2} & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \sqrt{2} & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \sqrt{2} & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \sqrt{2} & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & 1 & 1 & \cdot & \cdot & \cdot & \cdot & 2 & \cdot \cdot \\ \cdot & \cdot & 1 & 1 & \cdot & \cdot & \cdot & 2 & \cdot & \cdot \\ \vdots & \ddots \end{pmatrix}$$

$$(17)$$

Because we diagonalized the degenerate block by changing bases, we see that when we compute the +,- terms in the sum on line 15, we do not have to deal with a zero in the denominator because the zero in the numerator takes precedence. We have discovered that when we perturb the Hamiltonian, the first excited state splits into two states $|\psi_{-}\rangle$, $|\psi_{+}\rangle$, and we can compute the corrections to the energies and states using the matrix elements from line 17 and 12.

$$E_{-}^{(1)} = \langle \psi_{-} | \hat{H}' | \psi_{-} \rangle = -\frac{\hbar \lambda}{2m\omega} \tag{18}$$

$$E_{+}^{(1)} = \langle \psi_{+} | \hat{H}' | \psi_{+} \rangle = + \frac{\hbar \lambda}{2m\omega} \tag{19}$$

$$E_{-}^{(2)} = \frac{|\langle \psi_{21} | \hat{H}' | \psi_{-} \rangle|^{2}}{\hbar \omega (2 - 4)} + \frac{|\langle \psi_{12} | \hat{H}' | \psi_{-} \rangle|^{2}}{\hbar \omega (2 - 4)} = \frac{\hbar \lambda^{2}}{4m^{2}\omega^{3}} \left(\frac{|-1|^{2}}{-2} + \frac{|1|^{2}}{-2} \right) = -\frac{\hbar \lambda^{2}}{4m^{2}\omega^{3}}$$
(20)

$$E_{+}^{(2)} = \frac{|\langle \psi_{21} | \hat{H}' | \psi_{+} \rangle|^{2}}{\hbar \omega (2 - 4)} + \frac{|\langle \psi_{12} | \hat{H}' | \psi_{+} \rangle|^{2}}{\hbar \omega (2 - 4)} = \frac{\hbar \lambda^{2}}{4m^{2}\omega^{3}} \left(\frac{|1|^{2}}{-2} + \frac{|1|^{2}}{-2}\right) = -\frac{\hbar \lambda^{2}}{4m^{2}\omega^{3}}$$
(21)

$$|\psi_{-}^{(1)}\rangle = \frac{\langle \psi_{21}|\hat{H}'|\psi_{-}\rangle}{\hbar\omega(2-4)}|\psi_{21}\rangle + \frac{\langle \psi_{12}|\hat{H}'|\psi_{-}\rangle}{\hbar\omega(2-4)}|\psi_{12}\rangle = \frac{\lambda}{2m\omega^2}\left(\frac{-1}{-2}|\psi_{21}\rangle + \frac{1}{-2}|\psi_{12}\rangle\right) = \frac{\lambda}{4m\omega^2}\left(|\psi_{21}\rangle - |\psi_{12}\rangle\right)$$
(22)

$$|\psi_{+}^{(1)}\rangle = \frac{\langle \psi_{21}|\hat{H}'|\psi_{+}\rangle}{\hbar\omega(2-4)}|\psi_{21}\rangle + \frac{\langle \psi_{12}|\hat{H}'|\psi_{+}\rangle}{\hbar\omega(2-4)}|\psi_{12}\rangle = \frac{\lambda}{2m\omega^{2}}\left(\frac{1}{-2}|\psi_{21}\rangle + \frac{1}{-2}|\psi_{12}\rangle\right) = -\frac{\lambda}{4m\omega^{2}}\left(|\psi_{21}\rangle + |\psi_{12}\rangle\right)$$
(23)

Exact solution By changing coordinates, we can solve this problem exactly! If we use rotated coordinates u, v so that,

$$u = \frac{1}{\sqrt{2}}(x+y), \qquad v = \frac{1}{\sqrt{2}}(x-y),$$
 (24)

$$x = \frac{1}{\sqrt{2}}(u+v), \qquad y = \frac{1}{\sqrt{2}}(u-v),$$
 (25)

the perturbed Hamiltonian becomes.

$$\hat{H} = \hat{H}_0 + \hat{H}' = \frac{1}{2m} (\hat{p}_u^2 + \hat{p}_v^2) + \frac{1}{4} m\omega^2 \left((u+v)^2 + (u-v)^2 \right) + \lambda (u+v)(u-v)$$

$$= \frac{1}{2m} (\hat{p}_u^2 + \hat{p}_v^2) + \frac{1}{2} m\omega^2 \left[\left(1 + \frac{\lambda}{m\omega^2} \right) u^2 + \left(1 - \frac{\lambda}{m\omega^2} \right) v^2 \right], \tag{26}$$

so we have now have a non-isotropic two-dimensional harmonic oscillator. In the *u*-direction, $\omega_u = \omega \sqrt{1 + \lambda/(m\omega^2)}$, and in the *v*-direction $\omega_v = \omega \sqrt{1 - \lambda/(m\omega^2)}$. This means that the exact energies and their power series expansion are,

$$E_{n_u,n_v} = \hbar\omega \left[\sqrt{1 + \frac{\lambda}{m\omega^2}} \left(n_u + \frac{1}{2} \right) + \sqrt{1 - \frac{\lambda}{m\omega^2}} \left(n_v + \frac{1}{2} \right) \right]$$
 (27)

$$= \hbar\omega (1 + n_u + n_v) + \lambda \frac{\hbar(n_u - n_v)}{2m\omega} - \lambda^2 \frac{\hbar(1 + n_u + n_v)}{8m^2\omega^3} + O(\lambda^3), \tag{28}$$

where $(n_u, n_v) = (0, 0), (1, 0), (0, 1), (1, 1), \ldots$ are the quantum numbers. If we substitute quantum numbers $(n_u, n_v) = (0, 1)$ or $(n_u, n_v) = (1, 0)$ into the power series expansion, we see that we successfully recover our perturbative approximations!

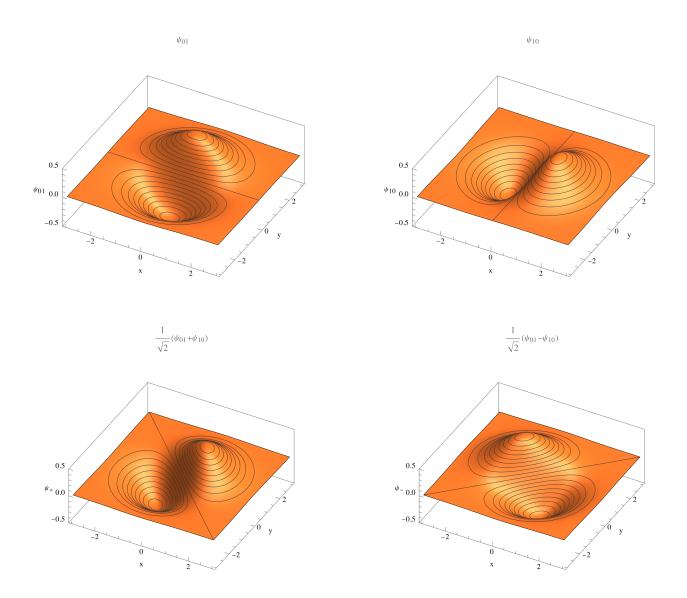


Figure 1: Number states $|\psi_{01}\rangle$, $|\psi_{10}\rangle$ and 'good' states $|\psi_{+}\rangle$, $|\psi_{-}\rangle$ in the position representation.