

PROBLEM SET 6 SOLUTIONS

WEBWORK:

①

(4 points) [setPS6/prob1.pg](#)

Consider the first excited state $|1\rangle$ of a harmonic oscillator with mass m and frequency ω . If we add a perturbation $H' = \lambda x^2$ to the Hamiltonian, the first order shift in the energy of the state (i.e. the correction at order λ) is $\lambda\hbar/(m\omega)$ times

The shift is $\langle 1 | H' | 1 \rangle$

$$= \lambda \langle 1 | x^2 | 1 \rangle$$

$$= \lambda \frac{\hbar}{2m\omega} \langle 1 | (a + a^\dagger)^2 | 1 \rangle$$

We have: $(a + a^\dagger)|1\rangle = |0\rangle + \sqrt{2}|2\rangle$ so

$$\begin{aligned} \langle 1 | (a + a^\dagger)(a + a^\dagger) | 1 \rangle &= (\langle 0 | + \sqrt{2}\langle 2 |)(|0\rangle + \sqrt{2}|2\rangle) \\ &= 1 + 2 = 3 \end{aligned}$$

Then $\Delta E = \lambda \frac{\hbar}{m\omega} \cdot \frac{3}{2}$.

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Consider the third excited state $|3\rangle$ of a harmonic oscillator. If we add a perturbation $H' = \lambda x^3$, the first order shift in the energy of the state (i.e. the correction at order λ) is .

The state shifts to

$$|3\rangle + \lambda|\delta\psi\rangle + \dots$$

where $|\delta\psi\rangle$ can be written as a superposition of energy eigenstates of the original harmonic oscillator:

$$|\delta\psi\rangle = \sum_n c_n |n\rangle$$

For how many n is the coefficient c_n nonzero?

$$\text{We have } \Delta E = \langle 3 | H' | 3 \rangle$$

$$= \lambda \langle 3 | x^3 | 3 \rangle$$

$$= \lambda \langle 3 | \left(\sqrt{\frac{k}{2m\omega}} \right)^3 (a + a^\dagger)^3 | 3 \rangle$$

But $(a + a^\dagger)^3 | 3 \rangle$ is a linear combination of $|6\rangle, |4\rangle, |2\rangle$, and $|0\rangle$, so the inner product with $\langle 3 |$ vanishes and $\Delta E = 0$.

For the state shift, we have:

$$c_n = \frac{\langle n | H' | 3 \rangle}{E_3 - E_n}$$

From the previous part, this will be nonzero for $n = 0, 2, 4, 6$, so four states contribute.

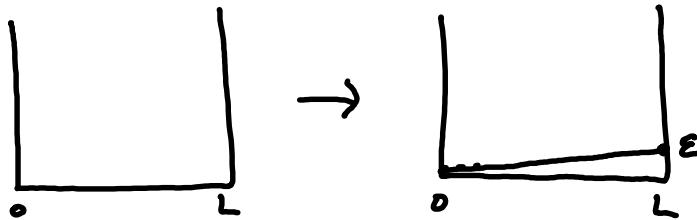
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The energy eigenstates for a particle in an infinite square well potential between $x = 0$ and $x = L$ have wavefunctions

$$\psi_n(x) = \sqrt{\frac{2}{L}} \sin\left(\frac{n\pi x}{L}\right).$$

Consider the $n = 37$ energy eigenstate. If we now modify the potential slightly by making it increase linearly from 0 at $x = 0$ to some potential energy ϵ at $x = L$, the change in energy of this state to leading order in ϵ will be ϵ times



For the n -th eigenstate, we have

$$\Delta E = \langle n | H' | n \rangle$$

Here $H' = \tilde{V}(x)$ where $\tilde{V}(x)$ is a function

$$\tilde{V}(x) = \begin{cases} \epsilon x/L & 0 \leq x \leq L \\ 0 & \text{elsewhere} \end{cases}$$

$$\begin{aligned} \text{Then } \langle n | H' | n \rangle &= \int_0^L dx \psi_n^*(x) \tilde{V}(x) \psi_n(x) \\ &= \frac{2}{L} \int_0^L dx \sin^2\left(\frac{n\pi x}{L}\right) \cdot \epsilon x/L \end{aligned}$$

$$\begin{aligned} \text{set } x = y \cdot L &= 2 \epsilon \int_0^1 dy \sin^2(n\pi y) \cdot y \\ &= \frac{1}{2} \epsilon \end{aligned}$$

2a) We need to compute

$$\Delta E_1 = \langle 1 | H_1 | 1 \rangle$$

$$= \frac{1}{100} \frac{\omega}{\hbar} \langle 1 | p^2 x^2 | 1 \rangle$$

We can write $p = -i \sqrt{\frac{\hbar \omega}{2}} (a - a^\dagger)$ $x = \sqrt{\frac{\pi}{2m\omega}} (a + a^\dagger)$
so $p^2 x^2 = -\frac{\hbar^2}{4} (a - a^\dagger)^2 (a + a^\dagger)^2$

So $\Delta E_1 = -\frac{\hbar \omega}{400} \cdot \langle 1 | (a - a^\dagger)^2 (a + a^\dagger)^2 | 1 \rangle$

Using $a|n\rangle = \sqrt{n}|n-1\rangle$, $a^\dagger|n\rangle = \sqrt{n+1}|n+1\rangle$, we get:

$$(a + a^\dagger)^2 |1\rangle = (a + a^\dagger) (|0\rangle + \sqrt{2}|2\rangle) = 3|1\rangle + \sqrt{6}|3\rangle$$

$$(a - a^\dagger)^2 |1\rangle = (a - a^\dagger) (|0\rangle - \sqrt{2}|2\rangle) = -3|1\rangle + \sqrt{6}|3\rangle$$

$\langle 1 | (a - a^\dagger)^2 (a + a^\dagger)^2 | 1 \rangle$ is the inner product between
these two so equals $-9 + 6 = -3$.

We conclude that the first order energy shift is

$$\Delta E_1 = \frac{3}{400} \hbar \omega$$

b) The correction to the state is:

$$\sum_{n \neq 1} |n\rangle \frac{\langle n | H_1 | 1 \rangle}{E_1 - E_n}$$

We have: $H_1 |1\rangle = -\frac{\hbar \omega}{400} (a - a^\dagger)^2 (a + a^\dagger)^2 |1\rangle$

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$$\begin{aligned}
& - \frac{\hbar\omega}{400} (a - a^\dagger)^2 \left[3|1\rangle + \sqrt{6}|3\rangle \right] \\
& = - \frac{\hbar\omega}{400} (a - a^\dagger) \left[3|0\rangle - 2\sqrt{6}|4\rangle \right] \\
& = - \frac{\hbar\omega}{400} \left[-3|1\rangle - 4\sqrt{6}|3\rangle + 2\sqrt{30}|5\rangle \right]
\end{aligned}$$

So $\langle 3 | H_1 | 1 \rangle = \frac{\sqrt{6}}{100} \hbar\omega$

$$\langle 5 | H_1 | 1 \rangle = - \frac{\sqrt{30}}{200} \hbar\omega$$

and all other matrix elements in our sum vanish.

$$\begin{aligned}
\text{Thus: } \delta |\psi_1\rangle &= |3\rangle \cdot \underbrace{\frac{\langle 3 | H_1 | 1 \rangle}{E_1 - E_3}}_{\approx 2\hbar\omega} + |5\rangle \cdot \underbrace{\frac{\langle 5 | H_1 | 1 \rangle}{E_1 - E_5}}_{\approx 4\hbar\omega} \\
&= \frac{\sqrt{6}}{200} \cdot |3\rangle - \frac{\sqrt{30}}{800} \cdot |5\rangle
\end{aligned}$$

Adding a perturbation $\lambda H_1 = \lambda p^3$, the first order shift in the ground state energy vanishes, since $\langle 0 | p^3 | 0 \rangle = 0$. The second order shift is:

$$\begin{aligned}\delta E_2 &= \lambda^2 \sum_{n \neq 0} \frac{|\langle n | p^3 | 0 \rangle|^2}{E_0 - E_n} \\ &= \lambda^2 \left(\sqrt{\frac{\hbar \omega m}{2}} \right)^6 \sum_{n \neq 0} \frac{|\langle n | (a - a^\dagger)^3 | 0 \rangle|^2}{E_0 - E_n}\end{aligned}$$

$$\begin{aligned}\langle (a - a^\dagger)^3 | 0 \rangle &= -(a - a^\dagger)^2 | 1 \rangle \\ &= (a - a^\dagger) (-| 0 \rangle + \sqrt{2}| 2 \rangle) \\ &= 3| 1 \rangle - \sqrt{6}| 3 \rangle\end{aligned}$$

$$S_o \quad \langle 1 | (a - a^\dagger)^3 | 0 \rangle = 3 \text{ and } \langle 3 | (a - a^\dagger)^3 | 0 \rangle = -\sqrt{6}$$

$$\begin{aligned}\text{Then } \delta E_2 &= \lambda^2 \cdot \left(\frac{\hbar \omega m}{2} \right)^3 \left\{ \frac{9}{(-\hbar \omega)} + \frac{6}{(-3\hbar \omega)} \right\} \\ &= -\frac{11}{8} \lambda^2 \cdot \hbar^2 \omega^2 \cdot m^3\end{aligned}$$