

Jan 2008

Problem 1: The Infinite Square Well: (10 Points)

A single particle is in a one dimensional infinitely deep potential well of width L where $V(x)$ is given by:

$$V(x) = \begin{cases} 0, & \text{if } 0 \leq x \leq L \\ \infty, & \text{otherwise} \end{cases}$$

1. Find the allowed energies (E_n) and the normalized eigenfunctions ($\Psi(x)$) to Schrodinger's Equation for this potential. Show all your work. **(2 Points)**
2. Sketch the wave functions for the first three stationary states for this potential. **(2 Points)**
3. Now, if four spin-1/2 identical particles of mass m are placed in this potential, calculate the three lowest values for the total energy of the system of particles. **(3 Points)**
4. Determine the degeneracy for each of the three energy states found in part 3. **(3 Points)**

Jan 2008

Quantum #1

a) $H\psi = E\psi$

$$-\frac{\hbar^2}{2m} \frac{\partial^2}{\partial x^2} \psi + 0(\psi) = E\psi$$

$$\frac{\partial^2}{\partial x^2} \psi = \frac{2mE}{-\hbar^2} \psi$$

* if $k = \sqrt{\frac{2mE}{\hbar^2}}$

$$\frac{\partial^2}{\partial x^2} \psi = -k^2 \psi$$

$$\hookrightarrow \psi = A \sin(kx) + B \cos(kx)$$

* We know that $\psi(0) = \psi(L) = 0$

$$\hookrightarrow 0 = A \sin(k \cdot 0) + B \cos(k \cdot 0)$$

$$0 = B$$

$$\Rightarrow \psi(x) = A \sin(kx); \text{ for this to be } 0 \text{ at } x=L, kx = n\pi \Rightarrow k_n = \frac{n\pi}{L}$$

$$\hookrightarrow \psi(x) = A \sin\left(\frac{n\pi x}{L}\right)$$

* Normalizing the wave function, we see:

$$1 = \int_{-\infty}^{\infty} |A \sin\left(\frac{n\pi x}{L}\right)|^2 dx$$

$$1 = A^2 \int_0^L \sin^2\left(\frac{n\pi x}{L}\right) dx$$

$$1 = A^2 \cdot \frac{L}{2}$$

$$\hookrightarrow A = \sqrt{\frac{2}{L}}$$

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

* Returning to k_n :

$$k_n = \frac{n\pi}{L} = \sqrt{\frac{2mE}{\hbar^2}}$$

$$\frac{n^2 \pi^2}{L^2} = \frac{2mE}{\hbar^2}$$

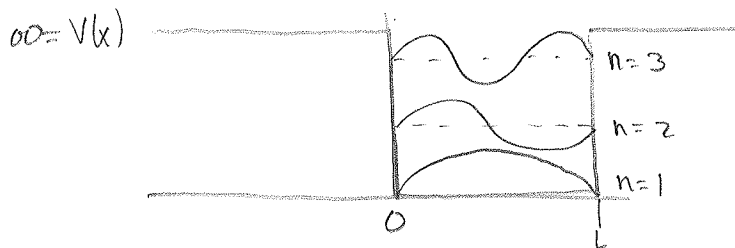
$$\hookrightarrow \boxed{E_n = \frac{n^2 \pi^2 \hbar^2}{2mL^2}}, n \in \mathbb{Z}^+ \text{ for non-trivial solutions}$$

#1 (cont.)

b) $\psi_1 = \sqrt{\frac{2}{L}} \sin\left(\frac{\pi x}{L}\right)$

$\psi_2 = \sqrt{\frac{2}{L}} \sin\left(\frac{2\pi x}{L}\right)$

$\psi_3 = \sqrt{\frac{2}{L}} \sin\left(\frac{3\pi x}{L}\right)$



c) * Since spin-1/2 particles are fermions, no more than one particle can occupy a single state

$$\hookrightarrow E_{\text{sys}} = \frac{(n_1^2 + n_2^2 + n_3^2 + n_4^2) \pi^2 \hbar^2}{2mL^2}$$

\Rightarrow our lowest energy configurations are:

$n = \{1, 2, 3, 4\}, E_{\text{sys}} = \frac{30\pi^2 \hbar^2}{2mL^2}$

$n = \{1, 2, 3, 5\}, E_{\text{sys}} = \frac{39\pi^2 \hbar^2}{2mL^2}$

$n = \{1, 2, 4, 5\}, E_{\text{sys}} = \frac{46\pi^2 \hbar^2}{2mL^2}$

(1)(2) (4) (5)

$1 + 4 + 16 + 25 = 46$

(1) (2) (3) (6)

$1 + 4 + 9 + 36 = 50$

d) Each state has 4! degeneracies, 24 overall for each state

Jan 2006

Problem 2: The Harmonic Oscillator (10 Points):

The normalized wave functions for the one-dimensional quantum harmonic oscillator can be written as,

$$\Psi_n(x) = \left(\frac{\sqrt{\alpha}}{2^n n! \sqrt{\pi}} \right)^{1/2} e^{-\alpha x^2/2} H_n(\sqrt{\alpha}x),$$

where n is the principle quantum number of the oscillator, H_n is the n^{th} order Hermite polynomial, $\alpha = \omega m/\hbar$, ω is the oscillator frequency, and m is its mass. The following equations may be useful,

$$H_{n+1}(q) + 2nH_{n-1}(q) - 2qH_n(q) = 0$$

$$\frac{dH_n(q)}{dq} = 2nH_{n-1}(q)$$

and

$$\begin{aligned}\langle H_n | q H_{n+1} \rangle &= 2^n (n+1)! \sqrt{\pi} \\ \langle H_n | q H_n \rangle &= 0 \\ \langle H_n | q H_{n-1} \rangle &= 2^{n-1} n! \sqrt{\pi}\end{aligned}$$

1. Calculate the expectation value of x and x^2 for the n^{th} state of the harmonic oscillator, where x is the position. **(2 Points)**
2. Calculate the expectation value of p and p^2 for the n^{th} state of the harmonic oscillator, where p is the momentum. **(2 Points)**
3. Calculate Δx and Δp for the n^{th} state. What is the uncertainty product ($\Delta x \Delta p$) for the oscillator? **(2 Points)**
4. Calculate the expectation value of the kinetic energy and the potential energy of the n^{th} state of the oscillator. Show that the sum of the expectation value of the kinetic and potential energies are equal to the total energy of the n^{th} state. **(2 Points)**
5. How does the uncertainty principle relate to the fact that the energy is not zero in the ground state? Explain and interpret your answer to receive credit. **(2 Points)**

Jan 2008

Quantum #2

a) Given: $\Psi_n(x) = \left(\frac{\sqrt{a}}{2^n n! \sqrt{\pi}} \right)^{1/2} \exp\left[-\frac{1}{2}ax^2\right] H_n(\sqrt{a}x)$

Find: $\langle x \rangle_n, \langle x^2 \rangle_n$

* Using raising/lowering operators, we know:

$$x = \sqrt{\frac{\hbar}{2m\omega}} (a + a^\dagger)$$

$$a|\psi_n\rangle = \sqrt{n}|\psi_{n-1}\rangle$$

$$\langle \psi_n | \psi_m \rangle = \delta_{nm}$$

$$p = -i\sqrt{\frac{\hbar m\omega}{2}} (a - a^\dagger)$$

$$a^\dagger|\psi_n\rangle = \sqrt{n+1}|\psi_{n+1}\rangle$$

$$\Rightarrow \langle x \rangle_n = \langle \psi_n | x | \psi_n \rangle$$

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

$$= \sqrt{\frac{\hbar}{2m\omega}} [\langle \psi_n | a | \psi_n \rangle + \langle \psi_n | a^\dagger | \psi_n \rangle]$$

$$= \sqrt{\frac{\hbar}{2m\omega}} [\langle \psi_n | \sqrt{n} | \psi_{n-1} \rangle + \langle \psi_n | \sqrt{n+1} | \psi_{n+1} \rangle]$$

$$= 0$$

$$\langle x^2 \rangle_n = \langle \psi_n | x^2 | \psi_n \rangle$$

$$= \langle \psi_n | \left(\frac{\hbar}{2m\omega} \right) (aa + aa^\dagger + a^\dagger a + a^\dagger a^\dagger) | \psi_n \rangle$$

$$= \frac{\hbar}{2m\omega} [\langle \psi_n | aa | \psi_n \rangle + \langle \psi_n | aa^\dagger | \psi_n \rangle + \langle \psi_n | a^\dagger a | \psi_n \rangle + \langle \psi_n | a^\dagger a^\dagger | \psi_n \rangle]$$

$$= \frac{\hbar}{2m\omega} [\langle \psi_n | \sqrt{n}\sqrt{n-1} | \psi_{n-2} \rangle + \langle \psi_n | \sqrt{n+1}\sqrt{n} | \psi_n \rangle + \langle \psi_n | \sqrt{n}\sqrt{n} | \psi_n \rangle + \langle \psi_n | \sqrt{n+1}\sqrt{n+2} | \psi_{n+2} \rangle]$$

$$= \frac{\hbar}{2m\omega} [2n+1]$$

b) Similarly to above:

$$\langle p \rangle_n = \langle \psi_n | p | \psi_n \rangle$$

$$= \langle \psi_n | -i\sqrt{\frac{\hbar m\omega}{2}} (a - a^\dagger) | \psi_n \rangle$$

$$= -i\sqrt{\frac{\hbar m\omega}{2}} [\langle \psi_n | a | \psi_n \rangle - \langle \psi_n | a^\dagger | \psi_n \rangle]$$

$$= -i\sqrt{\frac{\hbar m\omega}{2}} [\langle \psi_n | \sqrt{n} | \psi_{n-1} \rangle - \langle \psi_n | \sqrt{n+1} | \psi_{n+1} \rangle]$$

$$= 0$$

#2 (cont.)

b) $\langle p_n^2 \rangle = \langle \psi_n | p^2 | \psi_n \rangle$

$$= \langle \psi_n | -\frac{\hbar m \omega}{2} (aa - aa^\dagger - a^\dagger a + a^\dagger a^\dagger) | \psi_n \rangle$$

$$= -\frac{\hbar m \omega}{2} [\langle \psi_n | aa | \psi_n \rangle - \langle \psi_n | aa^\dagger | \psi_n \rangle - \langle \psi_n | a^\dagger a | \psi_n \rangle + \langle \psi_n | a^\dagger a^\dagger | \psi_n \rangle]$$

$$= -\frac{\hbar m \omega}{2} [\langle \psi_n | \sqrt{n-1} \sqrt{n} | \psi_{n-2} \rangle - \langle \psi_n | \sqrt{n+1} \sqrt{n+1} | \psi_n \rangle - \langle \psi_n | \sqrt{n} \sqrt{n} | \psi_n \rangle + \langle \psi_n | \sqrt{n+2} \sqrt{n+1} | \psi_{n+2} \rangle]$$

$$= \frac{\hbar m \omega}{2} [2n+1]$$

c) Generally $\Delta A = \sqrt{\langle A^2 \rangle - \langle A \rangle^2}$

$$\Rightarrow \Delta X = \sqrt{\langle X^2 \rangle - \langle X \rangle^2}$$

$$= \sqrt{\frac{\hbar}{2m\omega} [2n+1] - 0^2}$$

$$= \sqrt{\frac{\hbar}{2m\omega} [2n+1]}$$

$$\Delta P = \sqrt{\langle P^2 \rangle - \langle P \rangle^2}$$

$$= \sqrt{\frac{\hbar m \omega}{2} [2n+1] - 0^2}$$

$$= \sqrt{\frac{\hbar m \omega}{2} [2n+1]}$$

$$\Rightarrow \Delta X \Delta P = \sqrt{\frac{\hbar}{2m\omega} [2n+1]} \sqrt{\frac{\hbar m \omega}{2} [2n+1]}$$

$$= \frac{\hbar}{2} [2n+1]$$

d) $\langle T \rangle = \langle \psi_n | T | \psi_n \rangle$

$$= \langle \psi_n | \frac{p^2}{2m} | \psi_n \rangle$$

$$= \frac{\hbar \omega}{4} [2n+1]$$

$$\langle U \rangle = \langle \psi_n | U | \psi_n \rangle$$

$$= \langle \psi_n | \frac{1}{2} m \omega^2 x^2 | \psi_n \rangle$$

$$= \frac{\hbar \omega}{4} [2n+1]$$

$\hookrightarrow \langle T \rangle + \langle U \rangle = \frac{\hbar \omega}{2} [2n+1]$ which matches what we know to be the energy of the n^{th} state; $E_n = \hbar \omega (n + 1/2)$

e) * From the above formula, we know $E_0 = \frac{\hbar \omega}{2}$ and that $\Delta X \Delta P = \frac{\hbar}{2}$

\Rightarrow Rewriting the total energy in terms of the uncertainties, we see:

$$\Delta E = \frac{(\Delta P)^2}{2m} + \frac{1}{2} m \omega^2 (\Delta x)^2$$

#2 (cont.)

e) * but $\Delta p = \frac{\hbar}{2\Delta x}$

$$\Rightarrow \Delta E = \frac{\hbar^2}{8m(\Delta x)^2} + \frac{1}{2}m\omega^2(\Delta x)^2$$

$$\frac{d(\Delta E)}{d(\Delta x)} = 0 \quad \text{will give minimum of energy}$$

$$\Rightarrow 0 = \frac{-\hbar^2}{4m(\Delta x)^3} + m\omega^2(\Delta x)$$

$$\frac{\hbar^2}{4m(\Delta x)^3} = m\omega^2(\Delta x)$$

$$\frac{\hbar^2}{4m^2\omega^2} = \Delta x^4 \quad \Rightarrow \quad \Delta x = \sqrt{\frac{\hbar}{2m\omega}}$$

$$\begin{aligned} \Rightarrow \Delta E &= \frac{\hbar^2}{8m} \left(\frac{2m\omega}{\hbar} \right) + \frac{1}{2}m\omega^2 \left(\frac{\hbar}{2m\omega} \right) \\ &= \frac{\hbar\omega}{4} + \frac{\hbar\omega}{4} \\ &= \frac{\hbar\omega}{2} \end{aligned}$$

\Rightarrow The uncertainty principle directly implies a non-zero ground state energy

Jan 2008

Problem 3: The Variational Principle: (10 Points)

If the case where you would like to calculate the ground state energy (E_g) for a system described by the Hamiltonian H but you are unable to solve the Schrodinger equation, the variational principle will give you an upper bound for the ground state energy.

For any normalized function Ψ , the variational principle states:

$$E_g \leq \langle \Psi | H | \Psi \rangle$$

1. (2 Points) Prove the variational principle. i.e show that

$$E_g \leq \langle \Psi | H | \Psi \rangle$$

Hint (Write $\Psi = \sum_n c_n \phi_n$ where ϕ_n are the (unknown) eigenfunctions of H)

Now consider a specific case:

In the x-basis, a one-dimensional operator

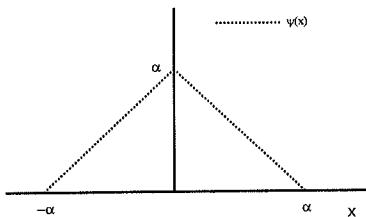
$$\Omega = -\frac{d^2}{dx^2} + |x|$$

has an eigenvalue λ and an eigenfunction $\psi(x)$ with $\psi(x) \rightarrow 0$ for $|x| \rightarrow \infty$.

Let us choose an *unnormalized* trial function

$$\psi(x) = \langle x | \psi \rangle = \begin{cases} \alpha - |x|, & \text{for } |x| < \alpha, \text{ and} \\ 0, & \text{for } |x| > \alpha \end{cases}$$

where α is the variational parameter.



2. (2 Points) Find $\langle \psi | \psi \rangle$.

3. (3 Points) Find the expectation value of the operator Ω .

4. (3 Points) Determine the **best** bound on the lowest eigenvalue (λ) of the operator Ω with the trial function $\psi(x)$. (Note your answer cannot depend on α .)

Jan 2006

Problem 4: Measurement of Hermitian Observables: (10 Points)

Consider a system with three Hermitian observables that are represented in a three-dimensional Hilbert space using the orthonormal basis $|e_1\rangle$, $|e_2\rangle$ and $|e_3\rangle$

with

$$|e_1\rangle = \begin{pmatrix} 1 \\ 0 \\ 0 \end{pmatrix}, |e_2\rangle = \begin{pmatrix} 0 \\ 1 \\ 0 \end{pmatrix}, |e_3\rangle = \begin{pmatrix} 0 \\ 0 \\ 1 \end{pmatrix}$$

and

$$A = \begin{pmatrix} 2 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}, B = \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 2i \\ 0 & -2i & 1 \end{pmatrix}, C = \begin{pmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 1 \end{pmatrix}$$

The system at time $t=0$ is in the state:

$$|\Psi(0)\rangle = \frac{1}{\sqrt{6}}|e_1\rangle - \frac{1}{\sqrt{6}}|e_2\rangle + \sqrt{\frac{2}{3}}|e_3\rangle$$

1. Find the eigenvalues and normalized eigenvectors of B and C . (1 Point)
2. Find the probability of measuring B at time $t = 0$ with the eigenvalue $b = 1$, and then immediately measuring C and finding the eigenvalue $c = 1$, i.e. find $P_{|\Psi(0)\rangle}(b = 1, c = 1)$. (2 Points)
3. Now find the probability if these measurements are performed in reverse order at $t = 0$, i.e. find $P_{|\Psi(0)\rangle}(c = 1, b = 1)$. (2 Points)
4. Are the probabilities obtained in part 1. and part 2. the same or different? Explain in detail. (2 Points)
5. Use the Generalized Uncertainty Principle to determine a lower bound on $\Delta B \Delta C$ for the system in the initial state $|\Psi(0)\rangle$. Discuss your results. (2 Points)
6. Discuss in detail, the conditions that would result in obtaining a lower bound of zero when using the Generalized Uncertainty Principle. Would you expect to get zero for a particular pair of the observables, A , B , and C in this problem? Or for other conditions? (1 Point)

Jan 2008

Quantum #4

a)

$$B = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 2i \\ 0 & -2i & 1 \end{bmatrix}$$

$$\det(B - \lambda I) = 0$$

$$\Rightarrow (1-\lambda)[(1-\lambda)^2 - (2i)(-2i)] = 0$$

$$0 = (1-\lambda)^3 - 4(1-\lambda)$$

$$= [(1 - 2\lambda + \lambda^2) - 4](1-\lambda)$$

$$= (\lambda^2 - 2\lambda - 3)(1-\lambda)$$

$$= (1-\lambda)(\lambda-3)(\lambda+1)$$

$$\Rightarrow \lambda = 1, 3, -1$$

$$C = \begin{bmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 1 \end{bmatrix}$$

$$\det(C - \lambda I) = 0$$

$$0 = -\lambda(-\lambda(1-\lambda) - 0) - 1 \cdot ((1-\lambda) - 0)$$

$$= \lambda^2(1-\lambda) - (1-\lambda)^2$$

$$= [\lambda^2 - (1-\lambda)](1-\lambda)$$

$$\Rightarrow \lambda = 1, 1, -1$$

$$Bx = \lambda x$$

* for $\lambda = 1$

$$\begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 2i \\ 0 & -2i & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix} = \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix}$$

$$x_1 = x_1$$

$$x_2 + 2ix_3 = x_2$$

$$-2ix_2 + x_3 = x_3$$

$$\Rightarrow \vec{x} = \begin{bmatrix} 1 \\ 0 \\ 0 \end{bmatrix}$$

* for $\lambda = 3$

$$\begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 2i \\ 0 & -2i & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix} = 3 \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix}$$

$$x_1 = 3x_1$$

$$x_2 + 2ix_3 = 3x_2$$

$$-2ix_2 + x_3 = 3x_3$$

$$2ix_3 = 2x_2$$

$$ix_3 = x_2$$

$$-2i(ix_3) + x_3 = 3x_3$$

$$-2x_3 + x_3 =$$

$$\Rightarrow \vec{x} = \begin{bmatrix} 0 \\ i \\ 1 \end{bmatrix} \cdot \frac{1}{\sqrt{2}}$$

* for $\lambda = -1$

$$\begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 2i \\ 0 & -2i & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix} = -1 \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix}$$

$$x_1 = -x_1$$

$$x_2 + 2ix_3 = -x_2$$

$$-2ix_2 + x_3 = -x_3$$

$$2ix_3 = -2x_2$$

$$ix_3 = -x_2$$

$$\Rightarrow \vec{x} = \begin{bmatrix} 0 \\ -i \\ 1 \end{bmatrix} \cdot \frac{1}{\sqrt{2}}$$

a) $Cx = \lambda x$

* for $\lambda = -1$

$$\begin{bmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix} = - \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix}$$

$$x_2 = -x_1$$

$$x_1 = -x_2$$

$$x_3 = -x_3$$

$$\Rightarrow \vec{x} = \frac{1}{\sqrt{2}} \begin{bmatrix} -1 \\ 1 \\ 0 \end{bmatrix}$$

* for $\lambda = 1$

$$\begin{bmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix} = \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix}$$

$$x_2 = x_1$$

$$x_1 = x_2$$

$$x_3 = x_3$$

$$\Rightarrow \vec{x} = \begin{bmatrix} 0 \\ 0 \\ 1 \end{bmatrix} \text{ or } \frac{1}{\sqrt{2}} \begin{bmatrix} 1 \\ 1 \\ 0 \end{bmatrix}$$

$|\lambda_c = 1, 2\rangle \quad |\lambda_c = 1, 1\rangle$

b) * Convert $|7(0)\rangle$ into B eigenbasis

$$\begin{aligned} |7(0)\rangle &= \begin{bmatrix} 1/\sqrt{6} \\ -1/\sqrt{6} \\ 2/\sqrt{6} \end{bmatrix} = \frac{1}{\sqrt{6}} \left(|\lambda=1\rangle + \frac{1}{2} [|\lambda=3\rangle - |\lambda=-1\rangle] + [|\lambda=3\rangle + |\lambda=-1\rangle] \right) \\ &= \frac{1}{\sqrt{6}} (|\lambda=1\rangle + (1 + \frac{1}{2})|\lambda=3\rangle + (1 - \frac{1}{2})|\lambda=-1\rangle) \end{aligned}$$

* To find probability

$$\begin{aligned} |\langle \lambda_b=1 | B | 7(0) \rangle|^2 &= |\langle \lambda_b=1 | \left(\frac{1}{\sqrt{6}} [1|\lambda=1\rangle + 3(1+\frac{1}{2})|\lambda=3\rangle - (1-\frac{1}{2})|\lambda=-1\rangle] \right)|^2 \\ &= \frac{1}{6} \end{aligned}$$

$$\begin{aligned} |\langle \lambda_c=1 | C | \lambda_b=1 \rangle|^2 &= |\langle \lambda_c=1 | C | \frac{1}{\sqrt{2}} (|\lambda_c=1,1\rangle - |\lambda_c=-1\rangle) |^2 \\ &= |\langle \lambda_c=1,1 | \frac{1}{\sqrt{2}} (|\lambda_c=1,1\rangle + |\lambda_c=-1\rangle) |^2 \\ &= \frac{1}{2} \end{aligned}$$

* Note: Only need $\langle \lambda_c=1,1 |$ case b/c of orthogonality of eigenkets, i.e. probability 0 in $\langle \lambda_c=1,2 |$ case

Overall probability: $\frac{1}{12}$

#4 (cont.)

c) * Reversing the order from part b, we see:

$$|7(0)\rangle = \begin{bmatrix} 1/\sqrt{6} \\ -1/\sqrt{6} \\ 2/\sqrt{6} \end{bmatrix} = \frac{1}{\sqrt{6}} (2|\lambda_c=1,2\rangle - |\lambda_c=-1\rangle)$$

$$|\langle \lambda_c=1 | C | 7(0) \rangle|^2 = |\langle \lambda_c=1 | \frac{1}{\sqrt{6}} (2|\lambda_c=1,2\rangle + |\lambda_c=-1\rangle)|^2$$

* only need $\langle \lambda_c=1 | = \langle \lambda_c=1,2 |$ case b/c of orthogonality of eigenvectors

$$= \left| \frac{2}{\sqrt{6}} \right|^2 = \frac{1}{4}$$

* Converting $|\lambda_c=1,2\rangle$ to B eigenbasis: $|\lambda_c=1,2\rangle = \begin{bmatrix} 0 \\ 0 \\ 1 \end{bmatrix} = \frac{1}{\sqrt{2}} (|\lambda_b=3\rangle + |\lambda_b=-1\rangle)$

$$|\langle \lambda_b=1 | B | \frac{1}{\sqrt{2}} (|\lambda_b=3\rangle + |\lambda_b=-1\rangle) \rangle|^2 = 0$$

Overall probability: 0

d) The probabilities in parts b + c are different because the two observables are not commutable. They have different eigenbasis and therefore the system is affected in different ways depending upon which operator is acted first

e) $\langle (\Delta B)^2 \rangle \langle (\Delta C)^2 \rangle \geq \frac{1}{4} | \langle [B, C] \rangle |^2$, where $\Delta A = \langle A^2 \rangle - \langle A \rangle^2$

~~$$B^2 = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 2i \\ 0 & -2i & 1 \end{bmatrix} \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 2i \\ 0 & -2i & 1 \end{bmatrix} = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 5 & -4i \\ 0 & 4i & 5 \end{bmatrix}$$

$$C^2 = \begin{bmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 1 \end{bmatrix} \begin{bmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 1 \end{bmatrix} = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix}$$

$$\langle B^2 \rangle = \frac{1}{6} \begin{bmatrix} 1 & -1 & 2 \end{bmatrix} \begin{bmatrix} 1 & 0 & 0 \\ 0 & 5 & -4i \\ 0 & 4i & 5 \end{bmatrix} \begin{bmatrix} 1 \\ -1 \\ 2 \end{bmatrix}$$

$$= \frac{1}{6} \begin{bmatrix} 1 & -1 & 2 \end{bmatrix} \begin{bmatrix} 1 \\ -5-8i \\ 10+4i \end{bmatrix} = \frac{1}{6} (1+5+20+8i-8i) = \frac{26}{6}$$

$$\langle C^2 \rangle = \frac{1}{6} \begin{bmatrix} 1 & -1 & 2 \end{bmatrix} \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix} \begin{bmatrix} 1 \\ -1 \\ 2 \end{bmatrix}$$

$$= \frac{1}{6} \begin{bmatrix} 1 & -1 & 2 \end{bmatrix} \begin{bmatrix} 1 \\ -1 \\ 2 \end{bmatrix} = \frac{1}{6} (6) = 1$$~~

#4 (cont.)

e) Taking square root of above equation yields: $(\Delta B)(\Delta C) \geq \frac{1}{2} |\langle [B, C] \rangle|$

$$\Rightarrow BC = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 2i \\ 0 & -2i & 1 \end{bmatrix} \begin{bmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 1 \end{bmatrix} = \begin{bmatrix} 0 & 1 & 0 \\ 1 & 0 & 2i \\ -2i & 0 & 1 \end{bmatrix}$$

$$CB = \begin{bmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 1 \end{bmatrix} \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 2i \\ 0 & -2i & 1 \end{bmatrix} = \begin{bmatrix} 0 & 1 & 2i \\ 1 & 0 & 0 \\ 0 & -2i & 1 \end{bmatrix}$$

Jan 2008

Problem 5: Perturbation Theory: (10 Points)

A single particle is in a one dimensional infinite well of length L . The potential $V(x)$ is given by:

$$V(x) = \begin{cases} 0, & \text{if } 0 \leq x \leq L \\ \infty, & \text{otherwise} \end{cases}$$

Suppose the potential energy inside the well is changed to

$$V(x) = \epsilon \sin \frac{\pi x}{L}$$

when $0 \leq x \leq L$.

Note you may use your results from Problem 1 for this problem.

1. Calculate the energy shifts for the perturbed well to first order in ϵ . **(2 Points)**
2. Which energy level is shifted the most to first order in ϵ ? **(1 Point)**
3. Calculate the second order (in ϵ) correction to the ground state energy **(2 Points)**
4. Calculate the corrections to the ground state wavefunction to first order in ϵ . **(2 Points)**
5. Suppose that ϵ is larger than the energy of the first excited state. Carefully sketch the wavefunction versus x for the ground state and for the first excited state. How many nodes, maxima, and minima does the wavefunction have in each state. **(2 Points)**
6. Suppose the wavefunction is a linear combination of the ground state and the first excited state from part 5. Describe how the maximum of the probability density depends on time. **(1 Point)**

Jan 2008

Problem 6: Spherically Symmetric States: (10 Points)

Consider eigenfunctions of the Hamiltonian of a particle in a three-dimensional central potential. In particular, consider those eigenfunctions that depend only on the electron's radial coordinate r , that is $\Psi_E = \Psi_E(r)$. States represented by such eigenfunctions are called "spherically symmetric states".

1. Derive an equation for a function $\chi_E(r)$ defined by:

$$\Psi_n(r) \equiv \frac{1}{r} \chi_n(r),$$

where n is the principle quantum number. **(2 Points)**

The remainder of this problem concerns a hydrogen atom in the approximation that we neglect all interactions except the Coulomb interaction and treat the proton as an infinitely massive point particle at the origin.

2. Sketch $\chi_n(r)$ for the lowest three spherical bound states of the hydrogen atom. Justify the qualitative features of each function. **(2 Points)**
3. **(2 Points)**. Consider the eigenfunction for the ground state. Prove that to be physically admissible this function must decay exponentially as r becomes infinite.

$$\chi_1(r) \rightarrow e^{-\alpha r}, \text{ when } r \rightarrow \infty$$

where α is a constant, and that therefore $\chi_1(r)$ must have the form.

$$\chi_1(r) = f(r)e^{-\alpha r}.$$

4. Use $f(r) = r$. Justify why this is an appropriate choice and show that the above equation is a solution of the equation you derived for $\chi_1(r)$ and determine the corresponding eigenvalue E_1 . **(2 Points)**
5. Derive an expression for the constant α in terms of fundamental constants. **(2 Points)**