

Quantum Mechanics Qualifying Exam

Fall 2021

August 18 (9:00 am - 12:00 pm)

1. Consider a physical system with a three-dimensional state space. An orthonormal basis of the state space is chosen. In this basis, the Hamiltonian is represented by the matrix:

$$H = \begin{pmatrix} 2 & 1 & 0 \\ 1 & 2 & 0 \\ 0 & 0 & 3 \end{pmatrix}$$

- (a) What are the possible results when the energy of the system is measured?
- (b) A particle is in the state $|\psi\rangle$ represented in this basis as

$$|\psi\rangle \sim \frac{1}{\sqrt{3}} \begin{pmatrix} i \\ -i \\ i \end{pmatrix}$$

Find the expectation values $\langle H \rangle$, $\langle H^2 \rangle$ and the uncertainty ΔH in this state.

2. A state for a simple harmonic oscillator of frequency ω starts evolving (at $t = 0$) from an arbitrary superposition of two number states

$$|\alpha, 0\rangle = \cos \theta |n\rangle + e^{i\phi} \sin \theta |l\rangle$$

where θ, ϕ are real and $n > l$.

- (a) Derive the state vector at arbitrary time t .
- (b) What is the energy expectation value at time t ? Is it a periodic function of time? If yes, what is the period?
- (c) Calculate the expectation value of potential energy at time t . Is it a periodic function of time? If yes, what is the period?

3. A molecule is rotating around its center of mass. The Hamiltonian is

$$H = \frac{L_x^2 + L_y^2}{2I_a} + \frac{L_z^2}{2I_b}$$

where I_a and I_b are the moments of inertia, and L_x, L_y, L_z are the orbital angular momentum operators.

- (a). Find the energy eigenvalues and eigenstates.

(b) Consider a state described by the normalized angular wave function

$$\psi(\theta, \phi) = \sqrt{\frac{3}{4\pi}} \sin \theta \cos \phi$$

where θ and ϕ the polar and azimuthal angles respectively. Compute the expectation value of L_z in the state ψ .

(c). Suppose L_z is measured in the above state ψ and the result is $+\hbar$. Immediately afterwards, L_x is measured. Find the uncertainty (standard deviation) ΔL_x in this measurement.

4. In a hydrogen atom, the wavefunction $\psi(\mathbf{r})$ describes the motion of the electron relative to the proton (i.e. $\mathbf{r} = \mathbf{r}_e - \mathbf{r}_p$, where \mathbf{r}_e and \mathbf{r}_p are the electron's and proton's coordinates respectively). If the atom is localized at the origin, show that the probability density of the proton is:

$$P_p(\mathbf{r}_p) = \left(\frac{m_e + m_p}{m_e} \right)^3 \left| \psi \left(\frac{m_e + m_p}{m_e} \mathbf{r}_p \right) \right|^2$$

[Hint: Since you are implicitly given $\psi(\mathbf{r})$, you'll need to work with the center-of-mass \mathbf{R} and relative \mathbf{r} coordinates. Start by constructing the expression for P_p in terms of the atom's *two-body* wavefunction. Use the fact that the atom is localized at the origin, i.e. the two-body wavefunction pins \mathbf{R} to zero.]