Two-state systems

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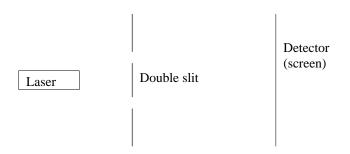
Quantum Mechanics 1 August 11, 2008

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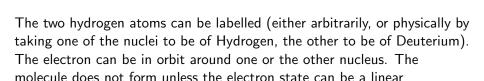
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The two-slit experiment



One can label the slits in many ways: the labelling can be arbitrary or physical). Clearly there is an identifiable quantum state corresponding to the photon coming from slit 1, and another for the photon coming from slit 2. The interference phenomenon is seen whenever the quantum state relevant to the measurement is a superposition of the two.

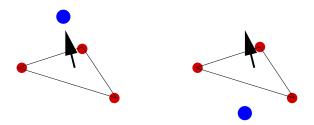
The H_2^+ molecule



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superposition of the two.

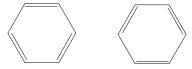
The ammonia molecule



The three hydrogen atoms define a plane. The ammonia molecule has a magnetic moment, and its orientation with respect to this plane defines an "up" and a "down" direction.: the nitrogen atom can be either "above" the plane or "below" it, or in a linear superposition of the two.

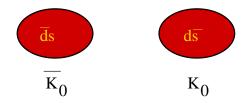
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The benzene molecule



The benzene ring must include both single and double (covalent) bonds between the carbon atoms. There are two possible arrangements of these kinds of bonds (described either by an arbitrary labelling of the carbon atoms or by physically labelling them in some way). Can there be linear superpositions of the electronic states of the valence electrons?

Neutral K meson oscillations



Mesons are made of a quark and an anti-quark. Uncharged strange mesons are called K_0 mesons. Since the strange quark (s) has charge -1/3, the same as a down quark (d), there are two possible kinds of K_0 —either one with a strange anti-quark and a down quark $(d\bar{s})$, called the K_0) or one with a strange quark and a down anti-quark $(\bar{d}s)$, called the \bar{K}_0). Are linear superpositions of these two states allowed?

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The Hilbert space

In all the examples above, we have started with two quantum states (call them $|1\rangle$ and $|2\rangle$) which are orthogonal, and we have claimed that quantum states which are arbitrary superpositions of these two can exist. It is useful to keep in mind a concrete representation of the two states

$$|1\rangle = \begin{pmatrix} 1 \\ 0 \end{pmatrix}, \quad \text{and} \quad |2\rangle = \begin{pmatrix} 0 \\ 1 \end{pmatrix}.$$

Since an overall phase is immaterial to all that follows, we could also have chosen, for example,

$$|1\rangle = rac{1}{\sqrt{3}} egin{pmatrix} 1+2i \ 0 \end{pmatrix}, \qquad ext{and} \qquad |2\rangle = rac{1}{\sqrt{3}} egin{pmatrix} 0 \ 1+2i \end{pmatrix}.$$

In quantum mechanics, these two choices are indistinguishable.

Hermitean operators for two-state systems

Take the most general 2×2 matrix

$$A = \begin{pmatrix} z_1 & z_2 \\ z_3 & z_4 \end{pmatrix}.$$

If $A^+ = A$ then z_1 and z_4 are real and $z_2 = z_3^*$. Hence four real parameters are sufficient to determine the most general Hermitean matrix,

$$H = \begin{pmatrix} a_0 + a_3 & a_1 - ia_2 \\ a_1 + ia_2 & a_0 - a_3 \end{pmatrix}.$$

If A and B are Hermitean, then aA + bB is Hermitean if a and b are real scalars.

Hence any (real) linear combination of the following matrices is Hermitean; the identity matrix 1 and the three Pauli matrices

$$\sigma_1 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \quad \sigma_2 = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}, \quad \sigma_3 = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}.$$

Algebra of Pauli Matrices

- The Pauli matrices are square roots of unity: $\sigma_1^2 = \sigma_2^2 = \sigma_3^2 = 1$.
- Since the Pauli matrices have vanishing trace and their squares are identity, the eigenvalues of each of them must be +1 and -1 (the eigenvalues are square roots of unity).
- The Pauli matrices do not commute with each other—

$$[\sigma_1, \sigma_2] = 2i\sigma_3, \quad [\sigma_2, \sigma_3] = 2i\sigma_1, \quad [\sigma_3, \sigma_1] = 2i\sigma_2.$$

In terms of the completely anti-symmetric symbol (the Levi-Civita symbol) we can write $[\sigma_j, \sigma_k] = 2i\epsilon_{jkl}\sigma_l$. Here we have chosen the convention that $\epsilon_{123} = 1$.

- The Pauli matrices anti-commute with each other: $\{\sigma_i, \sigma_j\} = 2\delta_{ij}$.
- Using the notation $\sigma_0 = 1$, any Hermitean matrix can be written in the form $H = a_i \sigma_i$. There are no restrictions on the components of a_i , and there are no relations between them.

Eigenvectors of Pauli matrices

- The states $|1\rangle$ and $|2\rangle$ are eigenvectors of σ_3 with eigenvalues 1 and -1 respectively.
- Since none of the Pauli matrices commute with each other, they cannot be simultaneously diagonalized.
- The eigenvectors of σ_1 are $|+\rangle$ and $|-\rangle 1$. Hence the unitary transformation that rotates the eigenbasis of σ_3 into that of σ_1 is

$$U_{31}=\frac{1}{\sqrt{2}}\begin{pmatrix}1&-1\\1&1\end{pmatrix}.$$

Check that this indeed diagonalizes σ_1 . What does it do to the other Pauli matrices?

• What are the eigenvectors of σ_2 ? Construct the unitary transformation which diagonalizes σ_2 , and check what it does to the other Pauli matrices.

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Functions of matrices

- If A is a diagonal matrix, and B = f(A), then B is diagonal and $B_{ii} = f(A_{ii})$.
- If A can be diagonalized, then f(A) can be found by first diagonalizing $A \to UAU^{\dagger}$, then constructing $B = f(UAU^{\dagger})$ in this basis, and finally transforming back, $B \to U^{\dagger}BU = U^{\dagger}f(UAU^{\dagger})U$.
- The exponentials of Pauli matrices can easily be found in this way.
 For example,

$$\exp(a\sigma) = \begin{pmatrix} e^a & 0 \\ 0 & e^{-a} \end{pmatrix}.$$

All other Pauli matrices look the same in their eigenbasis. (Exponentiate the other Pauli matrices)

- The eigenvalues of unitary matrices are complex numbers of unit modulus.
- Every unitary matrix, U can be written as exp(iH) where H is a Hermitean matrix.

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Time evolution

- **Planck's hypothesis**, when extended to all quantum states, contains the rule for time evolution of quantum states. All waves evolve through the phase factor $\exp(-i\omega t)$. Since $E=\hbar\omega$, a quantum state of definite energy evolves as $\exp(-iEt/\hbar)$.
- The time evolution of an arbitrary superposition of states must, therefore, be given by $|\psi(t)\rangle = U(t-t_0)|\psi(t_0)\rangle$, where $U(t) = \exp{(-iHt/\hbar)}$, and H is the Hamiltonian operator. The evolution operator is unitary since H is Hermitean.
- Using the definition of matrix exponentials, one sees that the time derivative of U(t) is $-iHU(t)/\hbar$. This gives us **Schrödinger's** equation

$$i\hbar \frac{d}{dt}|\psi\rangle = H|\psi\rangle.$$

The complete solution needs one initial condition, *i.e.*, $|\psi(t_0)\rangle$.

Evolution of density matrices

Since every density matrix can be written in terms of a complete basis in the form

$$\rho = \sum_{ij} \rho_{ij} |i\rangle\langle j|,$$

it's time evolution can be written in the form

$$\rho(t) = U(t - t_0)\rho(t_0)U^{\dagger}(t - t_0).$$

The chain rule of differentiation then gives the evolution equation for the density matrix—

$$i\hbar \frac{d}{dt}\rho(t) = [H, \rho(t)].$$

A complete solution needs the initial density matrix. This equation is called von Neumann's equation.

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The general case

The most general Hamiltonian for a two level system is just the most general Hermitean matrix we had written before: $H = \sum_{i=0}^{3} a_i \sigma_i$. The most general evolution operator is then $U = \exp(-it \sum_i a_i \sigma_i/\hbar)$. Write $\mathbf{a} = (a_1, a_2, a_3) = |a|\hat{\mathbf{n}}$, where $|a|^2 = \mathbf{a} \cdot \mathbf{a}$, and $\hat{\mathbf{n}}$ is a real unit

3-vector. The matrix $s = \hat{\mathbf{n}} \cdot \sigma$ has eigenvalues ± 1 since $s^2 = 1$. (Prove this)

Hence, the energy eigenvalues are $E_{\pm}=a_0\pm |a|$, and

$$U = V^{\dagger} egin{pmatrix} \exp(-iE_{+}t/\hbar) & 0 \ 0 & \exp(-iE_{-}t/\hbar) \end{pmatrix} V,$$

where V is the unitary matrix which diagonalizes s. Writing $\hat{\mathbf{n}} = (\sin \theta \cos \phi, \sin \theta \sin \phi, \cos \theta)$, one finds that

$$s = \begin{pmatrix} \cos \theta & \sin \theta \mathrm{e}^{-i\phi} \\ \sin \theta \mathrm{e}^{i\phi} & \cos \theta \end{pmatrix}, \qquad V = \begin{pmatrix} \cos(\theta/2) \mathrm{e}^{i\phi} & \sin(\theta/2) \\ -\sin \theta & \cos(\theta/2) \mathrm{e}^{-i\phi} \end{pmatrix}.$$

(Show this. What is the corresponding expression for U?)

Time evolution

If an initial quantum state is

$$|\psi(t_0)\rangle = \cos \alpha |E_+\rangle + \sin \alpha e^{i\beta} |E_-\rangle,$$

where the two basis states on the right are eigenstates of the Hamiltonian, then, after time evolution one has

$$|\psi(t)\rangle = \cos \alpha e^{-iE_{+}t/\hbar}|E_{+}\rangle + \sin \alpha e^{i(\beta - E_{-}t/\hbar)}|E_{-}\rangle.$$

The squared amplitude therefore oscillates in time.

How do you apply this argument to the double slit experiment?

A problem

The most general 2×2 unitary matrix is

$$U = \begin{pmatrix} z_1 & w_1 \\ z_2 & w_2 \end{pmatrix}, \quad \text{with} \quad U^{\dagger}U = 1 = UU^{\dagger}.$$

Choosing $z_1 = \cos\theta \exp(i\phi)$, $z_2 = \sin\theta \exp(i\phi')$, $w_1 = \cos\psi \exp(i\alpha)$, $w_2 = \sin\psi \exp(i\alpha')$ satisfies the normalization constraints from $U^{\dagger}U$. The normalization constraints from UU^{\dagger} further imply that $\cos^2\phi = \sin^2\theta$.

- Complete the evaluation using the normalization constraints.
- 2 What is the determinant of *U*?
- \odot Note that any overall phase of wavefunctions does not matter. Can this be used to restrict the class of unitary transformations required in quantum mechanics? (For example by putting a condition on the determinant of U)
- Compare the parametrization of a general complex matrix made here with the unitary evolution operator and find when the two are the same.

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References

- The Feynman Lectures in Physics (Vol 3), by R. P. Feynman *et al*. Examples pf two-state systems are discussed in detail in chapters 9, 10 and 11 of this book.
- Quantum Mechanics (Non-relativistic theory), by L. D. Landau and E. M. Lifschitz. The material on Pauli matrices can be found in chapter 8 of this book. There are also a couple of useful problems regarding the manipulation of the Pauli matrices.
- Quantum Mechanics (Vol 1), C. Cohen-Tannoudji, B. Diu and F. Laloë. Chapter 3 of this book discusses several two-state systems in the kind of detail that is the hallmark of this book.
- Mathematical Methods for Physicists, by G. Arfken. This book contains chapters on matrices and vector spaces which will be useful throughout this course.