

$$\Psi_n(x, t) = f_n(x) g_n(t) = \sqrt{\frac{2}{a}} \sin\left(\frac{n\pi x}{a}\right) \exp\left[\frac{-it}{\hbar} E_n\right]$$

does have a unique energy = E_n

example $c_1 = \sqrt{\frac{1}{2}} = c_2$, all others $c_n = 0$

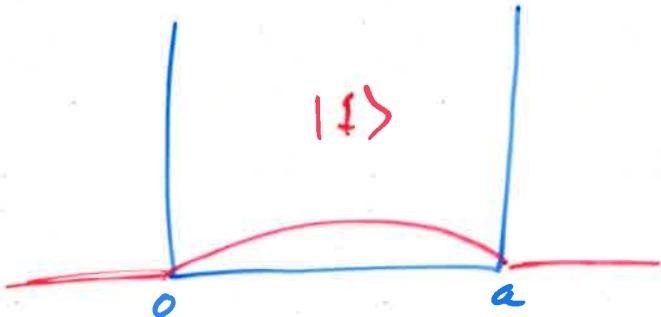
electron confined to $1\text{Å} = 10^{-10}\text{m}$

$$E_1 = 6.0 \times 10^{-18}\text{J}$$

1kg confined to 7m

$$E_1 = 1.1 \times 10^{-69}\text{J}$$

$$|1\rangle = \sqrt{\frac{2}{a}} \sin\left(\frac{2\pi x}{a}\right) \exp\left[-\frac{it}{\hbar}\left(\frac{4\pi^2 \hbar^2}{2ma^2}\right)\right]$$



$$|2\rangle = \sqrt{\frac{2}{a}} \sin\left(\frac{2\pi x}{a}\right) \exp\left[-\frac{it}{\hbar}\left(\frac{4\pi^2 \hbar^2}{2ma^2}\right)\right]$$

$$E_1 = \hbar \omega_1 = h \nu_1 = h f_1 \quad \hbar = \frac{h}{2\pi}$$

↑ ↑ ↑
 angular frequency linear frequency $\omega = 2\pi f$
 MKS $\rightarrow \text{rad/sec}$ $f_2 = \frac{\text{cycles}}{\text{sec}}$ $= 2\pi \nu$

Like guitar string $\lambda_2 = \frac{1}{2} \lambda_1$, $\lambda = \text{wavelength}$

Unlike a guitar string: $\nu_2 = 4\nu_1$,
 $\nu_3 = 9\nu_1$

Really unlike a guitar string $E_2 = 4E_1$,

guitar string has any energy you like.

$$\langle 1|1 \rangle = 1 = \int_{x=0}^{x=a} \sqrt{\frac{2}{a}} \sin\left(\frac{\pi x}{a}\right)^* e^{\frac{+itE_1}{\hbar}} \sqrt{\frac{2}{a}} \sin\left(\frac{\pi x}{a}\right) e^{\frac{-itE_1}{\hbar}} dx$$

$$= \frac{2}{a} \int_{x=0}^{x=a} \sin^2\left(\frac{\pi x}{a}\right) dx = 1 \quad \checkmark$$

$$\langle 2|1 \rangle = 0 \quad \langle 2|2 \rangle = 1$$

$$\langle n|p \rangle = \delta_{np}$$

$|1\rangle, |2\rangle, |3\rangle \dots |n\rangle \dots$ basis vectors for Hilbert space

Linear Superposition

$$|\Psi\rangle = \Psi_{(x,t)} = \sum_{n=1}^{\infty} c_n |n\rangle \quad \text{in general}$$

special combination

$$|\Psi\rangle = \underbrace{\frac{1}{\sqrt{2}}}_{c_1} |1\rangle + \underbrace{\frac{1}{\sqrt{2}}}_{c_2} |2\rangle$$

$$c_1 = \frac{1}{\sqrt{2}} = c_2 \\ \text{all other} \\ c_n = 0$$

restriction on coefficients: $\sum_{n=1}^{\infty} |c_n|^2 = 1$

$$|c_1|^2 + |c_2|^2 = \frac{1}{2} + \frac{1}{2} = 1 \quad \checkmark$$

$|\Psi\rangle$ is normalized

$$\begin{aligned}\langle \Psi | \Psi \rangle &= \left(\frac{1}{\sqrt{2}} \langle 1 | + \frac{1}{\sqrt{2}} \langle 2 | \right) \left(\frac{1}{\sqrt{2}} | 1 \rangle + \frac{1}{\sqrt{2}} | 2 \rangle \right) \\ &= \underbrace{\frac{1}{2} \langle 1 | 1 \rangle}_{1} + \underbrace{\frac{1}{2} \langle 2 | 2 \rangle}_{1} + \underbrace{\frac{1}{2} \langle 1 | 2 \rangle}_{0} + \underbrace{\frac{1}{2} \langle 2 | 1 \rangle}_{0} = 1.\end{aligned}$$

Expectation Value of the energy \rightarrow Hamiltonian operator

$$\langle \hat{H} \rangle = \langle \Psi | \hat{H} | \Psi \rangle = \langle \Psi | i\hbar \frac{\partial}{\partial t} | \Psi \rangle$$

$$\hat{H} | 1 \rangle = i\hbar \frac{\partial}{\partial t} | 1 \rangle = E_1 | 1 \rangle$$

↑ eigen vector
 matrix ↑ eigen value ↑ eigen vector

In this basis $|1\rangle, |2\rangle, |3\rangle, \dots |n\rangle$

$$\hat{H} = \begin{pmatrix} E_1 & 0 & 0 & \dots \\ 0 & E_2 & 0 & \dots \\ 0 & 0 & E_3 & \dots \\ \vdots & \vdots & \vdots & \ddots E_n \end{pmatrix} \quad \text{diagonal matrix}$$

$$H_{np} = \langle n | \hat{H} | p \rangle = E_n \delta_{np}$$

$$\hat{H} | 2 \rangle = E_2 | 2 \rangle \quad \hat{n} | n \rangle = E_n | n \rangle$$

The states $|1\rangle, |2\rangle \dots |n\rangle \dots$ are eigenstates (eigenkets) of the Hamiltonian operator. They are also called stationary states. They have definite energy E_n .

$$\begin{aligned}\langle -\hat{E} | \hat{H} | \psi \rangle &= \left(\frac{1}{\sqrt{2}} \langle 1 | + \frac{1}{\sqrt{2}} \langle 2 | \right) \hat{H} \left(\frac{1}{\sqrt{2}} | 1 \rangle + \frac{1}{\sqrt{2}} | 2 \rangle \right) \\ &= \frac{1}{2} E_1 \langle 1 | 1 \rangle + \frac{1}{2} E_2 \langle 2 | 2 \rangle + \cancel{\frac{1}{2} E_1 \langle 2 | 1 \rangle^{\circ}} \\ &\quad + \cancel{\frac{1}{2} E_2 \langle 1 | 2 \rangle^{\circ}} \\ &= \underline{\frac{1}{2} (E_1 + E_2)} = \frac{1}{2} \left(\frac{\pi^2 \hbar^2}{2ma^2} \right) (1 + 4) = \frac{5}{2} \frac{\pi^2 \hbar^2}{2ma^2}\end{aligned}$$

Meaning of expectation value

Prepare a large number ~ 1 billion identical systems, measure the energy.

~ 500 million you will get E_1

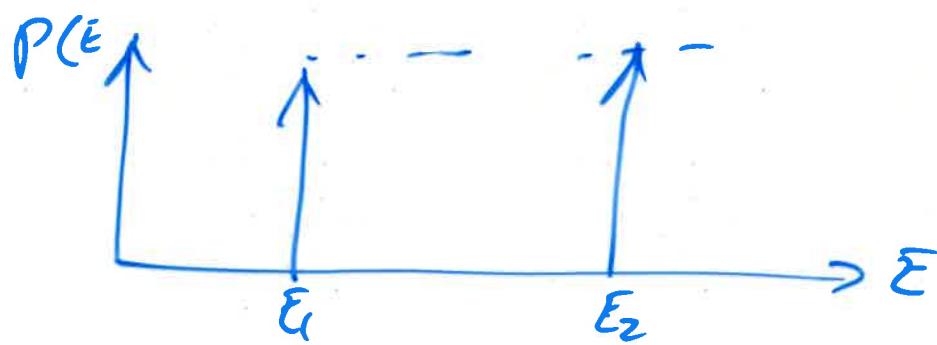
~ 500 million you will get E_2

Average is $\frac{500 \text{ million } E_1 + 500 \text{ million } E_2}{1 \text{ billion}} = \frac{E_1 + E_2}{2}$

but NONE of the measurements will ever give

$$\frac{\sum \pi^2 \hbar^2}{2ma^2}$$

Most likely energy measurement?
 E_1 and E_2 are both equally likely



After a measurement in which E_1 is obtained, the wavefunction collapses to $|1\rangle$ only. Stays $|1\rangle$ forever.

Measure energy and get E_2 , the wavefunction is pure $|2\rangle$. Stays $|2\rangle$ forever.

Measure x

$x|1\rangle$

Expectation value

$$\langle 1 | \chi | 1 \rangle = \frac{a}{2}$$

$$\langle 2 | \chi | 2 \rangle = \frac{a}{2}$$

$$\langle 1 | \chi | 2 \rangle = ?$$

$$\langle \mathcal{E} | \chi | \mathcal{E} \rangle = ?$$

Expect "beats"
sinusoidal time dependence