Once again, the Schrödinger equation:

\[
\frac{i\hbar}{\partial t} \frac{\partial \psi(x,t)}{\partial t} = -\frac{\hbar^2}{2m} \frac{\partial^2 \psi(x,t)}{\partial x^2} + V(x,t)
\]

(which can also be written \( \hat{H} \psi(x,t) \) if you like.)

And once again, assume \( V = V(x) \) (no \( t \) in there!)

We can start to solve the PDE by **SEPARATION OF VARIABLES**.

Assume (hope? wonder if?) we might find a solution of form \( \Psi (x,t) = u(x) \phi(t) \).

Griffiths calls \( u(x) \equiv \psi(x) \), but I can't distinguish a "small \( \psi \)" from the "capital \( \Psi \)" so easily in my handwriting. You'll find different authors use both of these notations…)

So

\[
\frac{\partial \Psi}{\partial t} = u(x) \frac{d\phi}{dt} \quad \text{Note full derivative on right hand side!}
\]

\[
\frac{\partial^2 \Psi}{\partial x^2} = \phi(t) \frac{d^2 u(x)}{dx^2}
\]

So Schrödinger equation reads (with \( \frac{d\phi}{dt} = \dot{\phi} \), and \( \frac{du}{dx} = u' \))

\[
i\hbar \left( u' \phi \right) = -\frac{\hbar^2}{2m} \left( \phi u'' \right) + V(u \cdot \phi)
\]

Now divide both sides by \( \Psi = u \cdot \phi \).

\[
i\hbar \left( \frac{\dot{\phi}(t)}{\phi(t)} \right) = -\frac{\hbar^2}{2m} \left( \frac{u''(x)}{u(x)} \right) + V(x)
\]

This is not possible unless both sides are **constants**.

Convince yourself; that is the key to the "method of separation of variables".

Let's name this constant "E".

[Note **units** of E are \( \frac{\hbar^2}{2m \cdot \text{(dist)}^2} \) or \( \frac{\hbar}{\text{time}} \) or simply \( V(x) \), either way, check, it's **Energy**!]

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting.
So (1)  \[ i\hbar \phi'(t) = E\phi(t) \]

(2)  \[ -\frac{\hbar^2}{2m}u''(x) + V(x)u(x) = E \cdot u(x) \]

These are ordinary O.D.E.'s

Equation (1) is about as simple as ODE's get!

Check  \[ \phi(t) = \phi_0 e^{-iEt/\hbar} \]

\(\hbar\) any constant, it's a linear ODE.

(1st order linear ODE is supposed to give one undetermined constant, right?)

This is "universal", no matter what \(V(x)\) is, once we find a \(u(x)\), we'll have a corresponding

\[ \Psi(x, t) = u(x)\phi(t) = u(x)e^{-iEt/\hbar} \]

But be careful, that \(u(x)\) depends on \(E\),

\[ -\frac{\hbar^2}{2m}u''(x) + V(x)u(x) = E \cdot u(x). \]

This is \[ \frac{d^2u(x)}{dx^2} \]
This is the "time independent Schrödinger equation".

You can also write this as  \[ \hat{H}[u(x)] = E \cdot u(x) \]
which is an "eigenvalue equation".

\(\hat{H} = "Hamiltonian"\) operator  \[ -\frac{\hbar^2}{2m} \frac{d^2}{dx^2} \]

\[ + V(x) \]

In general,  \[ \hat{H}[u(x)] = E \cdot u(x) \]
has many possible solutions.

\[ \psi_1(x), \psi_2(x), \ldots, \psi_n(x) \]
\[ \text{may all work, each corresponding to some } \text{particular } \text{eigenvalue} \]
\[ E_1, E_2, \ldots, E_n. \]
(What we will find is not any old \(E\) is possible if you want \(u(x)\) to be normalizable and well behaved. Only certain \(E\)'s, the \(E_n\)'s, will be ok.)

Such a \(u_n(x)\) is called a "stationary state of \(\hat{H}\)". Why? Let’s see…

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting.  Fall 2008
Notice that $\Psi_n(x,t)$ corresponding to $u_n$ is

$$\Psi_n(x,t) = u_n(x)e^{-iE_n t/\hbar} \quad \text{go back a page!}$$

So

$$\left|\Psi_n(x,t)\right|^2 = \left|u_n(x)\right|^2 \quad \text{no time dependence}$$

(for the probability density). It's not evolving in time; it's "stationary".

(Because $\left|e^{-iE_n t/\hbar}\right|^2 = 1$ ) Convince yourself!

(If you think back to de Broglie's free particle

$$\psi_{\text{free}} = Ae^{ikx}e^{-i\omega t} \quad \text{with } E = \hbar\omega,$$

it looks like we had stationary states $\Psi_n(x,t)$ with this (same) simple time dependence, $e^{i\omega t}$, with $\omega_n = E_n/\hbar$ This will turn out to be quite general)

If you compute $\langle \hat{Q} \rangle$ (the expectation value of any operator "Q" for a stationary state) the $e^{-iE_n t/\hbar}$ in $\Psi$ multiplies the $e^{iE_n t/\hbar}$ in $\Psi^*$, and goes away… (This is assuming the operator Q depends on x or p, but not time explicitly)

$$\langle \hat{Q} \rangle_{\text{in state } \Psi_n} = \int \Psi_n^*(x,t)\hat{Q}(x,\hbar i \frac{\partial}{\partial x})\Psi_n(x,t)dx$$

again, no time dependence.

Stationary state are dull, nothing about them (that's measurable) changes with time. (Hence, they are "stationary").

(Not all states are stationary … just these special states, the $\Psi_n$'s)

And remember, \[ \hat{H}u_n(x) = E_n u_n(x) \quad \text{Time Indep. Schröd. eq'n (2) } \]

and also \[ \hat{H}\Psi_n(x) = E_n \Psi_n(x,t) \]

This is still the Schrödinger Equation! (after plugging in our time solution for $\varphi(t)$.)
Check this out though: for a state $\Psi_n(x,t)$

$$\langle \hat{H} \rangle = \int \Psi_n^* \hat{H} \Psi_n \, dx = \int \Psi_n^* E_n \Psi_n \, dx$$

In stationary $\Psi_n$:

$$= E_n \int \Psi_n^* \Psi_n \, dx = E_n$$

constants come out of integrals  
this is normalization!

So the state $\Psi_n$ "has energy eigenvalue $E_n$" and has expectation value of energy $E_n$.

and

$$\langle \hat{H}^2 \rangle = \int \Psi_n^* \hat{H}^2 \Psi_n \, dx = \int \Psi_n^* \hat{H} ( \hat{H} \Psi_n ) \, dx$$

$$= \int \Psi_n^* \hat{H} ( E_n \Psi_n ) = E_n \int \Psi_n^* \hat{H} ( \Psi_n ) \, dx$$

$$= E_n^2$$

so

$$\sigma_H^2 = \langle H^2 \rangle - \langle H \rangle^2 = E_n^2 - E_n^2 = 0$$

Think about this – there is zero “uncertainty” in the measurement of $\hat{H}$ for this state.

Conclusion:

$\Psi_n$ is a state with a definite energy $E_n$. (no uncertainty!)

(that's really why we picked the letter $E$ for this eigenvalue!)

Remember, 

$$\hat{H} \Psi = i\hbar \frac{\partial \Psi}{\partial t}$$ is linear, so …

If $\Psi_1$ and $\Psi_2$ are solutions, so is $(a \Psi_1 + b \Psi_2)$.

But, this linear combo is not stationary! It does not have any "energy eigenvalue" associated with it!
The most general solutions of $\hat{H}\Psi_n = i\hbar \frac{\partial \Psi}{\partial t}$ (given $V= V(x)$) is thus

$$\Psi_{general}(x,t) = \sum_{n=1}^{\infty} c_n \Psi_n(x,t) = \sum_{n=1}^{\infty} c_n u_n(x)e^{-iE_n t / \hbar}$$

not an energy eigenstate
not stationary
It's a "mixed energy!"

energy eigenstate, or "stationary state"

any constants you like, real or complex,
Just so long as you ensure $\Psi$ is normalized.

Any/every physical state can be expressed in this way, as a combination of the more special stationary states, or "eigenstates".

If you measure energy on a stationary state, you get $E_n$, (definite value),

(But if you measure energy on a mixed or general state,…we need to discuss this further! Hang on …)
The Infinite Square Well: Let's make a choice for $V(x)$ which is solvable that has some (approximate) physical relevance.

Classically,

$$\bar{F} = -\frac{dV}{dx}$$

is 0 in the middle (free) but big at the edges. Like walls at the edges.

(Like an electron in a small length of wire:

```
  \hline
  |     | a     |
  \hline
  \hline
```

free to move inside, but stuck, with large force at the ends prevents it from leaving.)

Given $V(x)$, we want to find the special stationary states

(and then we can construct any physical state by some linear combination of those!)
Recall, we're looking for $u_n(x)$ such that

$$\hat{H}u_n(x) = E_n u_n(x)$$

(and then $\Psi_n(x,t) = u_n(x) e^{-iE_n t / \hbar}$)

(or, dropping "n" for a sec)

$$-\frac{\hbar^2}{2m} u''(x) + V(x) u(x) = E \cdot u(x)$$

(and then at last)

$$\Psi_{\text{general}}(x,t) = \sum_n c_n \Psi_n(x,t)$$

Inside the well, $0 < x < a$, and $V(x) = 0$, so in that region

$$\frac{d^2u}{dx^2} = -\frac{2mE}{\hbar^2} u(x) \equiv -k^2 u(x)$$

Here, I simply defined $k = \sqrt{\frac{2mE}{\hbar}}$

It's just shorthand, $k = \sqrt{\frac{2mE}{\hbar}}$ so $E = \frac{\hbar^2}{2m} k^2$

(However, I have used the fact that $E > 0$, you can't ever get $E < V_{\text{min}} = 0$! Convince yourself!)

I know this 2nd order ODE, and its general solution is

$$u(x) = A \sin kx + B \cos kx \quad \text{or} \quad \alpha e^{ikx} + \beta e^{-ikx}$$

But Postulate I says $u(x)$ should be continuous.

Now, outside $0 < x < a$, $V(x) \to \infty$. This is unphysical, the particle can't be there! So $u(x) = 0$ at $x = 0$ and $x = a$. This is a BOUNDARY CONDITION.

$u(x = 0) = A \cdot 0 + B \cdot 1 = 0$ so $B = 0$. (required!)

$u(x = a) = A \sin ka = 0$.

But now, I can't set $A = 0$ 'cause then $u(x) = 0$ and that's not normalized.

So $\sin ka = 0$, i.e. $k = n\pi / a$ with $n = 1, 2, 3, \ldots$

Ah ha! The boundary condition forced us to allow only certain $k$'s. Call them $k_n = \frac{n\pi}{a}$

Then since $E = \frac{\hbar^2}{2m} k^2$, we get $E_n = \frac{n^2\pi^2\hbar^2}{2ma^2}$

(Note: negative $n$'s just re-define "$x"$, it's not really a different state, $A \sin (kx)$ is the same function as $A \sin (-kx)$).

(Note: $n = 0$ no good, because $\sin (0x) = 0$ is not normalizable!)

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting. Fall 2008
Thus, our solutions, the "energy eigenstates" are

\[ u_n(x) = A \sin\left(\frac{n\pi x}{a}\right) \quad n = 1, 2, 3, \ldots \ (0 < x < a) \]

(and 0 elsewhere)

and \[ E_n = n^2 \left(\frac{\pi^2 \hbar^2}{2ma^2}\right) \]

For normalization, \[ \int_0^a u_n(x)^2 \, dx = 1 \Rightarrow |A|^2 \int_0^a \sin^2 \left(\frac{n\pi x}{a}\right) \, dx = 1 \]

Convince yourself, then, \(|A|^2 = \frac{2}{a}\) is required.

So \[ u_n(x) = \sqrt{\frac{2}{a}} \sin\left(\frac{n\pi x}{a}\right) \text{ (for } 0 < x < a) \]

(Note sign/phase* out front is not physically important. If you multiply \(u_n(x)\) by \(e^{i\theta}\), you have a wave function with the same \(|u(x)|^2\), it's physically indistinguishable. So e.g. \(u_n(x)\) and \(-u_n(x)\) are not "different eigenstates".)

- Energy is quantized (due to boundary condition on \(U\))
- Energies grow like \(n^2\)
- Lowest energy is not 0! (You cannot put an electron in a box and have it be at rest!)
Key properties to note (many of which will be true for most potentials, not just this one!)

Energy eigenstates \( u_n(x) \) are …

- "even" or "odd" with respect to center of box (\( n=1 \) is even, \( n=2 \) is odd, this alternates)
- oscillatory, and higher energy \( \leftrightarrow \) more nodes or zero crossings. (Here, \( u_n \) has \((n-1)\) intermediate zeros)
- orthogonal:

\[
\int_0^a u_n(x) u_m(x) dx = \delta_{nm} = \begin{cases} 0 & \text{if } n \neq m \\ 1 & \text{if } n = m \end{cases}
\]

check this, if you don't know why \( \sin \left( \frac{n\pi x}{a} \right) \) forms orthogonal states, work it out.

- complete: Dirichlet's Theorem, the basis of "Fourier series" says ANY function \( f(x) \) which is 0 at \( x = 0 \) and \( x = a \) can be written, uniquely,

\[
f(x) = \sum_{n=1}^{\infty} c_n u_n(x) \quad \text{can always find the } c_n's:
\]

(Recall, \( u_n = \sqrt{\frac{2}{a}} \sin \frac{n\pi x}{a} \) here)

Fourier's trick finds those \( c_n \)'s, given an \( f(x) \):

If \( f(x) = \sum_n c_n \psi_n(x) \) then do "the trick" …

Multiply both sides by "\psi*", \( f(x) \psi_m^*(x) = \sum_n c_n \psi_n(x) \psi_m^*(x) \)

Then integrate, so \( \int_0^a f(x) \psi_m^*(x) dx = \sum_n c_n \int_0^a \psi_n(x) \psi_m^*(x) dx \)

This tells us (flipping left and right sides)

\( c_m = \int_0^a f(x) \psi_m^*(x) dx \quad \text{this is how you figure out } c_m \)'s!

(The above features are quite general! Not just this problem!)

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And last but not least, this was all for \( u_n(x) \), but the full time-dependent stationary states are \( u_n(x)e^{-\frac{E_n t}{\hbar}} \), or writing it out:

\[
\psi_n(x,t) = \sqrt{\frac{2}{a}} \sin\left(\frac{n\pi x}{a}\right) e^{-i\frac{n^2\hbar^2}{2ma^2} \frac{t}{\hbar}}
\]

this particular functional form is specific to "particle in box"

This is still a stationary state, full solution to "Schröd in a box" with definite energy \( E_n \).

These are special, particular states, "eigenstates of \( \hat{H} \)."

The most general possible state of any "electron in a 1-D box", then, would be any linear combo of these:

\[
\psi(x,t) = \sum_n c_n \psi_n(x,t) = \sum_n c_n u_n(x)e^{-iE_n t h}\n
\]

You can pick any \( c_n \)'s you like (even complex!) and this will give you all possible physical states.

You might choose the \( c_n \)'s to give a particular \( \psi(x,t=0) \) that you start with, using Fourier's trick!

If \( \psi(x,0) = \sum_n c_n u_n \) is given, find the \( c_n \)'s,

and then the formula at top of page tells you \( \psi(x,t) \).

[So given initial conditions, we know the state at all times. That's the goal of physics!]

For other potentials, game is same: find \( u_n(x) \)'s and corresponding \( E_n \)'s, then form a linear combo at \( t=0 \) to match your starting state, and let it evolve.

\[
\psi(x,t) = \sum_n c_n u_n(x)e^{-iE_n t h} \ldots
\]
One last comment: If \( \psi(x,t) = \sum_n c_n u_n(x)e^{-iE_n t/\hbar} \)
then normalization says \( \int \psi^* \psi dx = 1 \)

But when you expand \( \Psi^* \) and \( \Psi \), all cross terms vanish after integration,
\[
\int u_n^*(x)u_m(x)dx = \delta_{nm}, \text{ leaving only the terms with } n=m, \text{ which integrate simply.}
\]

Work it out: \( \int \psi^* \psi dx = 1 \) tells you
\[
1 = \sum_n \sum_n \int c_n^* c_n u_n^*(x)u_n(x)e^{i(E_n - E_m)t/\hbar} dx
\]

Note: this collapses the whole integral to \( \delta_{nm} \), (thus \( e^{i(stuff)\cdots \rightarrow 1} \))
\[
= \sum_m \sum_n c_n^* c_n \delta_{nm}
\]
\[
= \sum_n |c_n|^2 \quad \rightarrow \text{ convince yourself.}
\]

Similarly (try it!) use \( \hat{H}\psi_n = E_n\psi_n \), to get
\[
\langle H \rangle = \int \psi^* \hat{H} \psi = \sum_n E_n |c_n|^2
\]

Interpretation (which we'll develop more formally)
\[
\psi(x,t) = \sum_n c_n u_n(x)e^{-iE_n t/\hbar} = \sum_n c_n \Psi_n(x,t)
\]
says "the state \( \Psi \) is a linear combination of the special stationary states, the \( \Psi_n \) 's."

The \( c_n \)'s carry information, \( \langle H \rangle = \sum_n |c_n|^2 E_n \) which looks like \( \sum_n P_n E_n \)

\[|c_n|^2 = \text{Probability that this particle's energy would be measured } E_n.\]

(so \( \sum_n |c_n|^2 = 1 \) is just "conservation of probability").

This is a central concept in QM, we'll come back to it often.
The Harmonic Oscillator:

Many objects experience a (classical) force \( F = -kx \), Hooke's Law. This is the harmonic oscillator, and \( V(x) = \frac{1}{2}kx^2 \).

Note that almost any \( V(x) \) with a minimum will look at least approximately like this, at least for small \( x \), so it's physically a very common situation!

Classically, \( x(t) = A \sin \omega t + B \cos \omega t \) for Harm. Osc. with \( \omega \equiv \sqrt{k/m} \) (so \( V = \frac{1}{2}m\omega^2x^2 \))

\[
\uparrow \text{ a defined quantity, but it tells you period } T = \frac{2\pi}{\omega}.
\]

So let's consider the Quantum Harmonic Oscillator. As before, separate \( x \) and \( t \), and look for solutions not 0.

\[
-\frac{\hbar^2}{2m}u''(x) + \frac{1}{2}m\omega^2x^2u(x) = E\ u(x) \quad \iff \text{these } u's \text{ will give us stationary states.}
\]

(then, if you find such a \( u \), \( \psi(x,t) = u(x)e^{-iEt/\hbar} \) gives the time dependence.)

Just like before, we will find that, if we insist \( u(x) \to 0 \) \( x \to \infty \), i.e. our "boundary cond."),
then we will find only certain \( E \)'s, called \( E_n \), will, and the corresponding \( u_n(x) \) will be unique. So just like in the box, we'll have \( \Psi_n(x,t) \)'s, each \( n \) corresponding to a stationary state with discrete energy.

This differential equation is a 2nd order ODE but that "\( x^2 \)" term makes it hard to solve. There are tricks. (*See aside, next page)

**Trick #1:** For large \( x \), \( x^2 \) dominates, so

\[
-\frac{\hbar^2}{2m}u''(x) = (E - \frac{1}{2}m\omega^2x^2)u \approx \frac{1}{2}m\omega^2x^2u
\]

Solution (check, it's not familiar)

\[
u(x) = Ae^{-\frac{m\omega^2}{2\hbar^2}x^2} + Be^{\frac{m\omega^2}{2\hbar^2}x^2} \quad \text{(Note: 2 undetermined constants for 2nd order ODE, nice!)}
\]

\[
\uparrow \text{ This } B \text{ term is very nasty as } x \to \infty, \text{ toss is as unphysical!}
\]

Next comes trick #1 part 2, which is to "factor out" this large \( x \) behavior:
Let's now assume (hope, wonder if?) \( u(x) = h(x)e^{-\frac{m\omega x^2}{2\hbar}} \), and hope that maybe \( h(x) \) will be "simple". Indeed, use the (love this name!) Method of Frobenius: try \( h(x) = a_0 + a_1x + a_2x^2 + \ldots \) Plug in, see what happens…

For now, I'm going to skip this algebraically tedious exercise and skip to the punchline: If you insist that \( u(x) \) stays normalizable, you find that series for \( h(x) \) needs to end, it must be a finite order polynomial. And that will happen if and only if \( E \) takes on special values, namely \( E_n = \left(n + \frac{1}{2}\right)\hbar\omega \) (where "n" is an integer)

If \( E \) is this, then the polynomial is "\( n \)th orther", and the \( h_n(x) \) functions that come out are Hermite Polynomials. (See Griffiths p. 56.)

\[
\begin{align*}
U_1(x) & \quad U_2(x) & \quad U_3(x) \\
\hline
\begin{array}{c}
\text{Graph of } U_1(x) \\
\text{Graph of } U_2(x) \\
\text{Graph of } U_3(x)
\end{array}
\end{align*}
\]

\[
u_n(x) = \left(\frac{m\omega}{\pi\hbar}\right)^{\frac{1}{4}} \frac{1}{\sqrt{2}} e^{-\frac{m\omega x^2}{2\hbar}} \quad \text{with } E_n = \frac{1}{2} \hbar\omega
\]

Note, here \( h_1(x)=1 \), basically, the stuff out front is just to normalize!

\[
u_2(x) = \left(\frac{m\omega}{\pi\hbar}\right)^{\frac{1}{4}} \frac{2}{\sqrt{8}} \sqrt{\frac{m\omega}{\hbar}} e^{-\frac{m\omega x^2}{2\hbar}} , \quad \text{with energy } E_2 = \frac{3}{2} \hbar\omega
\]

(and check out what's in front of the exponential besides all the constants - there's my 1st order polynomial, the first Hermite polynomial)

You could just verify if you want, just be TAKING the derivatives that, at least for these 1st 2 examples, \(- \frac{\hbar^2}{2m} u'' + \frac{1}{2} m\omega^2 x^2 u_n(x) = E_n u_n(x)\)

Remember, these give \( u_n(x) \), time dependence is simple:

Stationary states as a function of time look like \( \psi_n(x,t) = u_n(x)e^{-iE_nt/\hbar} \)

You can then form any combination of these to build any "particle in a well" state, and the time dependence follows: \( \Psi = \sum_n c_n u_n(x)e^{-iE_nt/\hbar} \).
Observations: Like particle in box, $u_n(x)$ are:

- even or odd
- oscillatory, with $u_n$ having $n - 1$ zero's
- orthogonal (really, orthonormal): $\int u_n(x)u_m(x)dx = \delta_{nm}$
- Complete: Any $f_n$ can be expanded via Fourier's trick as $\sum_{n=1}^{\infty} c_n u_n(x)$
  (if $f_n$ vanishes properly at $\infty$, anyway)
- Energy of each $u_n$ is discrete, grows like $n$ here.
- Lowest energy is not 0, it's $1/2 \hbar \omega$. (This means you cannot "stop" a mass on a quantum spring!)

For small $n$, $|u_n(x)|^2$ doesn't look classical at all,
but for large $n$, it sort of does if you "average over short wiggles"!
Here's a "real" picture from Shankar's text, fig 7.2, page 202:

This dashed curve is the classical $\rho(x)$
function. Note that classically $E$ can
be anything, but given $E$,
you "turn around" at $1/2 \hbar k x_{\text{max}}^2 = E$. 

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting.  Fall 2008
--- An Aside ---

Another method to know of, you might call it **Trick #2: Numerical solutions!**

How to solve \( u'' = -\frac{2m}{\hbar^2} (E - V(x))u(x) \) ? (without yet knowing \( E! \))

Even solutions: Guess an \( E \) (use a similar situation, or just units?)
For Harmonic Oscillator, \( \hbar \omega \) has units of energy, so maybe just try it as a starting guess, call it \( \tilde{E} \).

Let \( u(0) = 1, u'(0) = 0 \) \( \Rightarrow \) flat at origin, since I'm looking for the **even** solution!

\( u(0) \) may not be 1, but we'll normalize later!
Now, pick a "step-size" \( \varepsilon \) (tiny!) We'll "step along" in \( x \), finding \( u(x) \) point by point…

- Use S.E. to tell you \( u''(0) \). \[ u''(0) = -\frac{2m}{\hbar^2} (\tilde{E} - V(0))u(0) \]
- Use \( u'(0) = \frac{u(\varepsilon) - u(0)}{\varepsilon} \) to compute \( u(\varepsilon) \)
- Use \( u''(0) = \frac{u'(\varepsilon) - u'(0)}{\varepsilon} \) to compute \( u'(\varepsilon) \)

Now repeat, stepping across \( x \) in stops of \( \varepsilon \).

- If \( u \) starts to blow up, and start again.
  - you need **more** curvature at start, so raise \( \tilde{E} \)

- If \( u \) blows down
  - you need less curvature, **lower** \( \tilde{E} \)

For odd solutions, start with \( u(0) = 0, u'(0) = 1 \) (again, \( u'(0) \) may not be 1, but we'll renormalize later) Otherwise, it's the same game …

So basically you **pick** an \( E \), numerically "shoot" from the origin, and keep fixing \( E \) till you get a solution which behaves well at large \( x \).

In this way, you can **plot/compute** \( u(x) \) and find the energies, for essentially any \( V(x) \).
(In this sense, you don't have to feel bad that only a few \( V(x) \)'s yield analytic solutions for \( u_n(x) \) and \( E_n \). All \( V \)'s can be "solved" numerically.)

--- End Aside ---
Trick #3 for solving the Harmonic Oscillator.

This trick is way out there! I would never have come up with it (?) but it's cool, and turns out to be more general than you can imagine right now. It's the basis for a similar trick to understand angular momentum in 3-D, and then spin, and moving on to quantum field theory. So it's worth learning! It will also teach us some "operator methods" that are deeply central to QM. (So enjoy this "trickery". The math is fun!)

Here again is the 1-D Schrödinger equation we're studying:

\[
\frac{1}{2m} \left[ p^2 + (m\omega x)^2 \right] u = E u \quad \text{(with } p = \frac{\hbar}{i} \frac{\partial}{\partial x} \text{ of course!)}
\]

For numbers, \( c^2 + b^2 = (ic + b)(-ic + b) \), so this tempts us to rewrite the equation on left as \( \frac{1}{2m} \left( ip + m\omega x \right) \left( -ip + m\omega x \right) \)

(But this isn't right, because \( \hat{p} \) operates on any x's it hits...)

Still, inspired by my idea of rewriting the equation, let's define new operators

\[
\hat{a}_+ = \frac{1}{\sqrt{2\hbar\omega}} \left( \frac{\hat{p}}{i} + m\omega \hat{x} \right)
\]

\[
\hat{a}_- = \frac{1}{\sqrt{2\hbar\omega}} \left( -\frac{\hat{p}}{i} + m\omega \hat{x} \right)
\]

These are just definitions, and the names \( \hat{a}_+ \) and \( \hat{a}_- \) will make more sense in a sec. We introduced the constants out front just for units...

Now try writing what I had before (aside from some constant)

\[
\hat{a}_+ \hat{a}_- = \frac{1}{2\hbar\omega} \left( +i\hat{p} + m\omega \hat{x} \right) \left( -i\hat{p} + m\omega \hat{x} \right)
\]

\[
= \frac{1}{2\hbar\omega} \left( \left( \hat{p}^2 + (m\omega \hat{x})^2 \right) - im\omega \left( \hat{x} \hat{p} - \hat{p} \hat{x} \right) \right)
\]

The last terms looks like zero, since for numbers, \( x p - p x = 0 \), but not for operators!

We call \([\hat{A}, \hat{B}] = \hat{A}\hat{B} - \hat{B}\hat{A}\) the commutator of A and B.

My expression has \([\hat{x}, \hat{p}]\) in it. What's that do?
What does $[\hat{x}, \hat{p}]$ mean? It's an operator, so consider acting it on any old function $f(x)$:

$$[\hat{x}, \hat{p}] f(x) = (\hat{x}\hat{p} - \hat{p}\hat{x}) f = x - \frac{\hbar}{i} \frac{\partial}{\partial x} f - \frac{\hbar}{i} \frac{\partial}{\partial x} x f$$

but $\frac{\partial}{\partial x} (xf) = f + x \frac{\partial f}{\partial x}$. The term $x \frac{\partial f}{\partial x}$ explicitly cancels, leaving

(please check this for yourself) $[\hat{x}, \hat{p}] f = -\frac{\hbar}{i} f$.

Since $f(x)$ is totally arbitrary, this is true for ANY $f$, we say

$$[\hat{x}, \hat{p}] = -\frac{\hbar}{i}$$

This is very important, $[\hat{x}, \hat{p}] = i\hbar$, it's not zero! (we'll talk much more about this!)

OK, back to our problem: looking again at that product of $\hat{a}_-$ and $\hat{a}_+$, we have

$$\hat{a}_- \hat{a}_+ = \frac{1}{2\hbar \omega} (\hat{p}^2 + (m \omega \hat{x})^2) - \frac{im \omega}{2\hbar \omega} \cdot (i\hbar)$$

$$= \frac{1}{\hbar \omega} \hat{H} + \frac{1}{2}$$

or

$$\hat{H} = \hbar \omega (\hat{a}_- \hat{a}_+ - \frac{1}{2})$$

Now go through yourself and convince yourself that if instead we had written down that product in the opposite order, we would have gotten

$$\hat{H} = \hbar \omega (\hat{a}_+ \hat{a}_- + \frac{1}{2})$$

(note the other sign for the 1/2 term at the end)

And indeed, we can easily compute the commutator of $\hat{a}_-$ and $\hat{a}_+$ (which you could do directly by writing out $\hat{a}_-$ and $\hat{a}_+$ in detail, but it's easier to use the expressions just above):

$$[\hat{a}_-, \hat{a}_+] = \left( \frac{1}{\hbar \omega} \hat{H} + \frac{1}{2} \right) - \left( \frac{1}{\hbar \omega} \hat{H} - \frac{1}{2} \right)$$

$$= 1$$

OK, but why have we done all this?? Here comes the crux ( #1 ) . . .

Suppose $u(x)$ is a solution to $\hat{H} u = E u$

i.e. suppose we knew the $u$'s (which is what we're after, remember!)

Let's consider acting $\hat{H}$ on $\hat{a}_+$ which in turn is acting on $u$. (Why? You'll see!)

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting. Fall 2008
\[ \hat{H}(\hat{a}_+, u) = \hbar \omega (\hat{a}_+ \hat{a}_+ + \frac{1}{2}) \hat{a}_+ u \]

\[ = \hbar \omega (\hat{a}_+ \hat{a}_+ \hat{a}_+ + \frac{1}{2} \hat{a}_+) u \]

\[ = \hbar \omega \hat{a}_+ (\hat{a}_+ \hat{a}_+ + \frac{1}{2}) u \]

\[ = \hat{a}_+ \hbar \omega \left( \hat{a}_+ \hat{a}_+ + 1 + \frac{1}{2} \right) u \]

\[ \text{Constants slide past } \hat{a}_+ \text{, right? So } \text{a}_+ \text{-a}_- = 1 \]

\[ = \hat{a}_+ (\hbar \omega (\hat{a}_+ \hat{a}_+ + \frac{1}{2}) + \hbar \omega u \]

\[ \text{But, this is just } \hat{H} \text{ again!} \]

Look over all the steps, these tricks are not hard, but require some practice. This is "operator methods", we'll play this sort of game more often later in the course!

OK, one more step: now use \( \hat{H} \) \( u = Eu \)

Thus \( \hat{H}(\hat{a}_+, u) = \hat{a}_+ (E + \hbar \omega) u \)

or \( \hat{H}(\hat{a}_+, u) = (E + \hbar \omega) (\hat{a}_+ u) \)

So \( \hat{a}_+ u \) is an eigenfunction of \( \hat{H} \) with eigenvalue \( (E + \hbar \omega) \)

Given \( u \), I found a new, different eigenfunction with a different eigenvalue.

Now you try it, this time with \( \hat{H}(\hat{a}_- u) \). I claim you'll get \( (E - \hbar \omega) (\hat{a}_- u) \).

So now we see the use of \( \hat{a}_- \) or \( \hat{a}_+ \): they generate new solutions, different eigenvalues and eigenfunctions.

\( \hat{a}_+ = \) "raising operator" because it raises \( E \) by \( \hbar \omega \).

\( \hat{a}_- = \) "lowering operator" because it lowers \( E \) by \( \hbar \omega \).

They are "ladder operators" because \( E \) changes in steps of \( \hbar \omega \).
Now for crux #2 …

If you apply \( \hat{a}_- \hat{a}_- \ldots \hat{a}_- u \) you create a state with \( E - m \hbar \omega \).

You can always pile up enough \( a_-'s \) to get a negative energy... but there's a theorem that says \( E \) can never be lower than the minimum of \( V(x) \), which is 0.

ACK! This is bad ... unless there is a bottom state \( u_0 \), such that \( a_- u_0 = 0 \). (That's the ONLY way out! If there was no such "bottom state", then we could always find a physical state which is impossible - one with negative energy)

Note that 0 "solves" the S. Eq., it's just not interesting.
And from then on applying more \( a_- \)'s doesn't cause any new troubles: \( \hat{a}_- \hat{a}_- u_0 = \hat{a}_- 0 = 0 \ldots \)

OK, so \( \hat{a}_- u_0(x) = 0 \). Let's write this equation out in detail, because we know what \( \hat{a}_- \) is:

\[
\frac{1}{\sqrt{2m\hbar\omega}} \left( \hbar \frac{d}{dx} + m\omega x \right) u_0 = 0
\]

That's no problem to solve! See Griffiths, or just check for yourself:

\( u_0(x) = Ae^{-mx^2/2\hbar} \) works (for any \( x \)) And it's a 1st order ODE, so with one unedtermined coefficient (A) here, that's the most general solution.

We can even find A: pick it to normalize \( u_0(x) \) (use the handy relation \( \int_{-\infty}^{\infty} e^{-x^2} dx = \sqrt{\pi} \))

So \( u_0(x) = \left( \frac{m\omega}{\sqrt{\pi\hbar}} \right)^{\frac{1}{4}} e^{-mx^2/2\hbar} \).

This solves S.E., just plug it in! On the right, you'll see \( E_0 \) pop out, \( E_0 = \frac{1}{2} \hbar \omega \).

Indeed, there's a very cute operator trick to see this directly:

\[
\hat{H} u_0 = E_0 u_0 \Rightarrow \hbar \omega (\hat{a}_- + 1/2) u_0 = E_0 u_0
\]

so \( \frac{1}{2} \hbar \omega u_0 = E_0 u_0 \)
OK, we found the "bottom of the ladder". But now, \( \hat{\alpha}_+ u_0 \) gives another state, (call it \( u_1 \)), with \( E_1 = E_0 + \hbar \omega = \frac{3}{2} \hbar \omega \). (although, you may have to normalize again!)

And in general, climbing back up the ladder, \( u_n(x) = A_n (\hat{\alpha}_+)^n u_0(x) \) with \( E_n = (n + \frac{1}{2}) \hbar \omega \)
(Although you must normalize each one separately - see Griffiths or homework!)

Question: could we have missed some solutions in this way? Might there be a \( \tilde{u}(x) \) which is not "discovered" by repeated applications of \( \hat{\alpha}_+ \) to \( u_0 \)? Fortunately no.
If there were, \( (\hat{\alpha}_-)^n \tilde{u}(x) \) would still have to give 0 for some \( n \), and thus there'd be a "lowest" \( \tilde{u}(x) \) ... but we found the unique lowest state already!

Griffiths has a cute trick to normalize the states, see p. 47 - 48.
He also has an elegant trick to prove \( \int_{-\infty}^{\infty} \tilde{u}_n(x) u_n(x) dx = \delta_{mn} \) (see page 49)

But let's leave this for now:

Bottom line: By "Frobenius" or "ladder operators", we found all the stationary states \( u_n(x) \), with \( E_n = (n + \frac{1}{2}) \hbar \omega \).

(In general, potentials \( V(x) \) give "bound states", with discrete energies.)

But let's move on to a qualitatively different problem, where we're not bound!

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting. Fall 2008
**The Free Particle:**
Consider $V(x) = 0$. No PE, just a free particle. We talked about this at the start, (de Broglie's ideas). Now we can tackle it with the Schrödinger Equation. There is some funny business here, but clearly this too is an important physics case, we need to describe quantum systems that aren't bound!

$$\frac{-\hbar^2}{2m} u''(x) = E \cdot u(x)$$ That looks easy enough!

Define $k = \sqrt{\frac{2mE}{\hbar}}$, so $u''(x) = -k^2 u(x)$

I can solve that 2nd order ODE by inspection…

$$u_k(x) = Ae^{ikx} + Be^{-ikx}$$ (The label "k" identifies $u$, different k's $\rightarrow$ different u's)

There is no boundary condition, so $k$ (thus $E$) is not quantized. Free particles can have any energy!

And $\Psi_k(x,t) = (Ae^{ikx} + Be^{-ikx})e^{-iEt/\hbar}$

(Let's define $\omega = E / \hbar$, as usual) (with $E = \hbar^2 k^2 / 2m$)

Rewriting: $\Psi_k(x,t) = \frac{Ae^{ik(x-\frac{\omega t}{k})}}{\text{this is a fn of x-}vt \text{ right-moving wave!}} + \frac{Be^{-ik(x+\frac{\omega t}{k})}}{\text{this is a fn of x+}vt \text{ left moving wave!}}$

(The speed of a simple plane wave is also called "phase velocity". Go back to our chapter 1 notes on classical waves to see why $v=\omega/k$)

Here, $v = \frac{\omega}{k} = \frac{E}{\hbar} = \frac{\hbar^2 k^2}{2m} = \frac{\hbar k}{2m}$

Is that right? What did we expect? I'm thinking $p = \hbar k^*$ so $v = \frac{p}{m} = \frac{\hbar k}{m}$

* [ For the function $\Psi_k(x,t) = Ae^{ik(x-\frac{\hbar k^2}{2m}t)}$, you get $\hat{p}\Psi_k = \frac{\hbar}{i} \frac{\partial}{\partial x} \Psi_k = \hbar k \Psi_k$ So this $\Psi_k$ is an eigenfunction of $\hat{p}$ with eigenvalue $\hbar k$. Just exactly as our "de Broglie intuition" says. (The $\Psi_k$ at the top of the page, (with $k > 0$ only) is a mix of +k and -k, which means a mix of right moving and left moving plane waves….)]

But wait, What's with that funny factor of 2, then? Why did I get $v = \frac{\hbar k}{2m}$? We're going to need to consider this more carefully, coming soon!

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Summary

So we have solution \( \Psi_k(x,t) = Ae^{ikx-\omega t} \)

where \( k = \pm \sqrt{\frac{2mE}{\hbar}} \)  \( \text{(Note: 2 values of } k \text{ for each } E, \text{ one plus, one minus.)} \)

and \( \omega = \frac{E}{\hbar} = \frac{\hbar k^2}{2m} \) comes from the usual free particle energy (basically, \( \frac{p^2}{2m} \))

These waves have \( \lambda = \frac{2\pi}{|k|} \), because \( \Psi(x \pm \lambda) = \Psi(x) \)

The speed (phase \( \frac{\hbar |k|}{2m} = \frac{\hbar}{2m\lambda} \) velocity) (Small \( \lambda \) \( \Rightarrow \) faster speed)

There's that funny "2" we saw above, again. Classically, I expected \( v = \frac{p}{m} = \frac{\hbar k}{m} = \sqrt{\frac{2E}{m}} \) without the factor of 2.

Also, here's a wave function which is not renormalizable!

\[
\int_{-\infty}^{\infty} |\Psi_k(x,t)|^2 \ dx = A^2 \int_{-\infty}^{\infty} 1 \ dx \quad \text{yikes!}
\]

Conclusion from this: This \( \Psi_k \) is not representing a physical object!

There are no free (quantum) particles with definite energy!

But … not to worry, these \( \Psi_k \)'s are useful, even essential!
We can form linear combos that are physical.

So these “plane wave solutions” are eigenfunctions of momentum, they are eigenfunctions of the free-particle Hamiltonian (i.e. if \( V=0 \), they are easy to write down, they have a simple form. That’s all “good news”.
They are not normalizable, that’s “bad news”.
I would consider them to be idealizations of physical systems, a kind of limit (which cannot be actualized, but can be approximated) of an ideal free particle with definite momentum.
And as we will see, by the methods of Fourier transforms, they provide the mathematical basis for constructing perfectly physical free-particle states.
Since $k$ is a continuous variable, combining $\Psi_k$'s requires not a sum over $k$, but an integral over $k$!

Consider $\Psi_{\text{general}}^{\text{free}}(x,t) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} \phi(k) \Psi_k(x,t)\,dk = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} \phi(k)e^{ikx-E_kt}\,dk$

This looks a little scary, but it's just the continuous analogue of our old familiar "Fourier" formula:

$$\Psi_{\text{general}}^{\text{square well}}(x,t) = \sum_n c_n \Psi_n^{\text{square well}}(x,t),$$

where we basically replace $c_n \to \phi(k)\,dk$

- This $\Psi$ has many $k$'s (momenta, energies) superposed.

- If I give you $\Psi_{\text{general}}(x,0) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} \phi(k)e^{ikx}\,dk$ then you can figure out $\phi(k)$ (like Fourier's trick!) and given $\phi(k)$, you then know, from equation at top of page, $\Psi$ at all future times.

So we need to review "Fourier's trick" for continuous $k$!

Note $\Psi_k(x,0) = e^{ikx} \quad \leftarrow$ simple enough …

Suppose I give you $f(x)$, and ask "what $\phi(k)$ is needed" so that $f(x) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} \phi(k)e^{ikx}\,dx$

This is our problem: Given $\Psi_{\text{gen}}(x, t=0)$, find $\phi(k)$. Because once you've got it, you know $\Psi$ at all times.
The answer is Plancherel's Theorem, also known as the "Fourier Transform".

\[ \phi(k) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} f(x) e^{-ikx} \, dx \]  
(Note the minus sign in the exponential here!)

Recall that if \( f(x) = \sum_n c_n \sqrt{\frac{2}{a}} \sin \frac{n\pi x}{a}, \) then Fourier's trick gave \( c_n = \int_{-\infty}^{\infty} f(x) \sqrt{\frac{2}{a}} \sin \frac{n\pi x}{a} \, dx \)

It's rather similar; the coefficients are integrals of the desired function with our "orthogonal basis" functions.

- \( \phi(k) \) is the "Fourier transform" of \( f(x) \) here.
- If \( f(x) \) (here \( \Psi(x, t=0) \) is normalizable to start with, \( \Psi(x,t) \) will be too, as will \( \phi(k) \).
- You can "build" almost ANY FUNCTION \( f(x) \) you want this way!!

**Digression** – Some examples of Fourier transforms.

Griffiths (p. 62) has

\[ f(x) \]

\[ \begin{array}{c}
\text{f(x)} \\
\text{}\neg a \neg a \neg x
\end{array} \]

"width" is \( \approx 2a \).

then \( \phi(k) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} f(x) e^{-ikx} \, dx \) is easy.

you get \( \phi(k) = \frac{1}{\sqrt{\pi a}} \frac{\sin ka}{k} \)  
(Do it, try it yourself!)

If \( a \to 0 \),

Need many momenta to build a narrow wave packet

If \( a \to \infty \),

One (sharp) momentum \( \Rightarrow \) broad wave packet, where is it?

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Digression continued:

Let's do another: \( f(x) = Ae^{-x^2/4\alpha} \)

Here \( \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} f(x)e^{-ikx} \, dx \) can be done analytically (guess what's on the next homework?)

\[
\phi(k) = \frac{A}{\sqrt{2\pi}} \int_{-\infty}^{\infty} e^{\left(\frac{x^2}{4\alpha}+ikx\right)} \, dx = \frac{A}{\sqrt{2\pi}} e^{-k^2\alpha} \int_{-\infty}^{\infty} e^{-\left(\frac{x^2}{2\sqrt{\alpha}}-ik\sqrt{\alpha}\right)^2} \, dx
\]

(That little trick on the right is called "completing the square", you need to DO that algebra yourself to see exactly how it works, it's a common and useful trick)

Basically, letting \( x' = \frac{x}{2\sqrt{\alpha}} - ik\sqrt{\alpha} ; \quad dx' = \frac{dx}{2\sqrt{\alpha}} ; \quad \int e^{x^2} \, dx' = \sqrt{\pi} \)

you then get \( \phi(k) = \frac{2\sqrt{\alpha}}{\sqrt{2}} Ae^{-k^2\alpha} \)

Once again, narrow in \( f \rightarrow \) wide in \( \phi \), and vice versa.

- If \( f \) is centered around \( x_0 \), \( f(x) = Ae^{-(x-x_0)^2/4\alpha} \).

work it out… This will add (well, really multiply!) a phase \( e^{ikx_0} \) to \( \phi(k) \).

So the phase of \( \phi(k) \) does carry some important information (but not about momentum itself!)

(Puzzle for you: what does a phase \( e^{ikx} \) multiplying \( f(x) \) do? …)
So here's the central idea: instead of $\psi(x) = \text{Pure } e^{ikx}$, (which is not normalizable, and thus not a physical state) we will start with a more physical $\psi(x)$, a "wave packet".

Any $\psi(x)$ can be "built" like this, the math of Fourier transforms yields the following, in general:

- $\varphi(k)$ may very well turn out complex, (but might also be real.)
- If $\psi(x)$ is "mostly sinusoidal" with wavelength $\sim \lambda_0$ (like above) then $\varphi(k)$ will be centered around $k = 2\pi / \lambda_0$
- If $\psi(x)$ is localized with size $\Delta x$, then $\varphi(k)$ is localized with size $\Delta k \approx 1 / \Delta x$
- Given $\Psi(x, t=0) \Rightarrow \varphi(k) \Rightarrow \Psi(x, t)$ is determined for all times

Claim: If we start with a simple wave packet,

As time goes by, the ripples in the $\psi(x)$ wave packet will move with phase velocity $\omega/k$, but the envelope itself (which we interpret as "where the particle is located") moves with a different velocity, "group velocity" $v = \frac{d\omega}{dk}$. 

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting. Fall 2008
DIGRESSION on “phase and group velocity”, and the resolution of the “speed” puzzle.

Let’s look more at this last statement, about phase and group velocity. In our case (a free particle),

$$\omega = \frac{E}{\hbar} = \frac{\hbar k^2}{2m}, \quad \text{so} \quad \frac{d\omega}{dk} = \frac{\hbar k}{m}, \quad \text{no funny factor of 2 present!}$$

This resolves both our "free particle issues":

1) The packet moves with $v_{\text{classical}} = \frac{\hbar k}{m}$
2) The $\Psi(x, t)$ is perfectly normalizable.

- And something else happens: different $k$'s move with different speeds, so the ripples tend to spread out the packet as time goes by,

So $\Delta x$ grows with time!

Let me go back, and make a casual proof of the "claim" above about the time development of a wave packet:

$$\Psi_{\text{general}}(x, t) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} \phi(k)e^{i(kx - \omega t)} dk$$

Recall $\omega$ is itself dependent on $k$, $\omega = \frac{\hbar k^2}{2m}$ in this case of a free particle, it's just a consequence of $E = p^2/2m$.

Let's consider a wave packet where $\phi(k)$ is peaked near $k_0$.

So we have a "reasonably well-defined momentum, (this is what you think of when you have a real particle like an electron in a beam …) ( If you had many k's, each k travels at different speeds $\rightarrow$ packet spreads, not really a "particle" …)

So this is a special case, but a common/practical one …

So if k's are localized, consider Taylor expanding $\omega(k)$

$$\omega(k) = \omega(k_0) + \omega'(k_0)(k - k_0) + \ldots$$

This should be a fine approximation for the $\omega(k)$ in that integral we need to do at the top of the page!

Let's define $k' = k - k_0$ as a better integration variable, so $dk' = dk$, but

$$\omega \approx \omega(k_0) + \omega'(k_0)k' + \ldots$$
These are ripples, a function of (x - \( \omega_0 t \)). This is a function of (x - \( \omega_0 t \)).

They travel with \( \nu_{\text{phase}} = \frac{\omega_0}{k_0} \). That's the group velocity, \( \omega' = \frac{d\omega}{dk} \left|_{k_0} \right. \)

So the quantum free wave packet moves with \( \nu = \omega' \approx \frac{\hbar k_0}{m} \).

Just as you'd expect classically.

The ripples travel at \( \nu_{\text{phase}} \), but \( |\psi|^2 \) don't care about the ripples, really.

Wave packets do "spread out" in time, we could compute this (turns out it spreads faster if starts out narrower!)

END OF DIGRESSION. (Which showed how group and phase velocities are different, how wave packets DO behave classically, and how they represent physical particles)

At this point we're armed to solve a variety of general quantum mechanics problems: Given \( V(x) \), and \( \psi(t=0) \), find \( \Psi(x, t) \). That’s huge!

→ Bound states will generate discrete En's and un's (this is the story we started the term with – like the harmonic oscillator, or the infinite square well) The time dependence is especially interesting when you start with “mixed states”

→ Free particles will become "scattering states". This is what we’ve just gotten started. We'll talk more about this soon, at which point we’ll go back and tackle a few more simple problems, to get familiar with the \( \Psi \)'s for “scattering states”.

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting. Fall 2008
But first, let’s talk about the **INTERPRETATION** of our Fourier transform: what does \( \phi(k) \) **mean**? It plays an important role in Quantum Mechanics, and will lead us to even more powerful ways of thinking about wave functions. We’ve already danced around it - \( \phi(k) \) is the “momentum distribution”, it tells you about the particle in “momentum space”, it’s a **probability distribution for momentum**! Let’s see how this works:

To begin, consider again the general (math!) formulas for Fourier transform and inverse:

\[
\begin{align*}
\mathcal{F}(f(x)) &= \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} \phi(k) e^{ikx} \, dk \\
\mathcal{F}^{-1}(\phi(k)) &= \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} f(x) e^{-ikx} \, dx
\end{align*}
\]

Now, consider the Fourier transform of a \( \delta \) function, i.e. let \( f(x) = \delta(x) \), so (from above)

\[
\phi(k) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} \delta(x) e^{-ikx} \, dx = \frac{1}{\sqrt{2\pi}}
\]

(The last step is pure math - integrating \( \delta(x) \) gives you the rest of the integrand evaluated at the point where the argument of \( \delta(x) \) vanishes)

But now the inverse of this (just read it off from the Fourier formulas above!) says:

\[
f(x) = \delta(x) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} \phi(k) e^{ikx} \, dk = \frac{1}{2\pi} \int_{-\infty}^{\infty} e^{ikx} \, dk
\]

Let me “name” this relation (1)

It’s strange - the integral on the right looks awful. It oscillates forever, is it really defined? Well, it’s a little dicey (you do need to be careful how you apply it), but basically, if \( x \) is NOT 0 the “oscillating integrand” gives zero (think of real and imaginary parts separately, the area under a “sinusoidal function” is zero), but if \( x \) is zero, you integrate “1” over all space and get infinity. That’s the delta function!

OK, armed with this identity, consider *any* normalized wave function \( \Psi(x) \), for which

\[
\int_{-\infty}^{\infty} \Psi^*(x) \Psi(x) \, dx = 1.
\]

Let’s write \( \Psi(x) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} \phi(k) e^{ikx} \, dk \) (which you can ALWAYS do), so here \( \phi(k) \) is just the usual Fourier transform of \( \Psi(x) \).

We’ll put this formula for \( \Psi(x) \), into the normalization integral, (twice, once for \( \Psi(x) \) and once for \( \Psi^*(x) \).) Just be careful to change dummy indices for the second one!

\[
\int_{-\infty}^{\infty} \left( \int_{-\infty}^{\infty} \phi(k) e^{ikx} \, dk \right)^* \left( \int_{-\infty}^{\infty} \phi(k') e^{ik'x} \, dk' \right) \, dx = 1
\]

That’s a triple integral (ack!) but check it out, the only x-dependence is in TWO places, the two exponentials. So let’s pull that x dependence aside and do the x-integration first:

\[
\int_{-\infty}^{\infty} e^{-ikx} e^{ik'x} \, dx \left( \int_{-\infty}^{\infty} \phi(k) \, dk \right)^* \left( \int_{-\infty}^{\infty} \phi(k') \, dk' \right) = 1
\]

(Note the “–” sign in the ikx term came from the complex conjugation. Follow the algebra here, don’t take my word for anything!)
But I can DO that integral over x: just use our cute integral (1) to simplify. Let me first take integral (1) and rewrite it, flipping x’s and k’s:

\[ \delta(k) = \frac{1}{2\pi} \int_{-\infty}^{\infty} e^{ikx} \, dx \quad (1') \]  

(Convince yourself, treat x and k as pure symbols!)

But this is precisely what I have in my triple integral, except instead of k I have (k’-k). Substitute this in (getting rid of the dx integral, which “eats up” the Sqrt\[2\pi\] terms too!)

\[ \iint \delta(k' - k) \phi^* (k) \phi(k') \, dk \, dk' = 1 \]  

(Convince yourself, treat x and k as pure symbols!)

It’s easy to “glaze over” this kind of math. Pull out a piece of paper and DO it, the steps are not hard, and you’ll use these tricks many times. Check that you see what happened to the 2\pi’s and the integral over x)

There’s still a double integral over k and k’, but with the \delta(k’-k) it’s easy enough to do, I love integrating over \delta’s! Let’s integrate over k’, and collapse it with the \delta, giving

\[ \int_{-\infty}^{\infty} \phi^* (k) \phi(k) \, dk = 1 \]

Ah – this is important. We started with a normalized wave function \Psi(x), and what I just showed is that the Fourier transform, \phi(k) , is ALSO normalized. But there’s more! Consider now

\[ \langle p \rangle = \int_{-\infty}^{\infty} \Psi^* (x) \frac{\hbar}{i} \frac{\partial}{\partial x} \Psi(x) \, dx \]  

Play the exact same game as before, but with that derivative (d/dx) inside. All it will do will pull out an ik’ (convince yourself), everything else looks like what I just did, and we get

\[ \langle p \rangle = \int_{-\infty}^{\infty} (\hbar k) \phi^* (k) \phi(k) \, dk \]

Now, deBroglie identifies \hbar k with p. So just stare at this until you realize what it’s saying. I would expect (from week 1!) that \[ \langle p \rangle = \int_{-\infty}^{\infty} p \quad \text{Prob density}(p) \, dp \]

So here it is, apparently (other than annoying factors of \hbar, since p and k are essentially the same thing, p=\hbar k) we know what the “Probability density for momentum” is, it’s just \[ \phi^* (k) \phi(k) = |\phi(k)|^2 \]

To get the \hbar’s right, I would define \Phi(p) = \frac{1}{\sqrt{\hbar}} \phi(k), again with p=\hbar k. I’ll let you convince yourself that \Phi(p) is properly normalized too,

\[ 1 = \int_{-\infty}^{\infty} \Phi^* (p) \Phi(p) \, dp \quad \text{and} \quad \langle p \rangle = \int_{-\infty}^{\infty} p \Phi^* (p) \Phi(p) \, dp \]

We’ve just done a lot of math, so let’s pull this all together and summarize
Bottom line: We’d gotten used to thinking about $\Psi(x)$ as “the wave function”. When you square it, you learn the probability density to find a particle at position $x$. Expectation values tell us about OTHER things (any other operator/measurable you want.) $\Psi(x)$ contains ALL information about a quantum particle. Now we have the Fourier transform, $\varphi(k)$, or equivalently $\Phi(p) = \frac{1}{\sqrt{\hbar}} \varphi(k)$. This function depends on “$p$”.

It contains the SAME information as $\Psi(x)$!

It just lives in “momentum space” instead of “position space”. It’s a function of $p$.

• $\Phi(p)$ is normalized, just like $\Psi(x)$ was.
• $|\Phi(p)|^2$ tells the “probability density of finding the particle with momentum $p$”
• $\Phi(p)$ is the wave function too. There’s nothing special about $x$! $\Phi(p)$ is called the “momentum space wave function”.
• We can compute expectation values in momentum space to learn about anything, just like we could before. Let’s now consider this in more detail:

We already know $\langle p \rangle = \int_{-\infty}^{\infty} p \Phi^*(p) \Phi(p) dp$

In fact, it shouldn’t be hard to convince yourself that for any function of momentum $f(p)$,

$\langle f(p) \rangle = \int_{-\infty}^{\infty} f(p) \Phi^*(p) \Phi(p) dp$. 

How about $<x>$? Here again, start with the usual old formula in terms of $\Psi(x)$ and play the same math game as on the previous pages: expand $\Psi(x)$ as a Fourier transform, you get a triple integral, do the “$\delta$-function” trick, and (try this, it’s not so hard!) you find

$\langle x \rangle = \int_{-\infty}^{\infty} \Phi^*(p) \hbar i \frac{\partial}{\partial p} \Phi(p) dp$.

Look at that for a second. There’s a lovely symmetry here:

In position space, we’ve used $\hat{x} = x$ and $\hat{p} = -\hbar i \frac{\partial}{\partial x}$. Now we’re seeing that in momentum space, $\hat{x} = \hbar i \frac{\partial}{\partial p}$, and $\hat{p} = p$.

It’s almost perfect “symmetry” (except for a minus sign, but don’t forget, the inverse Fourier transform has an extra minus sign in front of the “$i\hbar$”, which is the source)

We will discover there are other “spaces” we might live in. Energy-space comes to mind, but pretty much any-operator-you-like-space can be useful too! The momentum-space example will be the first of many!

But for now we move back to our central problem, solving the Schrodinger equation. We’ve done two “bound states”, and the free particle. Next we will consider particles with positive energy (so, they’re still “free”) but where $V$ is not just simply 0 everywhere. This is still a lot LIKE the free particle, it’s a “free particle interacting with things”. This is getting to be real physics, like the beam of protons at LHC interacting with nuclei! We’ll still stick to one dimension (your beam can only go forward and backward, you can only transmit or reflect – no “scattering off at 27 degrees” just yet!) And, at the same time, we will tackle other “bound state” problems too - to keep the math simple, we’ll consider examples with “piecewise constant” $V(x)$, like in the picture on the next page:

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting. Fall 2008
Back to our problem, in general, of solving the TISE, Schrodinger's equation:

\[ \hat{H} u_n(x) = E_n u_n(x), \]

or writing it out:

\[ -\frac{\hbar^2}{2m} \frac{d^2}{dx^2} u(x) + V(x) u(x) = E u(x) \]

\[ u''(x) = -\frac{2m}{\hbar^2} (E - V(x)) u(x). \]

\[ \leftarrow \text{So } u \text{ and } u' \text{ need to be continuous, at least if } V \text{ is finite.} \]

If \( E > V(x) \), that means \( u'' = -(k^2) u \)

Classically, \( KE > 0 \)

\[ k^2 = \frac{2m(E - V)}{\hbar^2}, \quad \text{or } E - V = \frac{\hbar^2 k^2}{2m} = \frac{p^2}{2m} \text{ as usual!} \]

\[ u'' = -k^2 u \implies \text{curvature is towards axis. This is a wiggly function!} \]

If \( k^2 \) is constant, \( u = A \sin kx + B \cos kx \)

or \( = A'e^{ikx} + B'e^{-ikx} \)

Big K.E. \( \rightarrow \) Big \( k^2 \) \( \rightarrow \) more wiggly (smaller \( \lambda \))

Classically Big K.E. \( \rightarrow \) faster speed \( \rightarrow \) less likely to be found there.

This is very different, it corresponds in QM not to \( \lambda \), but amplitude. \( \leftarrow \) (well, \textit{usually}, but there are exceptions)

So as a specific example:

\( \text{I expect } \)

\[ \text{smaller } k \rightarrow \text{longer } \lambda \]

\[ \text{higher prob } \rightarrow \text{big } |\psi| \]

\[ \text{Big } k \rightarrow \text{small } \lambda \]

\[ \text{low prob } \rightarrow \text{small } |\psi| \]

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting.
What if $E < V$? Classically this never happens (negative K.E.?!)

In Q.M.,

$$u''(x) = -\frac{2m}{\hbar^2} (E - V(x)) \cdot u(x)$$

If $E < V$, define $\kappa = +\sqrt{\frac{2m(V-E)}{\hbar}}$, so $u'' = +\kappa^2 u$

This is not wiggly, curvature is away from axis.

If $\kappa^2 = $ constant (constant $V$ in a region)

$$u(x) = C \ e^{-\kappa x} + D \ e^{+\kappa x}$$

The other signs, e.g. $e^{+\kappa x}$ for $x > 0$ blow up at $\infty$, no good!

So in the "classically forbidden" regions, can have $u(x) \neq 0$.

But, $u \to 0$ exponentially fast. So e.g.

$$d = \text{"Penetration depth"} = \frac{1}{\kappa} = \frac{\hbar}{\sqrt{2m(V-E)}}$$

If $V \gg E$, rapid decay!
With just these ideas, you can guess/sketch \( u(x) \) in many cases:

- \( u, u' \) continuous always (if \( V, E \) finite)
- \( E > V \Rightarrow \) sinusoidal. Bigger \( E - V \Rightarrow \) smaller \( \lambda \) and usually smaller \( |\psi| \)
- \( E < V \Rightarrow \) decaying exponential. Bigger \( V - E \Rightarrow \) faster decay

Example: Finite square well \( \big\{ \) continuous spectrum (free)

\[
V(x) \bigg\} \text{ Discrete spectrum (bound)}
\]

\[
\begin{align*}
E_1 & \quad \text{Ground state, even about center} \\
E_2 & \quad \text{1st excited state, 1 node, odd}
\end{align*}
\]

Turns out, you can prove, if \( V(x) \) is even, \( V(x) = V(-x) \), then \( u(x) \) must be even or odd, so \( |u(x)|^2 \) is symmetric!

Let's work this out more rigorously to check, ok?

Let \( V(x) = \begin{cases} 
0 & \quad -a < x < a \\
+V_0 & \quad \text{outside} 
\end{cases} \)

Note: I shifted the zero w.r.t Griffiths

Assume \( 0 < E < V_0 \), i.e. bound state
Let's start with even solutions for \( u(x) \):

\[
V(x) = \begin{cases} \text{Region I} & & \text{Region II} & & \text{Region III} \\
\text{No good, blows up} & & \text{No good, looking for even solution right now} & & \text{Same } C, \text{ 'cause looking for even solution!}
\end{cases}
\]

Boundary conditions:

Continuity of \( u \) at \( x = +a \):
\[
\left. u_{III}(x) \right|_{x = +a} = \left. u_{II}(x) \right|_{x = a}
\]
Continuity of \( u' \) at \( x = +a \):
\[
\left. u'_{III}(x) \right|_{x = +a} = \left. u'_{II}(x) \right|_{x = a}
\]

(Conditions at \( x = -a \) add no new info, since we've already used fact that \( u \) is even!)

We have 3 unknowns: \( E \) and \( (A \text{ and } C) \)
The energy eigenvalue.

We have 2 boundary conditions and normalization, so we can solve for 3 unknowns!

Continuity of \( u \) says
\[
Ce^{-\kappa a} = A \cos k a
\]
Continuity of \( u' \) says
\[
-\kappa Ce^{-\kappa a} = -Ak \sin k a
\]

Dividing these \( \Rightarrow \)
\[
\frac{\kappa}{k} = \frac{\sqrt{V_0 - E}}{E}
\]
(A nasty, transcendental equation with one variable, \( E \))
Can solve that transcendental equation numerically, or graphically,

Let \( z = ka = \sqrt{\frac{2mE}{\hbar^2}a} \) \( \left\{ \right. \) so \( \frac{V_0}{E} = \left( \frac{z_0}{z} \right)^2 \)

Let \( z_0 = \sqrt{\frac{2mV_0}{\hbar^2}a} \) \( \left. \right\} \)

So we want \( \tan z = \sqrt{\left( \frac{z_0}{z} \right)^2 - 1} \)

I see there are solutions for \( z \) (which in turn means \( E \), remember, \( z \) and \( E \) are closely related, by definition, if you know one you know the other).

They are discrete. There is a finite number of them.

No matter what \( z_0 \) is, there's always at least solution. (Look at the graph!)

If \( V_0 \to \infty \), solutions are near \( \pi/2, 3\pi/2, \ldots \)

Giving \( E = \frac{\hbar^2}{2ma^2}Z^2 = \frac{\hbar^2 \pi^2}{2m(2a)^2} \) \( (1^2, \text{ or } 3^2, \ldots) \)

which is exactly the even solutions for our \( \infty \) well with width \( 2a \).

Knowing \( z \Rightarrow E \Rightarrow k \) and \( \mathcal{K} \), our B.C.'s give \( A \) in terms of \( C \), and normalization fixes \( C \).

I leave the odd solutions to you. The story is similar and you will get, as your transcendental equation, \( \cot(ka) = -\frac{\mathcal{K}}{k} \)

This time, it's not always the case there's even a single solution.

(The well needs to be deep enough to get a 1st excited state, but even a shallow well has one bound state.)

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting.

Fall 2008
What about $E > V_0$? This is the continuous region, also called the "scattering states", because $u$ will be nonzero off to $\infty$, it's like a particle comes in, interacts with our potential well, and is affected (scattered).

Lots of physics in this example (- like "shooting electrons" at some object it interacts with.)

Griffiths treats the square well, I'll do a simpler example:

We could build a wave packet, send it in from left with $E > V_0$, and solve for time dependence. That's the right way to make this physically realistic. But our approach will be simpler. Look for $u(x)$ which solves S.E. and which is physically interpreted as "something coming in from the left".

$$u_I(x) = Ae^{ikx} + Be^{-ikx} \quad \text{with} \quad k = \sqrt{\frac{2mE}{\hbar^2}}$$

$$u_{II}(x) = Ce^{ik'x} + De^{-ik'x} \quad \text{with} \quad k' = \sqrt{\frac{2m(E - V_0)}{\hbar^2}}$$

Remember our interpretation of free particle solutions from awhile back, the $Ae^{ikx}$ term \(\Rightarrow\) incoming wave, traveling rightwards with $p = \hbar k$

$Be^{-ikx}$ represents a left moving wave, but it's in region I, so it must physically represent reflected wave (in steady state, remember!)

$Ce^{ik'x}$ represents "transmitted wave", going right in region II.

We set $D = 0$, because otherwise we'd have "incoming wave from the right", and our physics condition was "waves coming from left".

Boundary Conditions: Continuity of $u$ and $u'$ at the origin gives:

$$u_I(x) \bigg|_{x=0} = u_{II}(x) \bigg|_{x=0}$$

$$u'_I(x) \bigg|_{x=0} = u'_{II}(x) \bigg|_{x=0}$$

(S. Pollock, taken from M. Dubson) with thanks to J. Anderson for typesetting.
What about normalization? These are not normalizable u's, (like the free particle, you must make a "wave packet" if you want it to be normalized.)

But wait, what are we after? Well, given A (= amplitude entering) I want to know, T, "Transmission Coefficient" and R, "Reflection Coefficient", how much of an incident wave moves on, and how much "bounces"?

**Intuitively,** B/A will tell us about R
and C/A will tell us about T. Really, R, and T are "the physics" we're after.

We need to be careful, I said "tell us about", but it's not "equal", we'll have to come back to R and T in a minute…Let's first compute these ratios, though.

Boundary condition on u(x) at x=0 gives A + B = C  (Because $e^{\pm i0} = 1$)

B.C. on u' gives $k (A - B) = k' C$

Now divide by k, add the two equations,

So I get (check!) $2A = C(1 + k'/k)$ or $\frac{C}{A} = \frac{2}{1 + k'/k}$

Put this back into 1st equation: $A + B = C = \frac{2}{1 + k'/k}A$, which then means (again, check that you follow this algebra, it's not hard, and worth doing)

$$\frac{B}{A} = \frac{2}{1 + k'/k} - 1 = \frac{1 - k'/k}{1 + k'/k}$$

with $\frac{k'}{k} = \sqrt{\frac{E - V_0}{E}}$

OK, so we know what B/A and C/A are. Now we can talk about interpreting these coefficients of (non-normalizable) "plane waves".
**Probability Current** $J(x,t)$.

$|\psi|^2$ evolves with time, smoothly. So probability "flows" in (and out) of this region!

I claim $\frac{\partial}{\partial t} |\psi(x,t)|^2 = - \frac{\partial}{\partial x} (J(x,t))$, and this $J$ function is called "Probability current"*

where $J(x,t) = \frac{i\hbar}{2m} \left( \psi \frac{\partial \psi^*}{\partial x} - \psi^* \frac{\partial \psi}{\partial x} \right)$

You need to do the algebra in prev. equation!

Let's rewrite the first "claimed" equation as $\frac{\partial P}{\partial t} + \frac{\partial J}{\partial x} = 0$ where $P = \text{Prob. density} = |\psi|^2$

or, integrating $a \rightarrow b$

$$\frac{\partial}{\partial t} \int_a^b P \cdot dx = -\int_a^b \frac{\partial J}{\partial x} \, dx = - (J(b) - J(a)) = J(a) - J(b)$$

rate of "buildup" = "current in at a" - "current out at b" in $(a,b)$ region

* Just like electric charge, the name "current" is good!

Note that $J(x,t) = \frac{i\hbar}{2m} (-2i) \cdot \text{Im} \left( \psi^* \frac{\partial \psi}{\partial x} \right) = \frac{\hbar}{m} \text{Im} \left( \psi^* \frac{\partial \psi}{\partial x} \right)$

For a plane wave, $\Psi = Ae^{i(kx-\omega t)}$

$\frac{\partial \Psi}{\partial x} = ik\Psi$, so $J = \frac{\hbar}{m} \text{Im} (ik\Psi^*\Psi)$, which means (again, for a plane wave),

$J = \frac{\hbar}{m} |A|^2 k$

This is key: $\frac{\hbar k}{m}$ looks like $\frac{p}{m} = v$

So "flow of probability" in a plane wave is like $v \cdot |A|^2$

Think of electric current, where $j = \rho \cdot v$ speed of charge

charge density

large amp. $\Rightarrow$ lots of flow

fast movement $\Rightarrow$ lots of flow

Very similar! $|A|^2$ plays the role of "charge density" here.
So back to our scattering problem, with particles coming in from the left and hitting a "step up" potential, but with $E>V$:

\[ J = \text{"probability current"}, \]

If $J = \text{"probability current"}$, then

\[
R = \text{reflection coefficient} = \frac{J_{\text{ref}}}{J_{\text{inc}}}, \quad \text{which is just } |B|^2/|A|^2
\]

Remember that from the formula for $J$ we got on the previous page?

\[
T = \text{Transmission coefficient} = \frac{J_{\text{trans}}}{J_{\text{inc}}}, \quad \text{which is NOT } |C|^2/|A|^2
\]

(Think a little about the presence of the $k'/k$ term there - I might not have expected it at first, thinking maybe $T = |C|^2/|A|^2$, but when you look at our probability current formula, you realize it makes sense, the waves have a different speed over there on the right side, and probability current depends not JUST on amplitude, but ALSO on speed!)

We solved for $B/A$ and $C/A$, so here

\[
R = \left( \frac{1 - \frac{k'}{k}}{1 + \frac{k'}{k}} \right)^2
\]

Convince yourself then, that $R + T = 1$ (!)  Convince yourself

\[
T = \frac{4}{(1 + \frac{k'}{k})^2} \cdot \frac{k'}{k}
\]

Note: If $V_0 \to 0$, \( \frac{k'}{k} = \frac{E - V_0}{E} \to 1 \) (Convince yourself!) and then check, from the formulas above, that $R \to 0$ and $T \to 1$. This makes sense to me - there's no barrier in this limit!

On the other hand, if $E \to V_0$, I claim $\frac{k'}{k} \to 0$ (again, do you see why?), and in THIS limit, the formulas above tell us that $R \to 1$ and $T \to 0$.

Again, this makes sense - in this limit, it is "forbidden" to go past the barrier. (Classically, if $E > V_0$, $R = 0$. QM gives something new here.)
Other examples:

\[ \begin{align*}
V & \quad V_0 \\
E & \quad \text{Potential barrier}
\end{align*} \]

Pain, but can compute R, T.

\( E < V_0 \): T is \textbf{not} 0, even though classically it would be. This is \textit{tunneling}.

\( E > V_0 \): Classically, T=1, but QM \(\Rightarrow\) \textit{some} reflection.