Quantum Mechanics

Numerical solutions of the Schrodinger equation

- Integration of 1D and 3D-radial equations
- Variational calculations for 2D and 3D equations
- Solution using matrix diagonalization methods
- Time dependence

Brief review of quantum mechanics

In classical mechanics, a point-particle is described by its position x(t) and velocity v(t)

• Newton's equations of motion evolve x,v as functions of time

In quantum mechanics, x and v cannot be precisely known simultaneously (the uncertainty principle). A particle is described by a wave function $\Psi(x,t)$

• the probability of the particle being in a volume dx is

 $P(x,t)dx \propto |\Psi(x,t)|^2 dx$

- The Schrödinger equation evolves $\Psi(x,t)$ in time
- There are energy eigenstates of the Schrodinger equation
 - for these, only a phase changes with time

 $\Psi_n(x,t) = \Psi_n(x,0)e^{-itE_n/\hbar}, \quad \hbar \approx 1.05 \cdot 10^{-34} Js$

⇒ Finding the energy eigenstates (stationary states) is an important task

Time dependent Scrodinger equation

$$i\hbar \frac{\partial}{\partial t} \Psi(x,t) = H \Psi(x,t) \rightarrow \text{stationary} \qquad H \Psi(x) = E \Psi(x)$$

Stationary Scrodinger equation in three dimensions

$$-\frac{\hbar}{2m}\nabla^2\Psi(\vec{x}) + V(\vec{x})\Psi(\vec{x}) = E\Psi(\vec{x})$$

Spherical symmetric potentials; separable

$$\Psi_{L,L_{z},n}(\vec{x}) = R_{L,n}(r)Y_{L,L_{z}}(\phi,\Theta) = \frac{1}{r}U_{L,n}Y_{L,L_{z}}(\phi,\Theta)$$

Radial wave function

$$\left(-\frac{\hbar^2}{2m}\frac{d^2}{dr^2} + \frac{L(L+1)h^2}{2mr^2} + V(r)\right)U_{L,n}(r) = E_{L,n}U_{L,n}(r)$$

Similar to purely one-dimensional problems

$$-\frac{\hbar^2}{2m}\frac{d^2\Psi(x)}{dx^2} + V(x)\Psi(x) = E\Psi(x)$$

Numerov's method (one dimension)

Stationary Schrodinger equation

$$-\frac{\hbar^2}{2m}\frac{d^2\Psi(x)}{dx^2} + V(x)\Psi(x) = E\Psi(x)$$

Can be written as (also radial function in three dimensions)

 $\Psi''(x) = f(x)\Psi(x)$

Discretization of space: Δ_x . Consider Taylor expansion

$$\Psi(\Delta_x) = \Psi(0) + \sum_{n=1}^{\infty} \frac{\Delta_x^n}{n!} \Psi^{(n)}(0)$$

Add expansions for $\pm \Delta_x$

$$\Psi(\Delta_x) + \Psi(-\Delta_x) = 2\Psi(0) + \Delta_x^2 \Psi''(0) + \frac{1}{12}\Delta_x^4 \Psi^{(4)}(0) + O(\Delta_x^6)$$

Second derivative determined by the Schrodinger equation How to deal with the fourth derivative? Central difference operator

$$\delta g(0) = g(\Delta_x/2) - g(-\Delta_x/2)$$

$$\delta^2 g(0) = \delta[\delta g(0)] = g(\Delta_x) - 2g(0) + g(-\Delta_x)$$

$$g''(x) = \frac{1}{\Delta_x^2} \delta^2 g(x) + O(\Delta_x^2)$$

We can rewrite the previous equation

$$\Psi(\Delta_x) + \Psi(-\Delta_x) = 2\Psi(0) + \Delta_x^2 \Psi''(0) + \frac{1}{12}\Delta_x^4 \Psi^{(4)}(0) + O(\Delta_x^6)$$

1

using the second central difference, giving

$$\delta^2 \Psi(0) = \Delta_x^2 \Psi''(0) + \frac{1}{12} \Delta_x^4 \Psi^{(4)}(0) + O(\Delta_x^6)$$

Approximate the fourth derivative

$$\Delta_x^4 \Psi^{(4)}(0) = \Delta_x^4 [\Psi''(0)]'' = \Delta_x^2 \delta^2 \Psi''(0) + O(\Delta_x^6)$$

leads to the general result

$$\delta^2 \Psi(0) = \Delta_x^2 \Psi''(0) + \frac{1}{12} \Delta_x^2 \delta^2 \Psi''(0) + O(\Delta_x^6)$$

 $\delta^2 \Psi(0) = \Delta_x^2 \Psi''(0) + \frac{1}{12} \Delta_x^2 \delta^2 \Psi''(0) + O(\Delta_x^6)$ Schrodinger equation $\Psi''(x) = f(x)\Psi(x)$ gives $\delta^2 \Psi(0) = \Delta_x^2 f(0) \Psi(0) + \frac{1}{12} \Delta_x^2 \delta^2 [f(0) \Psi(0)] + O(\Delta_x^6)$ More compact notation: $g_n = g(n\Delta_x)$ $\Psi_1 - 2\Psi_0 + \Psi_{-1} = \Delta_x^2 f_0 \Psi_0 + \frac{1}{12} \Delta_x^2 [f_1 \Psi_1 + f_{-1} \Psi_{-1} - 2f_0 \Psi_0] + O(\Delta_x^6)$ Introduce function $\phi_n = \Psi_n (1 - \Delta_r^2 f_n / 12)$

 $\phi_1 = 2\phi_0 - \phi_{-1} + \Delta_x^2 f_0 \Psi_0 + O(\Delta_x^6)$

Julia implementation

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for n=2:nx
    phi2=dx2*fn1*psi(n-1)+2*phi1-phi0
    phi0=phi1; phi1=phi2
    fn1=2*(potential(dx*n)-energy)
    psi(n)=phi1/(1-dx2*fn1)
    end
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Boundary-value problems

The Schrodinger equation has to satisfy boundary conditions ➤ **quantization**, as not all energies lead to valid solutions

Example: Particle in a box (infinite potential barrier)

 $V(x) = 0 \ (|x| < 1), \quad V(x) = \infty \ (|x| \ge 1)$

Using $\hbar = 1, m = 1$

 $\Psi''(x) = 2[V(x) - E]\Psi(x)$

Boundary conditions: $\Psi(\pm 1) = 0$

 $\Psi(x) = N \cos(n\pi x/2), \quad (n \text{ odd}) \qquad E = \frac{\pi^2 n^2}{8}$ $\Psi(x) = N \sin(n\pi x/2), \quad (n \text{ even}) \qquad E = \frac{\pi^2 n^2}{8}$

How do we proceed in a numerical integration?

Choose valid boundary conditionas at x=-1

 $\Psi(-1) = 0, \quad \Psi(-1 + \Delta_x) = A$

A is arbitrary (not 0); normalize after solution found

E = -5.0000000"Shooting method" Pick an energy E > Integrate to x=1 > Is boundary condition at x=1 satisfied? > If not, adjust E, х integrate again > Use bisection to refine

Solving an equation using bisection (general)

We wish to find the zero of some function

f(E) = 0

First find E_1 and E_2 bracketing the solution

 $f(E_1) < 0, \quad f(E_2) > 0$

Then evaluate the function at the mid-point value

 $E_3 = \frac{1}{2}(E_1 + E_2)$

Choose new bracketing values:

if $f(E_3) < 0$, then $E'_1 = E_3$, $E'_2 = E_2$ if $f(E_3) > 0$, then $E'_1 = E_1$, $E'_2 = E_3$

Repeat procedure with the new bracketing values - until $f(E_3) < \epsilon$

Bisection search for the ground state

- > First find E1, E2 giving different signs at x=+1
- \succ Then do bisection within these brackets



More complicated example:

Box with central Gaussian potential barrier



Ground state Search

First excited state



Potential well with non-rigid walls

Looking for bound state; $\Psi(x \to \pm \infty) \to 0$

Asymptotic solution: $\Psi(x) = Ae^{\alpha x} + Be^{-\alpha x}$. $\alpha = \sqrt{2(V_{\infty} - E)}$

A=0 for x>0 B=0 for x<0

Use the asymptotic form for two points far away from the center of the well

Find E for which the solution decays to 0 at the other boundary



Ground state search

Using criterion: $\Psi(1) = \Psi(-1)$

