

Hints for Solution of the Time-Dependent Schrödinger Equation

Consider the one-dimensional, time-dependent Schrödinger equation, on the domain $0 \leq x \leq 1$, $t \geq 0$, where we have chosen units such that $\hbar = m = 1$:

$$i \frac{\partial \psi(x, t)}{\partial t} = -\frac{\partial^2 \psi(x, t)}{\partial x^2} + V(x) \psi(x, t) \quad (1)$$

Use the following finite difference approximation (*Crank Nicholson* approximation) which is (a) second order in both space and time, and (b) *implicit*, meaning a *system* of linear equations must be solved at each time step:

$$i \frac{\psi_j^{n+1} - \psi_j^n}{\Delta t} = -\frac{1}{2} \left(\frac{\psi_{j+1}^{n+1} - 2\psi_j^{n+1} + \psi_{j-1}^{n+1}}{\Delta x^2} + \frac{\psi_{j+1}^n - 2\psi_j^n + \psi_{j-1}^n}{\Delta x^2} \right) + V_j \frac{1}{2} (\psi_j^{n+1} + \psi_j^n) \quad (2)$$

where $\psi_j^n \equiv \psi(j\Delta x, n\Delta t)$, $V_j \equiv V(j\Delta x)$, etc. Rewrite the above equations in the form

$$c_j^+ \psi_{j+1}^{n+1} + c_j^0 \psi_j^{n+1} + c_j^- \psi_{j-1}^{n+1} = S_j \quad j = 2, 3, \dots, \text{nx} - 1 \quad (3)$$

which, along with the boundary conditions (representing an infinite potential barrier at $x = 0$ and $x = 1$)

$$\psi_1^{n+1} = \psi_{\text{nx}}^{n+1} = 0 \quad (4)$$

constitute a *complex, tridiagonal, linear* system for the advanced values ψ_j^{n+1} . You will need to determine precise formulae for the c_j^+ , c_j^0 , c_j^- and S_j yourself, and note that these quantities will generally be complex.

Use `complex*16` (not `complex`) arithmetic, and the LAPACK solver `zgtsv` to solve the tridiagonal system. See the source code at

http://laplace.physics.ubc.ca/People/matt/410/Doc/linsys/src_lapack/zgtsv.f

for usage details, but note that the routine is available via

```
$(F77_LOAD) ... -llapack $(LIBBLAS) -o ...
```

as with the rest of the LAPACK routines.

Use

$$\Delta t = \lambda \Delta x \quad (5)$$

where $\lambda \sim 1$ —you can experiment with various values, but something like $\lambda = 0.5$ should work well.

Note that an excellent way of checking your implementation follows from the fact that there is a conservation law for $\psi(x, t)$ (conservation of total probability):

$$I \equiv \int_0^1 \psi(x, t) \psi^*(x, t) dx = \text{constant}. \quad (6)$$

You can approximate the integral using the following second order approximation:

$$I \approx \sum_{j=1}^{\text{nx}-1} \psi_{j+1/2}^n \psi_{j+1/2}^{*n} \Delta x = \sum_{j=1}^{\text{nx}-1} \frac{1}{2} (\psi_j^n + \psi_{j+1}^n) \frac{1}{2} (\psi_j^{*n} + \psi_{j+1}^{*n}) \Delta x \quad (7)$$