

Homework 1: Quantum Harmonic Oscillator

DUE: Friday, October 31, 2008

The probability density evolution in a one-dimensional harmonic trapping potential is governed by the partial differential equation:

$$i\hbar\psi_t + \frac{\hbar^2}{2m}\psi_{xx} - V(x)\psi = 0, \quad (1)$$

where ψ is the probability density and $V(x) = kx^2/2$ is the harmonic confining potential. A typical solution technique for this problem is to assume a solution of the form

$$\psi = \sum_1^N a_n \phi_n(x) \exp\left(i\frac{E_n}{2\hbar}t\right) \quad (2)$$

which is called an eigenfunction expansion solution (ϕ_n =eigenfunction, E_n =eigenvalue). Plugging in this solution ansatz to Eq. (1) gives the boundary value problem:

$$\frac{d^2\phi_n}{dx^2} - [Kx^2 - \varepsilon_n] \phi_n = 0 \quad (3)$$

where we expect the solution $\phi_n(x) \rightarrow 0$ as $x \rightarrow \pm\infty$ and ε_n is the quantum energy. Note here that $K = km/\hbar^2$ and $\varepsilon_n = E_n m/\hbar^2$. In what follows, take $K = 1$ and always normalize so that $\int_{-\infty}^{\infty} |\phi_n|^2 dx = 1$.

(a) Calculate the first five *normalized* eigenfunctions (ϕ_n) and eigenvalues (ε_n) using a shooting scheme. For this calculation, use $x \in [-L, L]$ with $L = 4$ and choose $xspan = -L : 0.1 : L$. Save the absolute value of the eigenfunctions in a 5-column matrix (column 1 is ϕ_1 , column 2 is ϕ_2 etc.) and the eigenvalues in a 1x5 vector.

ANSWERS: Should be written out as A1.dat (eigenfunctions) and A2.dat (eigenvalues)

(b) Calculate the first five *normalized* eigenfunctions (ϕ_n) and eigenvalues (ε_n) using a direct method. Be sure to use a forward- and backward-differencing for the boundary conditions (HINT: $3 + 2\Delta x\sqrt{KL^2 - \varepsilon_n} \approx 3$). For this calculation, use $x \in [-L, L]$ with $L = 4$ and choose $xspan = -L : 0.1 : L$. Save the absolute value of the eigenfunctions in a 5-column matrix (column 1 is ϕ_1 , column 2 is ϕ_2 etc.) and the eigenvalues in a 1x5 vector. NOTE: This procedure solves for the *interior* points. So be sure at the end to include your first and last point.

ANSWERS: Should be written out as A3.dat (eigenfunctions) and A4.dat (eigenvalues)

(c) There has been suggestions that in some cases, nonlinearity plays a role such that

$$\frac{d^2\phi_n}{dx^2} - [\gamma|\phi_n|^2 + Kx^2 - \varepsilon_n] \phi_n = 0. \quad (4)$$

Depending upon the sign of γ , the probability density is focused or defocused. Find the first two *normalized* modes for $\gamma = \pm 0.05$ using shooting. For this calculation, use $x \in [-L, L]$ with $L = 2$ and choose $xspan = -L : 0.1 : L$. Save the absolute value of the eigenfunctions in a 2-column matrix (column 1 is ϕ_1 and column 2 is ϕ_2) and the eigenvalues in a 1x2 vector.

ANSWERS: For $\gamma = 0.05$, should be written out as A5.dat (eigenfunctions) and A6.dat (eigenvalues)

ANSWERS: For $\gamma = -0.05$, should be written out as A7.dat (eigenfunctions) and A8.dat (eigenvalues)

(d) For a fixed value of the energy (specifically, take $\varepsilon_n = 1$ with $x \in [-L, L]$ and $L = 2$. For initial launch conditions, take $\phi = 1$ and $\phi_x = \sqrt{KL^2 - 1}$), perform a convergence study of the shooting method by controlling the error tolerance, TOL, in ODE45 and ODE23:

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TOL=1e-4;
OPTIONS = odeset('RelTol',TOL,'AbsTol',TOL);
[T,Y] = ODE45('F',TSPAN,YO,OPTIONS);
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Show that indeed the schemes are fourth order and second order respectively by running the computation across the computational domain and adjusting the tolerance. Use specifically the following values of $TOL = [1e-4 \ 1e-5 \ 1e-6 \ 1e-7 \ 1e-8 \ 1e-9 \ 1e-10]$. In particular, plot on a log-log scale the average step-size (x-axis) using the *diff* and *mean* command versus the tolerance (y-axis) for the above tolerance values. What are the slopes of these lines? Use the POLYFIT command to get the slopes. Note that the local error should be $O(\Delta t^5)$ and $O(\Delta t^3)$ respectively. What are the local errors for ODE113 and ODE15s? Save the slopes computed from the POLYFIT command with the log-log of the data as a 4x1 vector with the slopes of ODE45, ODE23, ODE113 and ODE15s respectively.

ANSWERS: Should be written out as A9.dat

(e) Compare your solutions in both (a) and (b) with the exact Gauss-Hermite polynomial solutions for this problem (See wikipedia.com, for instance). Compute the error between your numerical solution and the exact solution for the values of the eigenfunctions and eigenvalues computed above. Specifically, calculate the following quantity for each eigenfunction $\| |\phi_n^{numerical}| - |\phi_n^{exact}| \|$ where $\|f(x)\| = \int_{-L}^L f(x)^2 dx$. For the eigenvalues, simply calculate the relative percent error $100 \times (|\varepsilon_n^{numerical} - \varepsilon_n^{exact}| / \varepsilon_n^{exact})$. The error vectors associated with the eigenfunctions and eigenvalues should be 5x1 vectors.

ANSWERS: For part (a), should be written out as A10.dat (eigenfunctions) and A11.dat (eigenvalues)

ANSWERS: For part (b), should be written out as A12.dat (eigenfunctions) and A13.dat (eigenvalues)

(f) For a double-well potential (replace kx^2 in (3) with the potential below so that $kx^2 \rightarrow 2V(x)$.)

$$V(x) = \begin{cases} -10 & 1 < x < 2 \text{ and } -2 < x < -1 \\ 0 & \text{otherwise} \end{cases}, \quad (5)$$

calculate the symmetric ground-state ϕ_1 and first, anti-symmetric state ϕ_2 using the method of part (b). Note that $|\varepsilon_1 - \varepsilon_2| \ll 1$, thus the difficulty in this problem. Make a 3D plot $(x, t, |\psi|)$ of the solution (2) for the initial condition $\psi = \phi_1 + \phi_2$ which shows the tunneling between potential wells. Note that for plotting purposes, take $\hbar = 1$ and $m = 1$. For this calculation, use $x \in [-L, L]$ with $L = 4$ and choose $xspan = -L : 0.1 : L$. Save the absolute value of the two eigenfunctions in a 2-column matrix (column 1 is ϕ_1 and column 2 is ϕ_2) and the eigenvalues in a 1x2 vector.

ANSWERS: Should be written out as A14.dat (eigenfunctions) and A15.dat (eigenvalues)

NOTE: For the tolerances in both the convergence (shooting method) of the ε_n in (a) and (c), and in the area for part (c), use a tolerance of 10^{-4} . As a final remark, my code takes 70 seconds.