

Assignment 4

Quantum Harmonic Oscillator

November 5th 2024

Exercise: Solve the time independent Schroedinger equation for the quantum harmonic oscillator

$$H = \frac{\hat{p}^2}{2m} + \frac{m\omega^2 \hat{x}^2}{2}$$

Scaling of the matrix-matrix multiplication. Consider the code developed in the Exercise 3 from Assignment 1 (matrix-matrix multiplication):

- (a) Write a program to compute the first E_k eigenvalues and $|\Psi_k\rangle$ eigenfunction.
- (b) How would you rate your program in terms of the priorities we introduced in class for good scientific software development (Correctness, Stability, Accurate discretization, Flexibility, Efficiency)?

Setting:

Units: $m = \hbar = 1$

1D: $\hat{p} = -i\hbar \nabla_x$

The initial equation reads:

$$H = \frac{1}{2}(-\nabla_x^2 + \omega^2 x^2)$$

One needs to solve the eigenvalue equation:

$$H|\Psi_k(x_i)\rangle = E_k|\Psi_k(x_i)\rangle$$

$$\left[\frac{1}{2}(-\nabla_x^2 + \omega^2 x^2)\right]|\Psi_k(x_i)\rangle = E_k|\Psi_k(x_i)\rangle$$

Method: **FINITE DIFFERENCE METHOD**

$$\left[\frac{1}{2}(-\nabla_x^2 + \omega^2 x^2)\right] |\Psi_k(x_i)\rangle = E_k |\Psi_k(x_i)\rangle$$

-find approximate solutions by transforming the differential equation into a matrix equations;

- x-axis discretization: divide the interval [a,b] in N intervals: $\Delta x = \frac{b-a}{N}$

$$\left[\frac{1}{2}(-\nabla_x^2 + \omega^2 x^2)\right] = \hat{K} + \hat{V}$$

$$\text{Kinetic: } \frac{-\Psi(x_{i+1}) + 2\Psi(x_i) - \Psi(x_{i-1}))}{2\Delta x^2} + \mathcal{O}(\Delta x^2)$$

$$\text{Potential: } V(x_i) = \frac{\omega^2 x_i^2}{2}$$

$$\hat{K} = \frac{1}{2\Delta x^2} \begin{pmatrix} 2 & -1 & 0 & \dots & 0 & 0 \\ -1 & 2 & -1 & \dots & 0 & 0 \\ 0 & -1 & 2 & \dots & 0 & 0 \\ \vdots & \vdots & \vdots & \dots & \vdots & \vdots \\ 0 & 0 & 0 & \dots & 2 & -1 \end{pmatrix}$$

TRIDIAGONAL

$$\hat{V} = \frac{\omega^2}{2} \begin{pmatrix} x_1^2 & 0 & 0 & \dots & 0 & 0 \\ 0 & x_2^2 & 0 & \dots & 0 & 0 \\ 0 & 0 & x_3^2 & \dots & 0 & 0 \\ \vdots & \vdots & \vdots & \dots & \vdots & \vdots \\ 0 & 0 & 0 & \dots & 0 & x_N^2 \end{pmatrix}$$

Method: **FINITE DIFFERENCE METHOD** (higher order)

$$\left[\frac{1}{2}(-\nabla_x^2 + \omega^2 x^2)\right] |\Psi_k(x_i)\rangle = E_k |\Psi_k(x_i)\rangle$$

-find approximate solutions by transforming the differential equation into a matrix equations;

- x-axis discretization: divide the interval [a,b] in N intervals

$$\hat{K} = \frac{1}{2\Delta x^2} \begin{pmatrix} 5/2 & -4/3 & 1/12 & \dots & 0 & 0 \\ -4/3 & 5/2 & -4/3 & \dots & 0 & 0 \\ 1/12 & -4/3 & 5/2 & \dots & 0 & 0 \\ \vdots & \vdots & \vdots & \dots & \vdots & \vdots \\ 0 & 0 & 0 & \dots & 5/2 & -4/3 \end{pmatrix}$$

$$\left[\frac{1}{2}(-\nabla_x^2 + \omega^2 x^2)\right] = \hat{K} + \hat{V}$$

$$\text{Kinetic: } \frac{-\Psi(x_{i+2}) + 16\Psi(x_{i+1}) - 30\Psi(x_i) + 16\Psi(x_{i-1}) - \Psi(x_{i-2}))}{12\Delta x^2} + \mathcal{O}(\Delta x^4)$$

$$\text{Potential: } V(x_i) = \frac{\omega^2 x_i^2}{2}$$

PENTADIAGONAL

$$\hat{V} = \frac{\omega^2}{2} \begin{pmatrix} x_1^2 & 0 & 0 & \dots & 0 & 0 \\ 0 & x_2^2 & 0 & \dots & 0 & 0 \\ 0 & 0 & x_3^2 & \dots & 0 & 0 \\ \vdots & \vdots & \vdots & \dots & \vdots & \vdots \\ 0 & 0 & 0 & \dots & 0 & x_N^2 \end{pmatrix}$$

Discussion: ANALYTICAL COMPARISON

$$\hat{H} |\psi_n\rangle = \left(\frac{\hat{p}^2}{2m} + \frac{1}{2}m\omega^2 \hat{x}^2 \right) |\psi_n\rangle = E_n |\psi_n\rangle$$

$$E_n = \hbar\omega \left(n + \frac{1}{2} \right) \quad \text{eigenvalues}$$

$$\psi_n(x) = \frac{1}{\sqrt{2^n n!}} \left(\frac{m\omega}{\pi\hbar} \right)^{1/4} e^{-\frac{m\omega x^2}{2\hbar}} H_n \left(\sqrt{\frac{m\omega}{\hbar}} x \right), \quad n = 0, 1, 2, \dots$$

eigenfunctions

$$H_n(x) = (-1)^n e^{z^2} \frac{d^n}{dz^n} (e^{-z^2})$$

Hermite polynomials

Discussion: SELF EVALUATION

Correctness —> comparison with analytical solution

Stability —> is the solution stable on different runs?

Accurate discretization —> what does it happen with smaller discretization step?

Flexibility —> can you easily add new functionalities? i.e. higher order

Efficiency —> runtimes?