

Quantum Information and Computing

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Saverio Monaco
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Solution of the Time Independent
Schrödinger Equation for
an harmonic oscillator

Theory

$$\left\{ \begin{array}{l} \hat{H} = \hat{p}^2 + \omega^2 \hat{x}^2 \quad \text{where } \hbar \equiv m \equiv 1 \\ \hat{H}\psi = E\psi \\ \hat{p} \rightarrow -i\hbar\partial/\partial x, \quad \hat{x} \rightarrow x \end{array} \right. \rightarrow \left\{ \begin{array}{l} \left(-\frac{1}{2} \frac{\partial^2}{\partial x^2} + \frac{1}{2} \omega^2 x^2 \right) \psi(x) = E_n \psi(x) \\ \psi_k'' = \frac{\psi_{k+1} - 2\psi_k + \psi_{k-1}}{dx^2} \end{array} \right.$$

$$\equiv -\frac{1}{2} \left[\frac{\psi_{k+1} - 2\psi_k + \psi_{k-1}}{dx^2} \right] + \frac{1}{2} \omega^2 x_k \psi_k = E \psi_k$$

$$H = \frac{1}{2} \begin{pmatrix} \frac{2}{dx^2} + \omega^2 x_1^2 & -\frac{1}{dx^2} & 0 & \cdots & 0 \\ -\frac{1}{dx^2} & \frac{2}{dx^2} + \omega^2 x_2^2 & -\frac{1}{dx^2} & \cdots & 0 \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & 0 & \cdots & \frac{2}{dx^2} + \omega^2 x_N^2 \end{pmatrix} \longleftrightarrow H\psi = E\psi$$

Code Development

qho_H_init() in module qho:

Initialize the complex matrix to represent the hamiltonian of the system given L and N

```
function qho_H_init(L,N,omega) result(H)
    real                                :: L,omega,dx
    integer                             :: N, ii
    real*16, dimension(:,,:), allocatable :: elem_real
    type(cmatrix)                       :: H
    dx = L/N

    allocate(elem_real(N+1,N+1))
    elem_real = 0 * elem_real ! Initialize everything to 0

    do ii=1, N+1, 1 ! ! diagonal elements
        elem_real(ii,ii) = ( 2 / (dx**2) ) + (omega**2)*((ii-1)*dx - L/2)**2
    end do

    do ii=2, N+1, 1 ! tridiagonal elements
        elem_real(ii,ii-1) = - 1/(dx**2)
        elem_real(ii-1,ii) = - 1/(dx**2)
    end do

    elem_real = 0.5* elem_real ! Everything must be divided by 2

    H = cmatrix_init(cmplx(X=elem_real,KIND=8))
end function qho_H_init
```

ZGEEV() function in

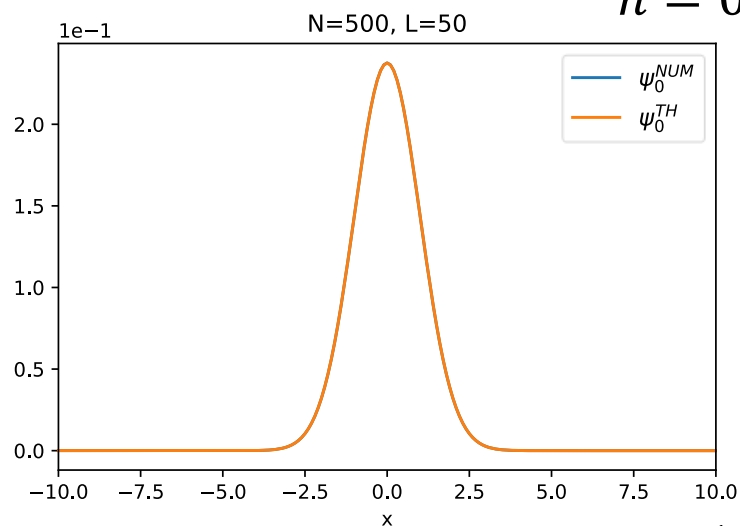
LLAPACK library:

For eigenvalues and eigenvectors

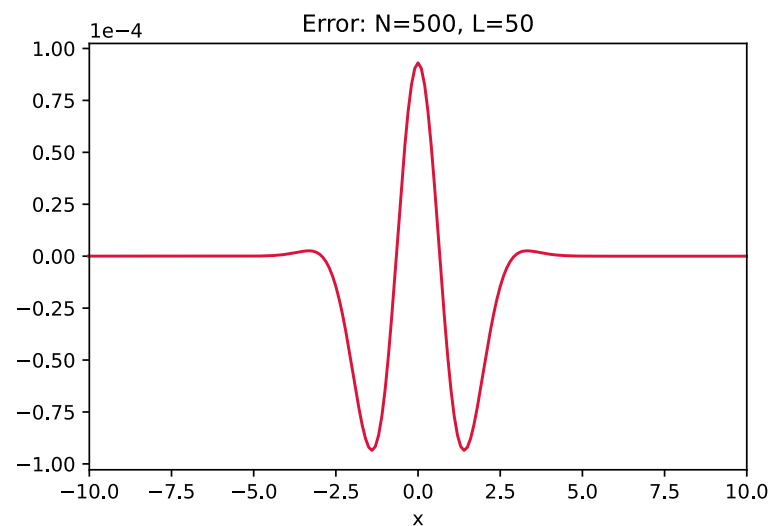
```
call ZGEEV(! To compute both eigenvalues and -vectors
            'V', 'V', &
            ! The order of the matrix A
            N, &
            ! The matrix
            A, &
            ! The leading dimension of A.
            N, &
            ! Where to store eigenvalues and eigenvectors
            eigenv, eigenh, &
            ! Other parameters
            N, VR, N, WORK,LWORK,RWORK, &
            ! Output parameter, if INFO == 0 -> successful exit
            INFO )
```

Results

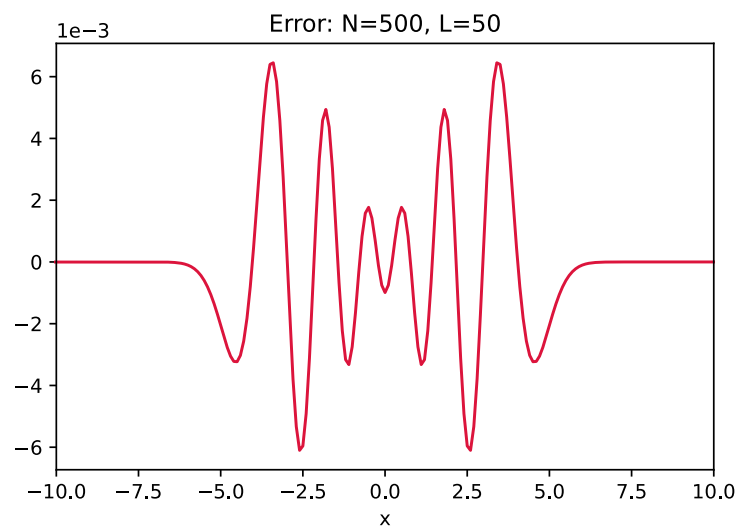
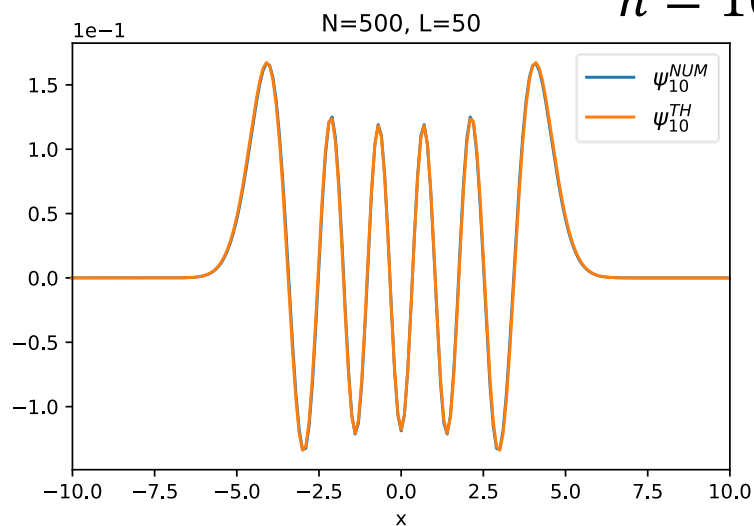
$n = 0$



$$err \equiv \psi_i^{num} - \psi_i^{th}$$



$n = 10$



Self evaluation

Correctness: Hugely depends on the choice of the parameters. It can be still improved by using a better formula for the derivative

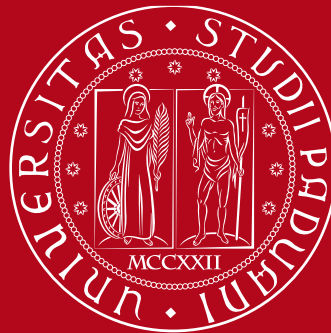
Stability: Various checks have been implemented to make the program as stable as possible

Accurate discretization: For the right parameters (for example $L=50$ and $N=1000$) the programs gives accurate results

Flexibility: Based on ZGEEV and not DSTEMR for generalizing to other problems

Efficiency: Can be improved by considering DSTEMR and real-only matrices

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DI PADOVA

Thanks for the attention
