

FYS4150 - Computational Physics

Project 2

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1 Introduction

In this project, we will have a look at the so called *Eigenvalue problem*. We will use the well known Schrödinger equation from quantum mechanics and apply it to a system consisting of a single electron in a harmonic oscillator potential. To achieve this, we will implement Jacobi's method to solve the Schrödinger equation and then find the eigenvalues, which will physically be the energies of the electrons.

We will also look at the case where we have two interacting electrons in a three dimensional harmonic oscillator, with different frequencies.

2 Method

2a) Mathematical intermezzo

We first want to look at the orthogonality properties of a specific basis. Let us consider the following basis of vectors

$$\mathbf{v}_i = \begin{pmatrix} v_{i1} \\ v_{i2} \\ \vdots \\ v_{in} \end{pmatrix} \quad (1)$$

We will also assume that this is an orthogonal basis

$$\mathbf{v}_j^T \mathbf{v}_i = \delta_{ij} \quad (2)$$

We want to look at the orthogonality properties for a orthogonal and unitary vector transformation in the form

$$\mathbf{w}_i = \mathbf{U}\mathbf{v}_i \quad (3)$$

For an orthogonal matrix, we know that $\mathbf{U}^T\mathbf{U} = \mathbf{I} = 1$, where \mathbf{I} is the identity matrix, so we have that

$$\mathbf{w}_j^T \mathbf{w}_i = (\mathbf{U}\mathbf{v}_i)^T (\mathbf{U}\mathbf{v}_j) = \mathbf{v}_j^T \mathbf{U}^T \mathbf{U} \mathbf{v}_i = \mathbf{v}_j^T \mathbf{v}_i = \delta_{ij} \quad (4)$$

For a unitary matrix, we have that $\mathbf{U}^*\mathbf{U} = \mathbf{I}$ and $\mathbf{U}^\dagger\mathbf{U} = \mathbf{I}$, where $\mathbf{U}^\dagger = (\mathbf{U}^*)^T$ is the hermitian conjugate. Doing the hermitian conjugate on the transformation vector, we can show that

$$\mathbf{w}_j^\dagger \mathbf{w}_i = (\mathbf{U}\mathbf{v}_j)^\dagger (\mathbf{U}\mathbf{v}_i) = \mathbf{v}_j^\dagger \mathbf{U}^\dagger \mathbf{U} \mathbf{v}_i = \mathbf{v}_j^\dagger \mathbf{v}_i = \delta_{ij} \quad (5)$$

Here we have assumed that the basis vector is real, so $\mathbf{v}_i^* = \mathbf{v}_i$.

2b) Implementing Jacobi's method for the non-interacting case

3 Implementation

4 Results

5 Conclusion

6 References

M. Hjort-Jensen, 2015, *Computational Physics*, accessible at course GitHub repository; <https://github.com/CompPhysics/ComputationalPhysics/tree/master/doc/Lectures> (as of 14.09.16), 551 pages.