

# Homework 0: LaTeX & GitHub Demonstration:

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I do not have an abstract here.

## I. LINKS TO CODE

This is the link to the github repository with the code. [Click here](#)

### A. Split Step Algorithm

This report is a short overview on solving the nonlinear Schrodinger equation. In the Split Operator method, we split the Hamiltonian into momentum space components,  $H_p = \frac{-\hbar^2}{2m} \nabla_z^2$  and position space components,  $H_r = V(x)$ . If we assume a somewhat general solution to our quantum system:

$$\psi(x, t + dt) = e^{-\frac{i\hat{H}dt}{\hbar}} \psi(x, t) = e^{\frac{-i(\hat{H}_p + \hat{H}_r)dt}{\hbar}} \psi(x, t) \quad (1)$$

and assume we are simulating our system by a series of small timesteps ( $dt$ ), we can perform similar splitting by

using the Baker-Campbell-Hausdorff formula:

$$\psi(x, t+dt) = e^{\frac{-i(\hat{H}_p)dt}{\hbar}} e^{\frac{-i(\hat{H}_r)dt}{\hbar}} e^{\frac{-i([\hat{H}_p, \hat{H}_r])dt^2}{2\hbar}} \psi(x, t) + \mathcal{O}(dt^2) \quad (2)$$

This accrues a small amount of error ( $dt^2$ ) related to the commutation of the real and momentum-space components of the Hamiltonian. In order to change the ( $dt^2$ ) error to ( $dt^3$ ), we can split the system by performing a half-step in position space before doing a full-step in momentum space, through a process called Strang Splitting:

$$\psi(x, t + dt) = e^{\frac{-i(\hat{H}_r)dt}{2\hbar}} e^{\frac{-i(\hat{H}_p)dt}{\hbar}} e^{\frac{-i(\hat{H}_r)dt}{2\hbar}} \psi(x, t) + \mathcal{O}(dt^3) \quad (3)$$

We can then address each part of this solution in parts, first in position space, then in momentum space, then in position space again by using Fourier Transforms.

$$\psi(x, t+dt) = \left[ \hat{U}_r(dt/2) \mathcal{F}^{-1} \left[ \hat{U}_p(dt) \mathcal{F} \left[ \hat{U}_r(dt/2) \psi(x, t) \right] \right] \right] + \mathcal{O}(dt^3) \quad (4)$$

where  $\hat{U}_r(dt) = e^{\frac{-i(\hat{H}_r)dt}{\hbar}}$  and  $\hat{U}_p(dt) = e^{\frac{-i(\hat{H}_p)dt}{\hbar}}$ .