

Report for movingFreeSoliton

Simulated with: lib.managers.crankNicolson.dimensionless

Simulation constants:

baseDensity: 1.000	chemicalPotential: 0.010	dt: 0.005
dx: 0.200	g: -0.010	hbar: 1.000
healingLength: 7.071	mass: 1.000	plotFPS: 1000.000
plotPause: 0.001	plotStep: 10	plotYMax: 2
plotYMin: -2	r: 0.125	tCount: 1000
tMax: 5	tMin: 0	velocity: 1.000
x0: 0.000	xCount: 100	xMax: 10
xMin: -10		

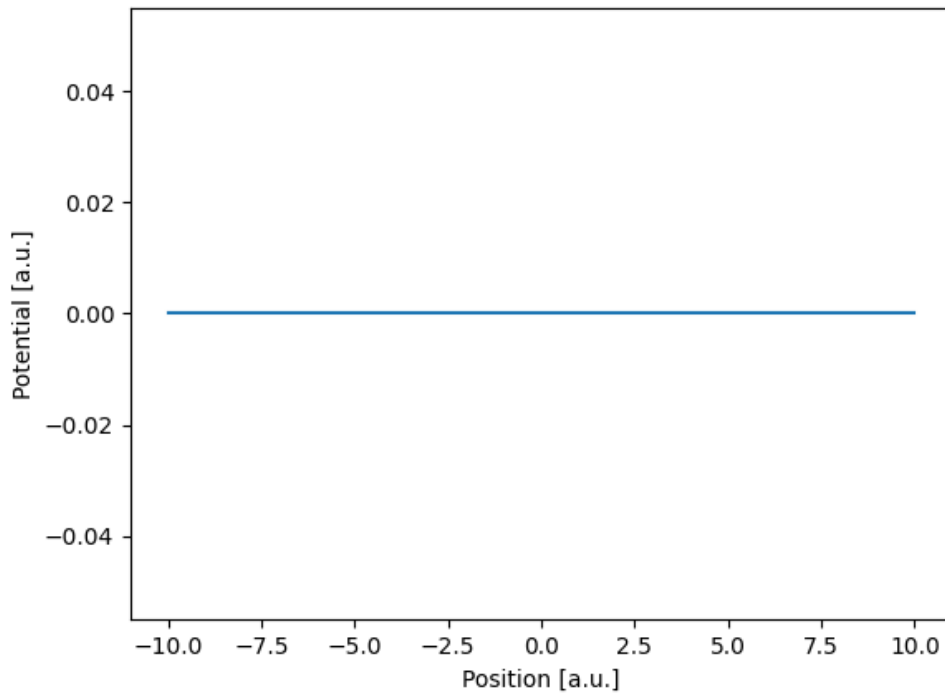
Wave function:

```
def brightSoliton(x, t, constants):    v = constants["velocity"]    g =
constants["g"]    eta = jnp.sqrt((v**2 + 2) / (-2 * g))    kappa = jnp.sqrt(2
/ (v**2 + 2))    spacePart = eta / jnp.cosh((x - v * t) / kappa) * jnp.exp(1j
* x * v)    timePart = jnp.exp(1j * (1 / 2 - v**2 / 4) * t)    return
spacePart * timePart
```

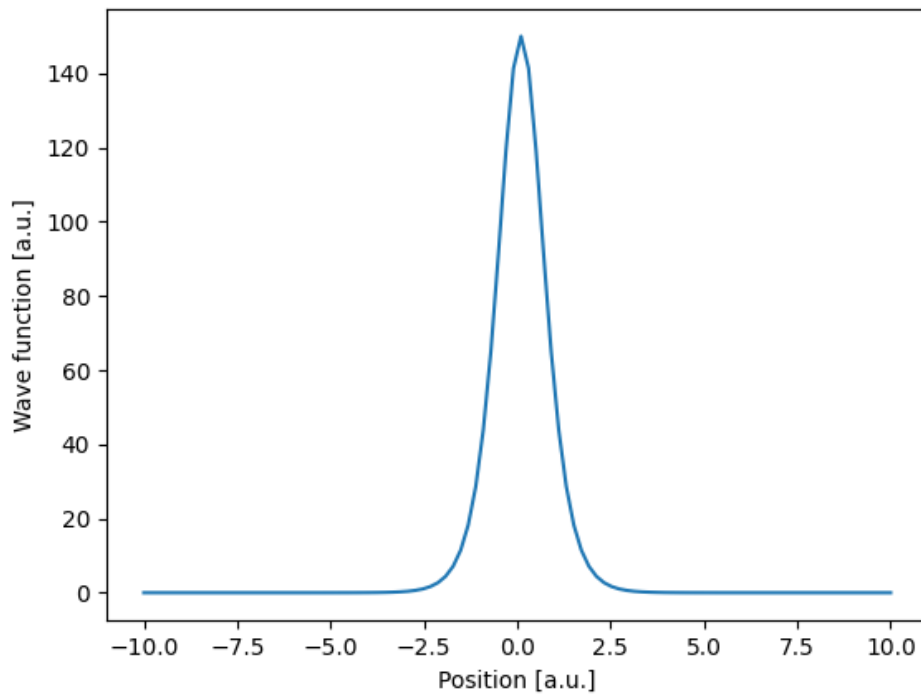
Potential function:

```
def V(x, t, constants):    return jnp.zeros_like(x)
```

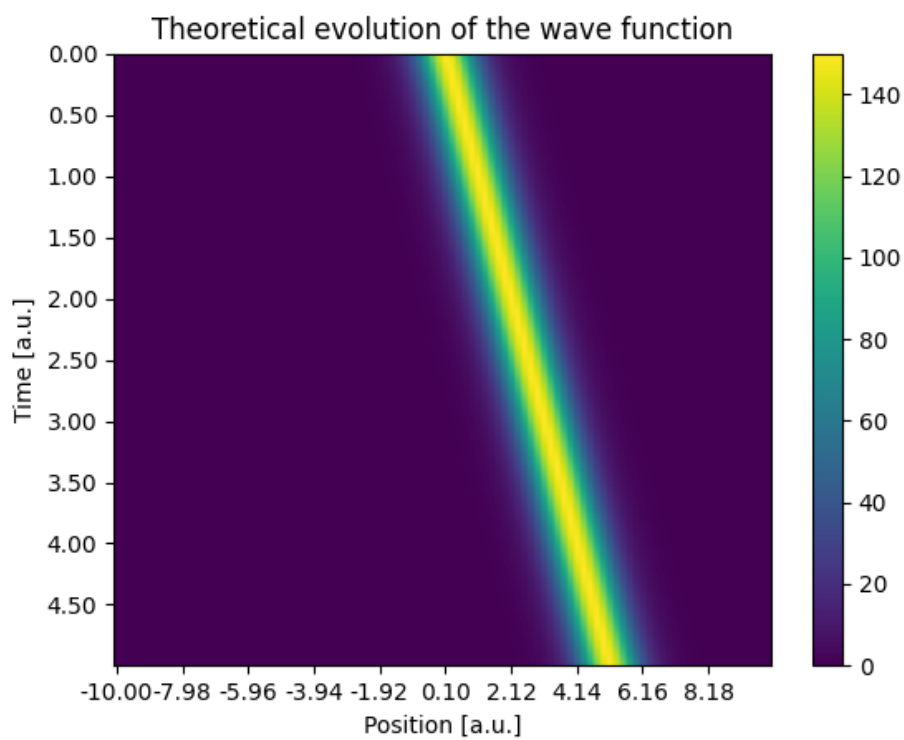
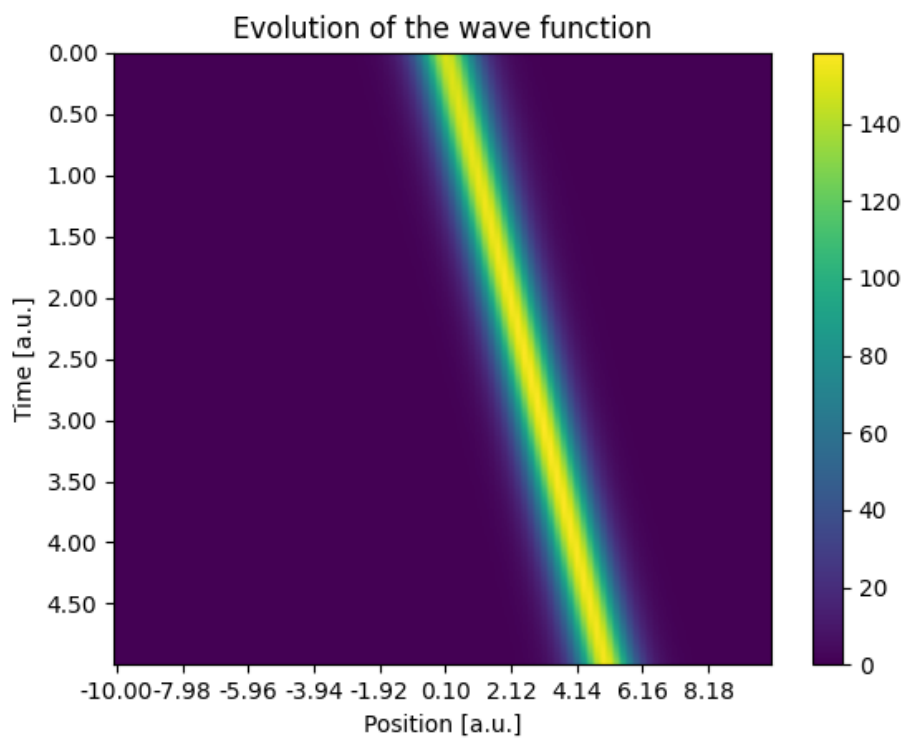
Potential at time $t=0$



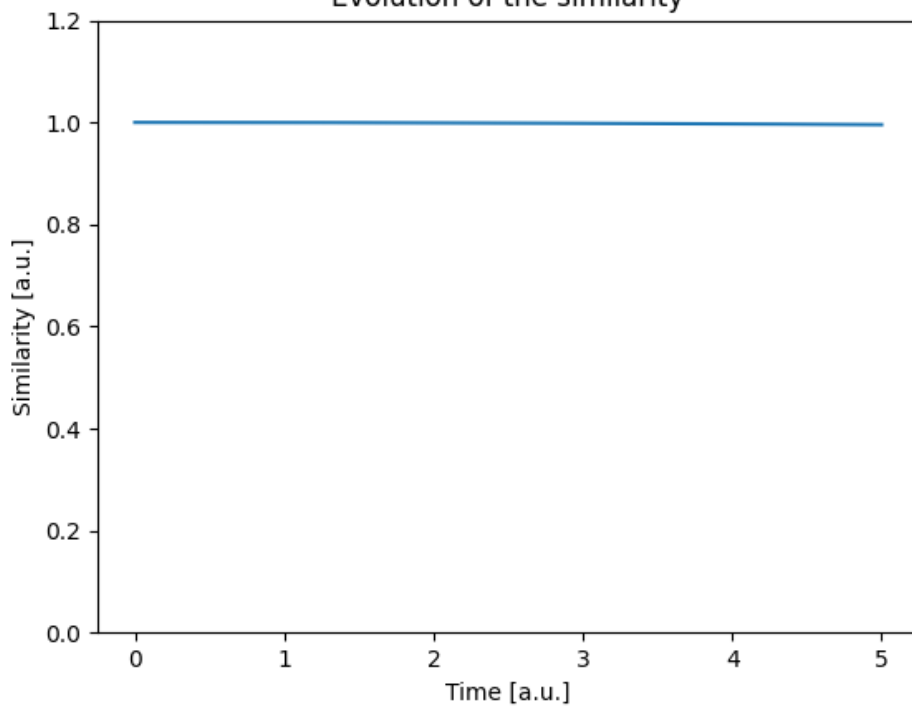
Wave function at time $t=0$



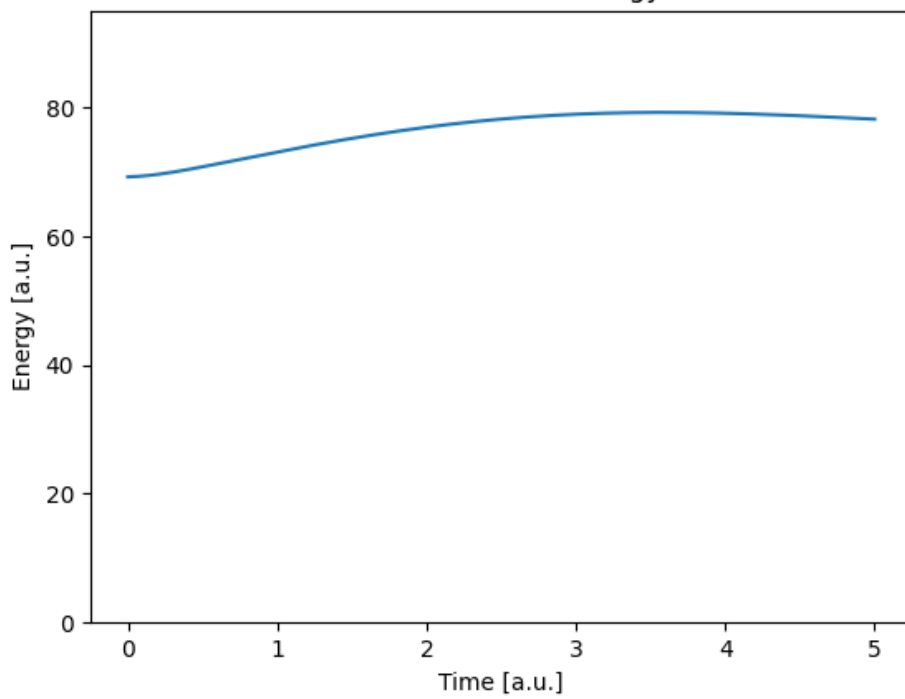
Results



Evolution of the similarity



Evolution of the energy



Evolution of the norm

