

1D Fast oscillations (Numerical)

$$C(t) = \bar{C} + A \sin\left(\frac{2\pi t}{T}\right)$$

$$T \gg \tau_{linear} \sim \bar{C}^{-1}$$

where we do not expect different dynamics if $A > \bar{C}$ or $A < \bar{C}$.

Kink dynamics

$$\dot{d}(t) \simeq -24\sqrt{2}\bar{C}^{\frac{1}{2}}e^{-2^{\frac{1}{2}}\bar{C}^{\frac{1}{2}}d}$$

Notice: It is the same formula in [1D Constant \(Numerical\)](#) but now $C \rightarrow \bar{C}$.

The variation of the distance over a period (assuming the distance to be constant inside the integrand)

$$\Delta d(d) = \int_0^T (\partial_t d) dt = -24\sqrt{2} \int_0^T e^{d\sqrt{2}\bar{C}^{1/2}} \bar{C}^{1/2} dt$$

Simulation

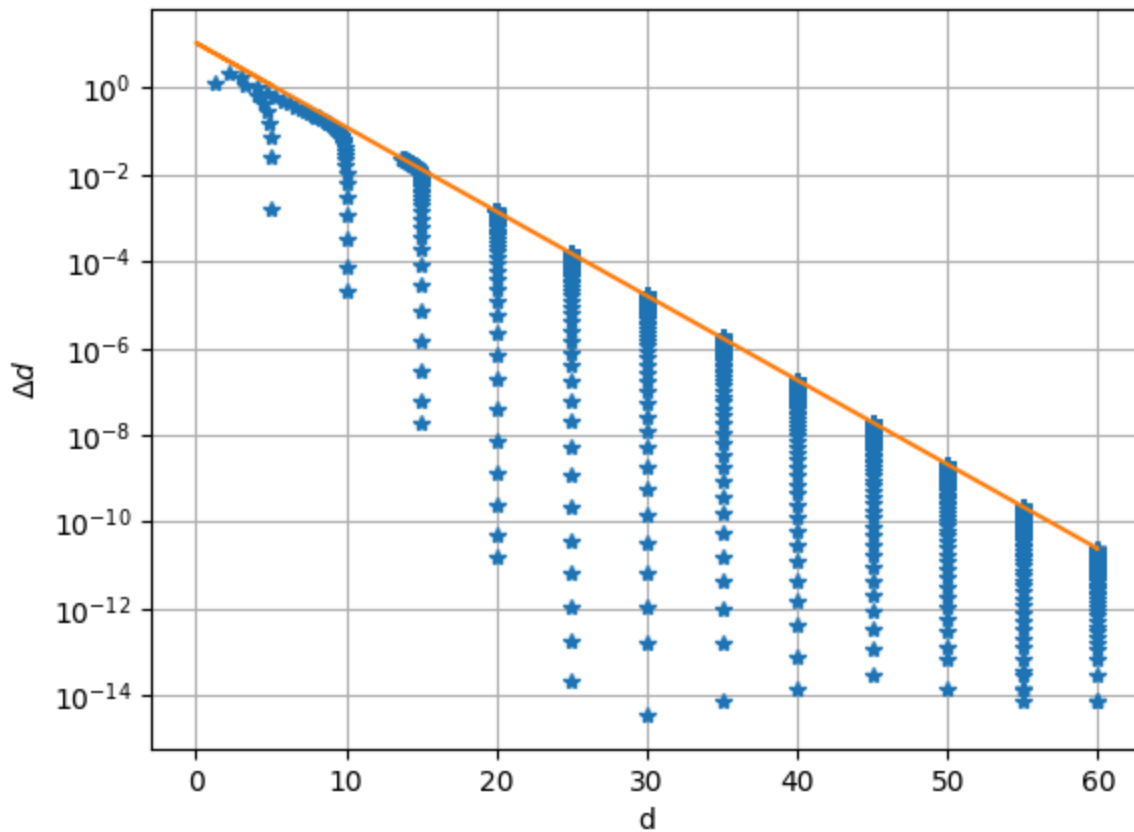
In all the following simulations, we adopted $T = 10\bar{C}^{-1}$, such that $T \gg \bar{C}^{-1}$.

The orange line is the formula above, while the blue points are experimental values. Those values are obtained running a simulation of $100T$ "seconds" for many different values of the initial distance. Δd is calculated for each period, for each simulation.

Variation of the distance of twokinks over one period T

tspan=100*T; dt = T/100

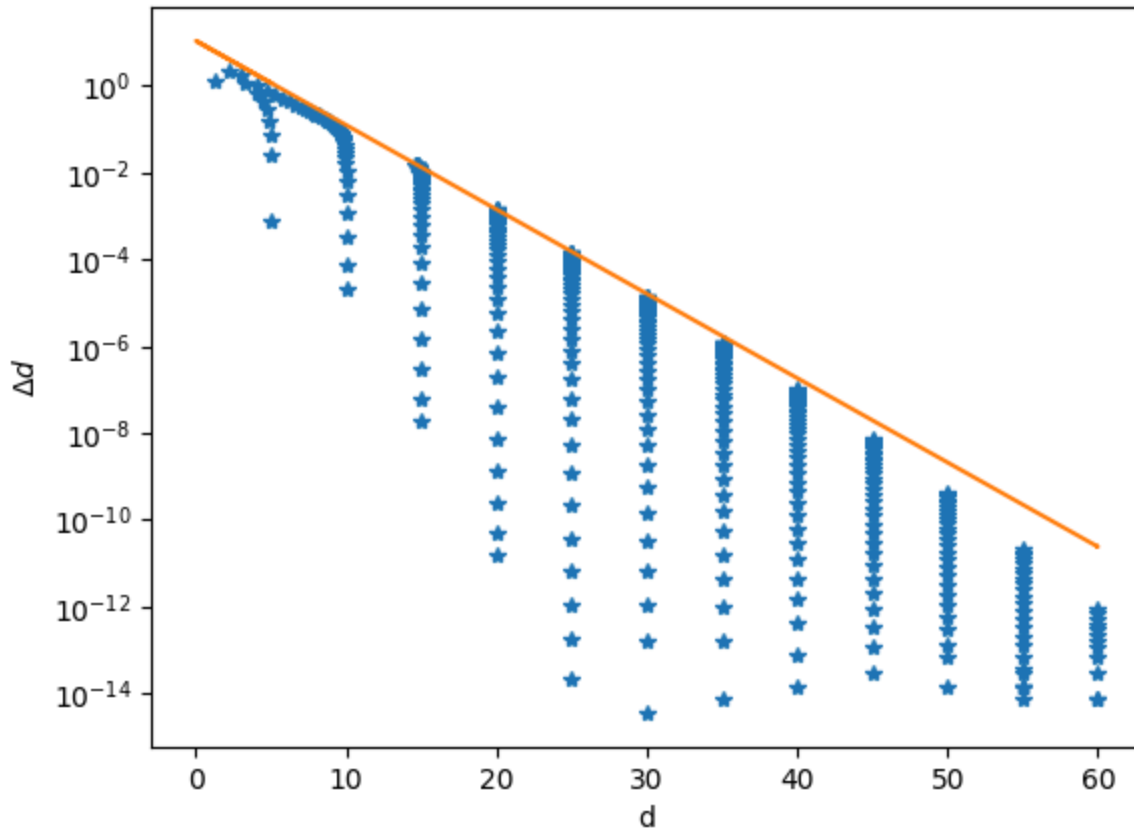
$$C(t) = 0.1 + 1\sin(2\pi t/1)$$



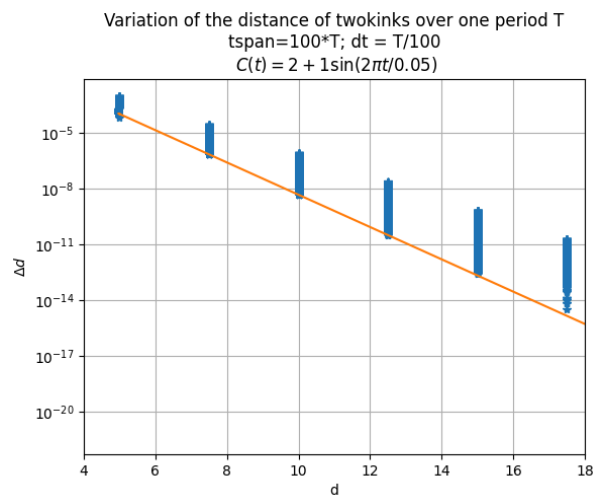
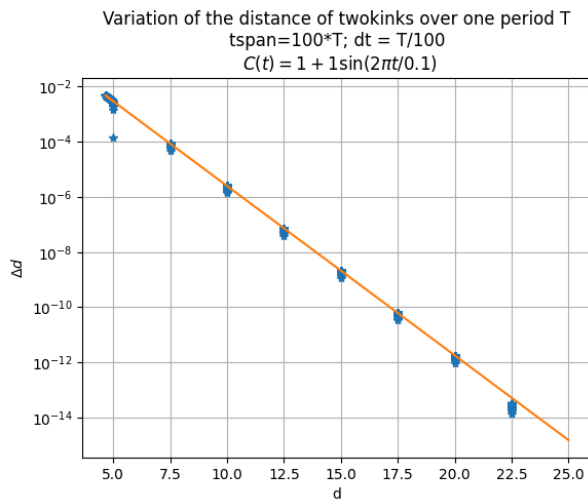
Notice: The way the blue dots "reach" the model tells something: We need to wait some periods before the simulation matches the model. Probably this is related to the **initial state preparation**. Here we show the same plot above, but each simulation lasts 50T instead of

100T.

Variation of the distance of twokinks over one period T
 $tspan=50*T$; $dt = T/100$
 $C(t) = 0.1 + 1\sin(2\pi t/1)$



Other simulations



Linear dynamics

$$\ell = \frac{2\pi}{\langle q^2 \rangle^{1/2}} \sim t^{1/2}$$

$$\tau_{linear} \sim \bar{C}^{-1}$$

Starting from random initial state

$$\ell = 2\pi / \langle q^2 \rangle^{1/2}$$

