

# Large-Scale Numerical Investigations into the Dynamics of Nonlinear Classical Systems

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Introduction

Numerical simulations

- Choosing an integrator

- The maximal Lyapunov exponent

Conclusions

# Acknowledgements

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- The author has been supported by the research project PN-III-P4-ID-PCE-2016-0792 funded by the Romanian Ministry of Research and Innovation.
- All numerical simulations were performed on the computing cluster of Department of Computational Physics and Information Technologies, “Horia Hulubei” National Institute for Physics and Nuclear Engineering.

# Introduction

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# The model

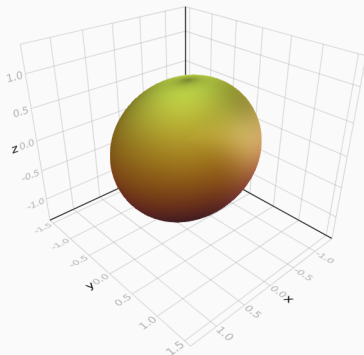
- The physical system that we model is the surface of heavy nuclei.
- The Hamiltonian describes the constrained motion of the vibrational quadrupole degrees of freedom of nuclear surface.

# The model

The Hamiltonian of the system

$$H = \frac{A}{2} (p_0^2 + p_2^2) + \frac{A}{2} (q_0^2 + q_2^2) + \frac{B}{\sqrt{2}} q_0 (3q_2^2 - q_0^2) + \frac{D}{4} (q_0^2 + q_2^2)^2$$

- Harmonic oscillator part
- Integrable part
- Non-integrable term

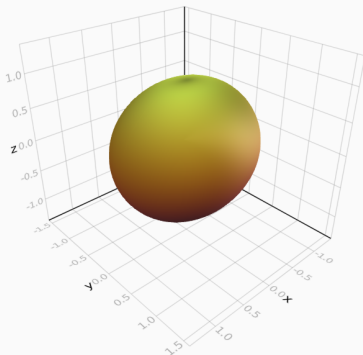


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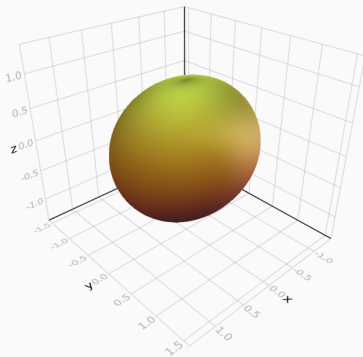


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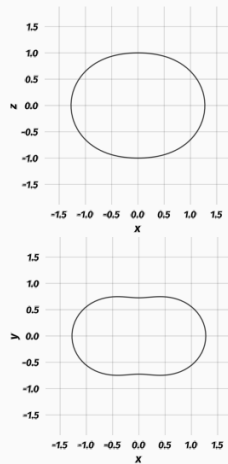
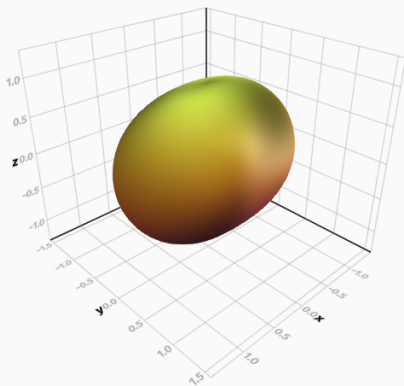
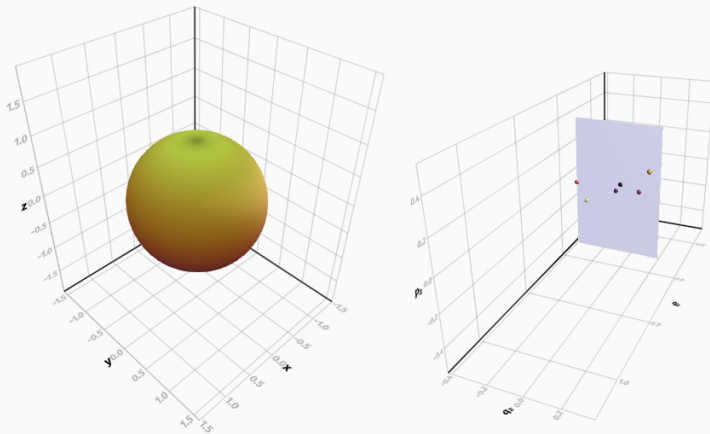


Figure 1: The nuclear surface and its sections

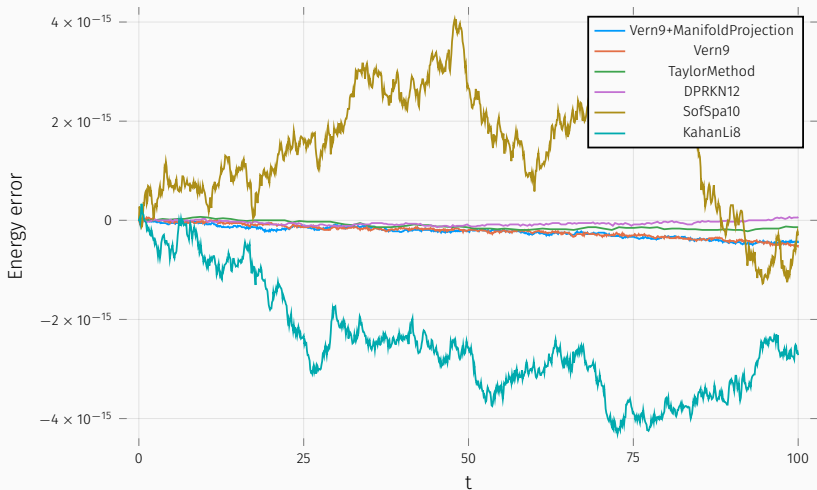


**Figure 2:** The nucleus and the corresponding trajectory in the phase space for a chaotic trajectory with  $B = 0.5, E = 0.3$

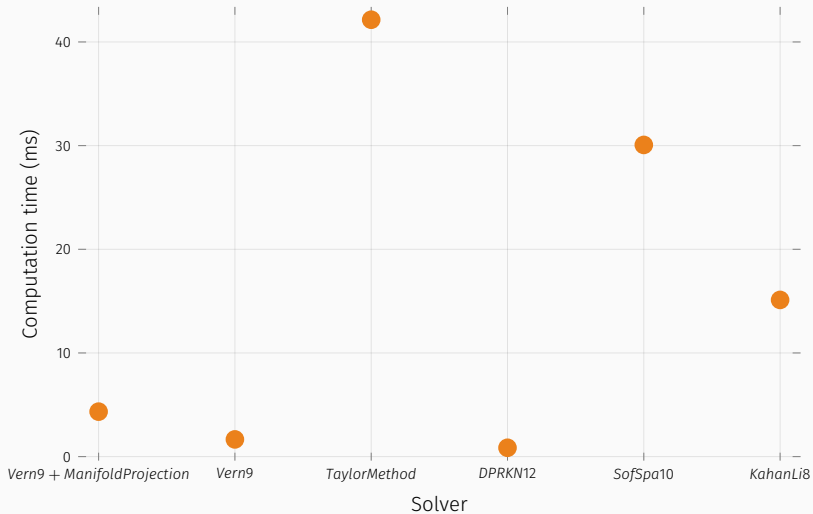
# Numerical simulations

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- Numerical simulations and the visualizations of the results was done in Julia (using `DifferentialEquations.jl` and `DynamicalSystems.jl`).
- Having access to the implementations of a large number of integrators helps us taking an informed decision for the choice of the integration algorithm.



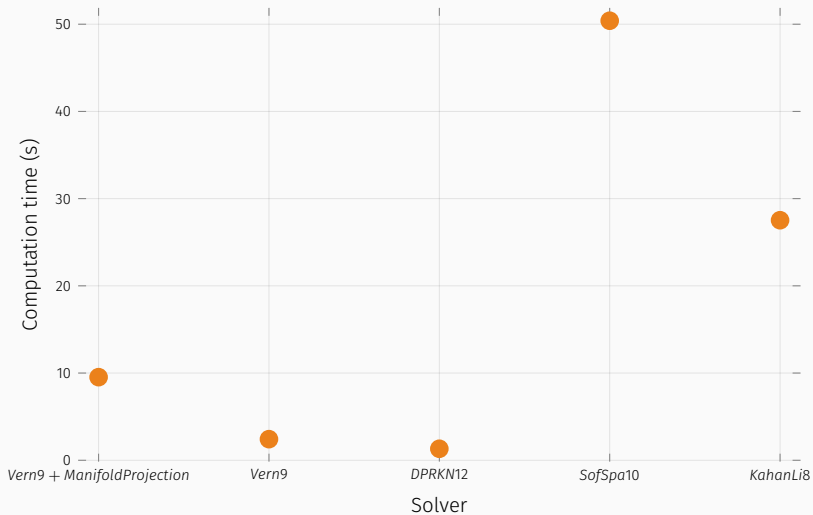
**Figure 3:** Energy error benchmark for short integration time



**Figure 4:** Computational time benchmark for short integration time



**Figure 5:** Energy error benchmark for long integration time



**Figure 6:** Computational time benchmark for long integration time

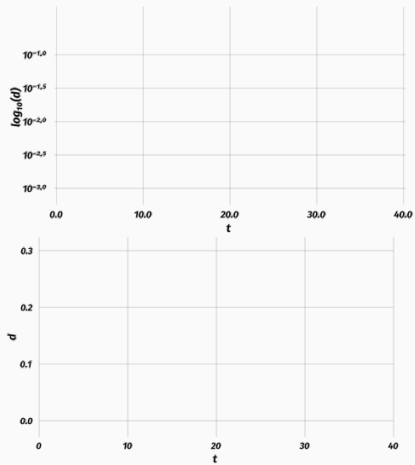
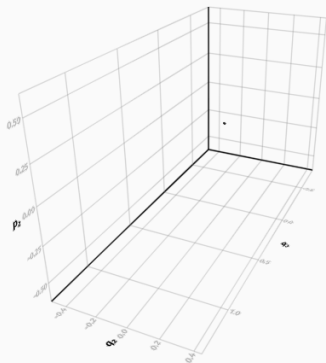


# Quantifying chaoticity

- The maximal Lyapunov exponent is defined as

$$\lambda = \lim_{t \rightarrow \infty} \frac{1}{t} \ln \frac{dx(t)}{dx(0)}$$

- One of the signature characteristics of chaos is the sensitivity to the initial conditions ( $\lambda > 0$ ).
- We can get some intuitive insight for the sensitivity to the initial conditions by following the evolution of two nearby initial conditions.



**Figure 7:** The distance between nearby trajectories for  $B = 0.5, E = 0.3$

- For a given set of parameters ( $B$  and  $E$ ) we have a set of compatible initial conditions.
- Poincaré sections give us a global picture of the dynamics.
- For a better (visual) understanding of the dynamics we can show the Lyapunov exponents on the Poincaré map.

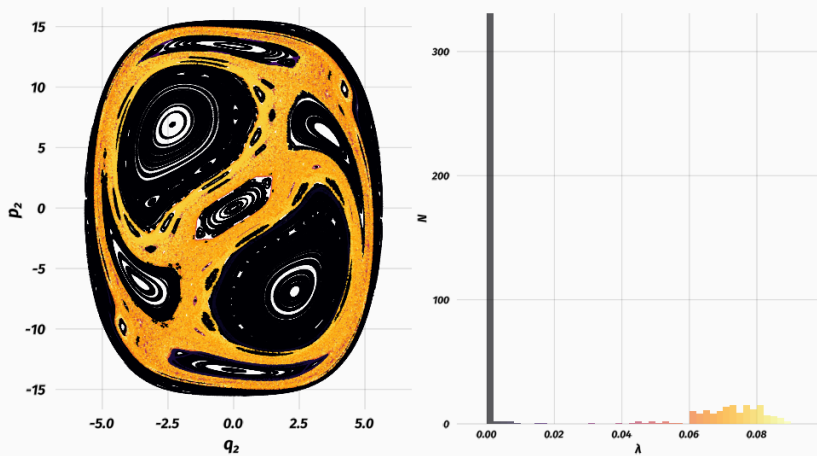
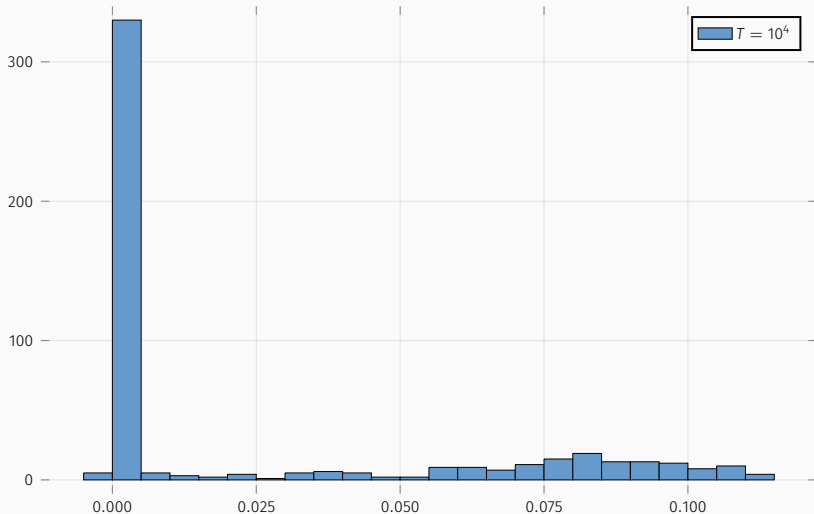


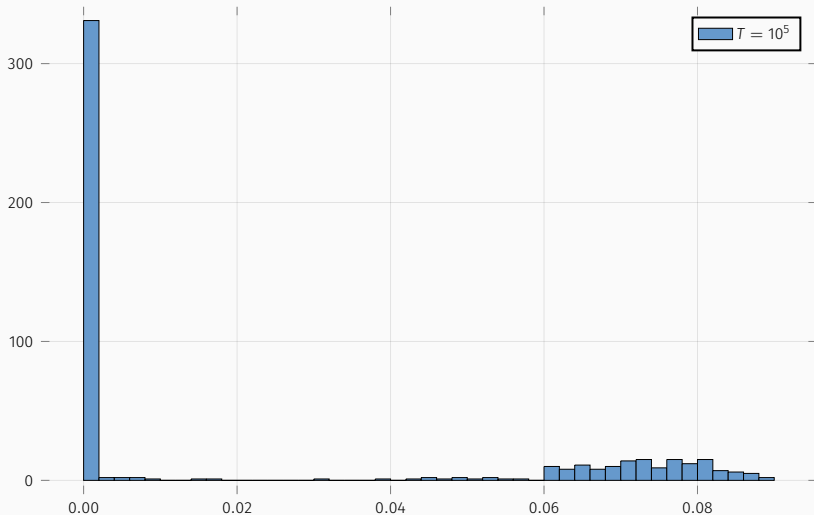
Figure 8: A Poincaré section at  $B = 0.5, E = 120$

## A note regarding the $\lambda$ histogram

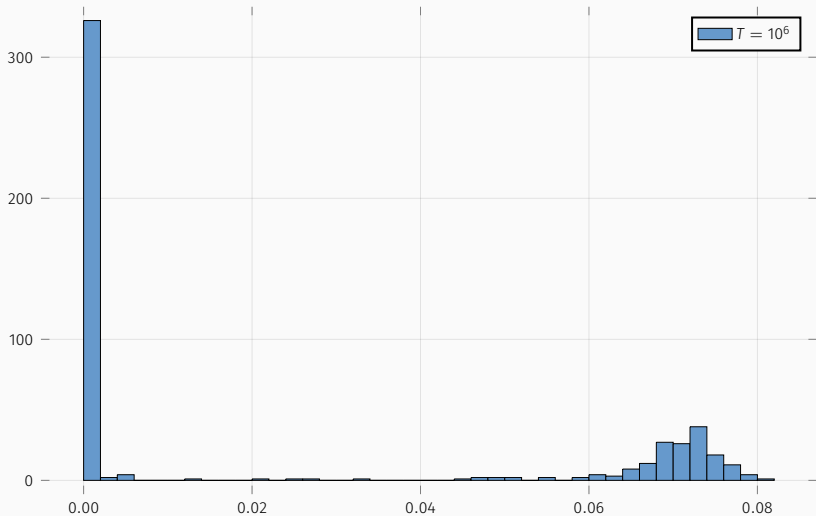
- Theoretically the Lyapunov exponent histogram should have two sharp peaks: one for the regular part and one for the chaotic one.
- The spread in the chaotic part is given by finite time effects.
- To better understand this we will take a look at how the integration time affects the results.



**Figure 9:** Maximal Lyapunov coefficient histogram for  $B = 0.5, E = 120$ .

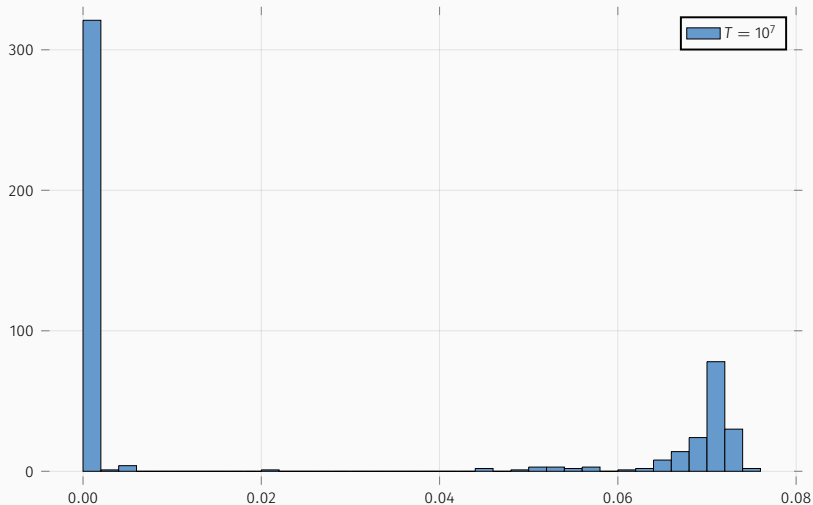


**Figure 10:** Maximal Lyapunov coefficient histogram for  $B = 0.5, E = 120$ .



**Figure 11:** Maximal Lyapunov coefficient histogram for  $B = 0.5, E = 120$ .

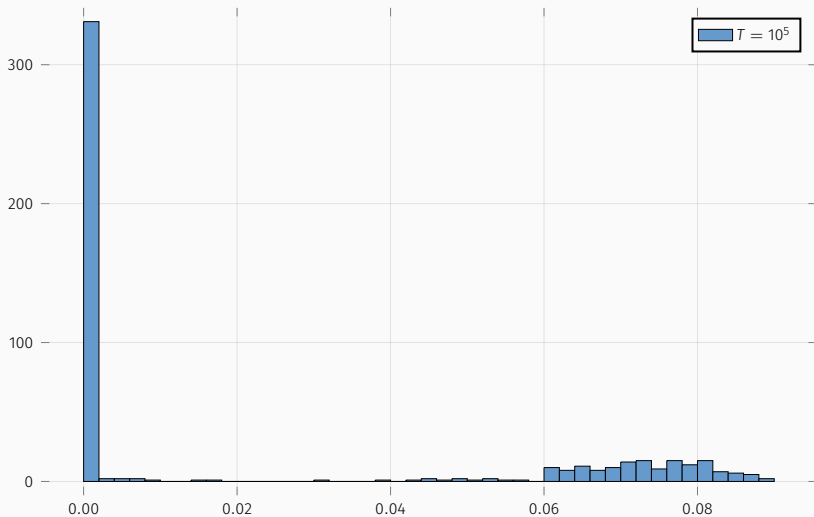




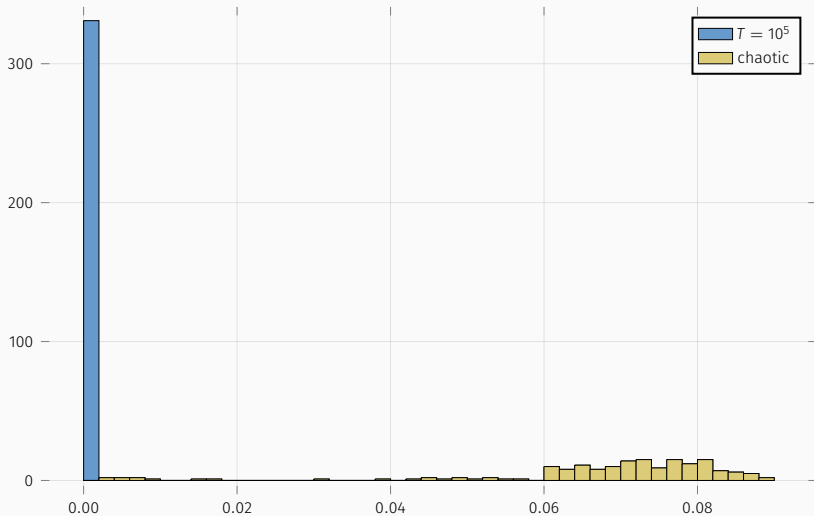
**Figure 12:** Maximal Lyapunov coefficient histogram for  $B = 0.5, E = 120$ .

## Averaging $\lambda$ over the initial conditions

- We define the averaged Lyapunov coefficient as the mean of the maximal Lyapunov exponents in the chaotic region.
- We need a sufficiently robust method of selecting the  $\lambda$ s in the chaotic region.
- We consider as chaotic everything after the first local maxima in the histogram.



**Figure 13:** Selecting the chaotic trajectories for  $B = 0.5, E = 120$ .



**Figure 14:** Selecting the chaotic trajectories for  $B = 0.5, E = 120$ .

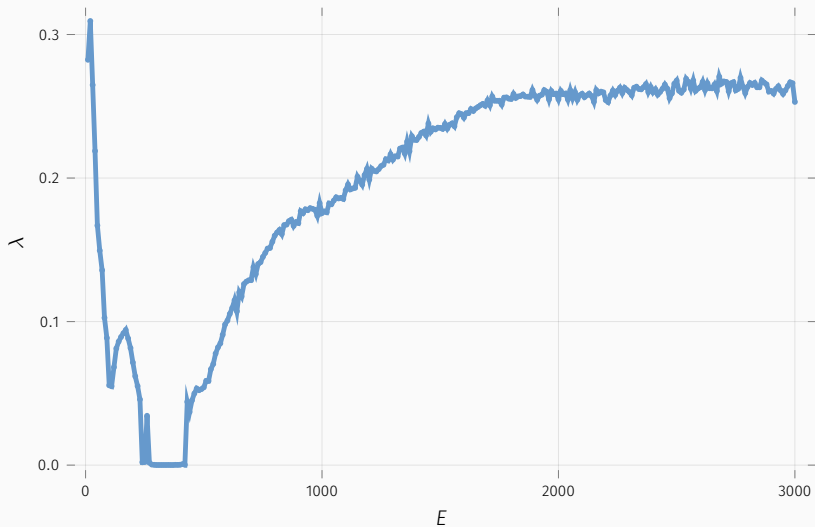
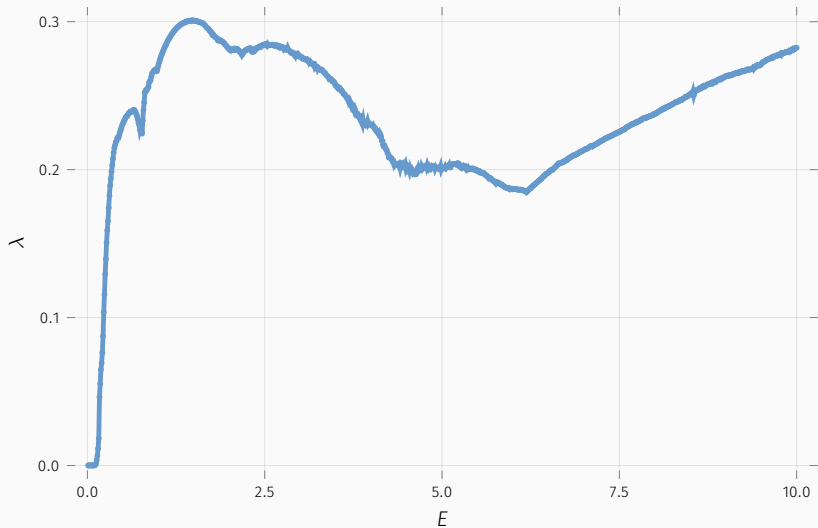


Figure 15: Averaged  $\lambda$  for  $B = 0.5, E \in (10, 3000)$ .



**Figure 16:** Averaged  $\lambda$  for  $B = 0.5, E \in (0.01, 10)$ .

## Other indicators

- We can look at the distance between two nearby trajectories in the limit of  $T \rightarrow \infty$  in order to get an estimate of the phase space volume.
- We define  $\Gamma$  as

$$\Gamma = \frac{e^\lambda - 1}{d_\infty}$$

- We can apply similar averaging techniques as for  $\lambda$ .

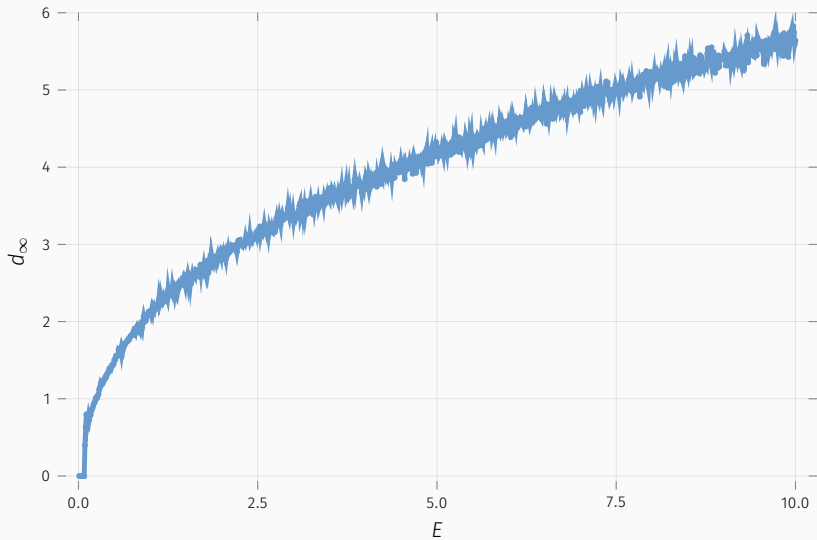
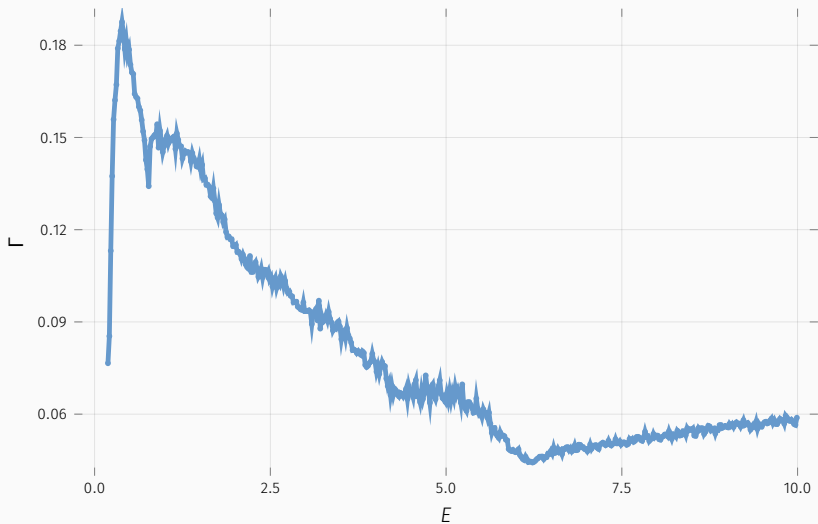


Figure 17: Averaged  $d_{\infty}$  for  $B = 0.5, E \in (0.01, 10)$ .

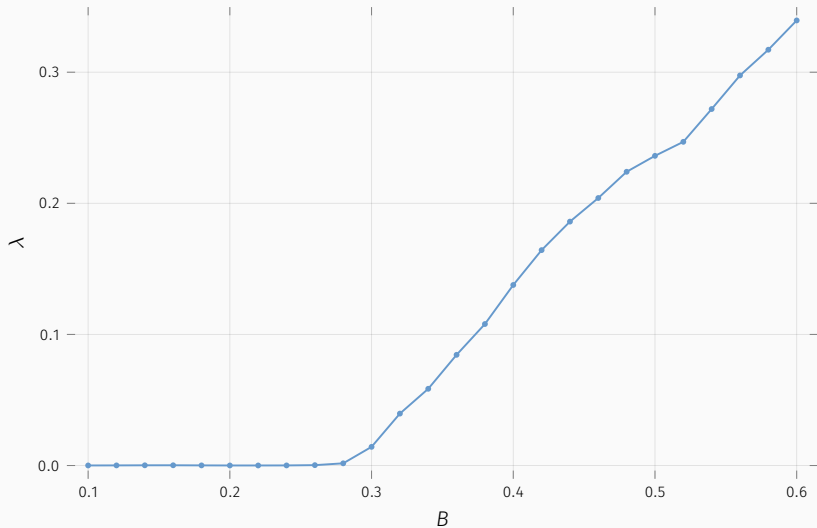




**Figure 18:** Averaged  $\Gamma$  for  $B = 0.5, E \in (0.01, 10)$ .

## Averaging $\lambda$ over the initial conditions

- We can get a global picture of the system by integrating the averaged Lyapunov coefficient over an energy interval.
- For a large energy interval we have a non-trivial dependence of the averaged Lyapunov exponent with respect to the non-integrability parameter  $B$ .



**Figure 19:** Averaged  $\lambda$  for  $B = 0.5, E \in (10, 3000)$  and  $B \in (0.1, 0.6)$ .

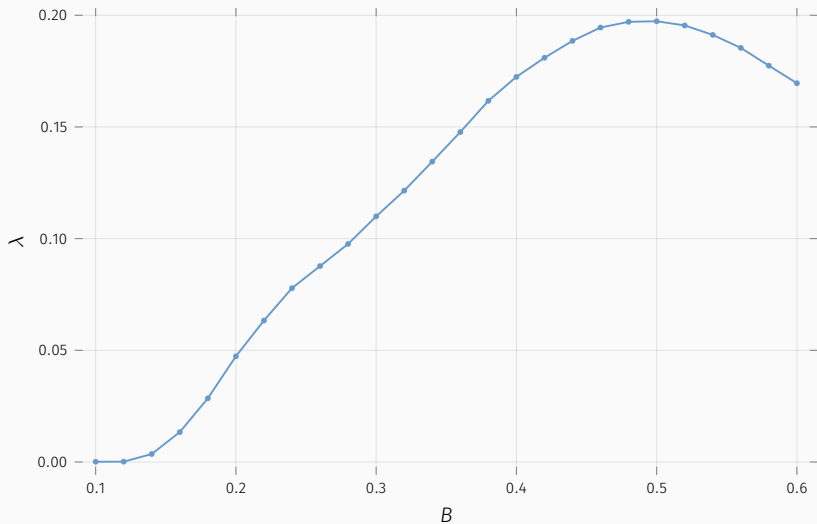


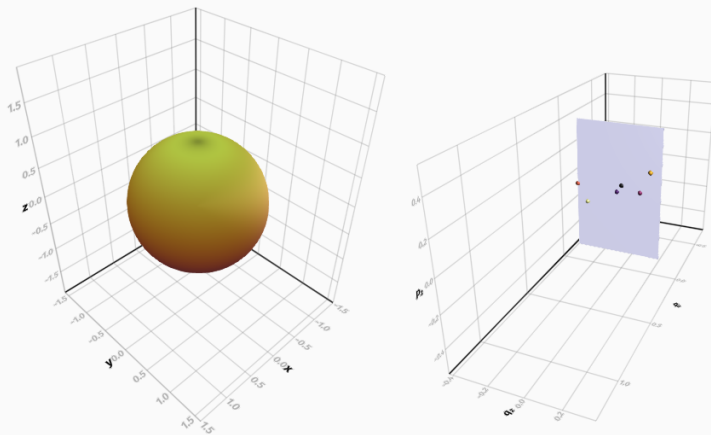
Figure 20: Averaged  $\lambda$  for  $B = 0.5, E \in (0.01, 10)$  and  $B \in (0.1, 0.6)$ .

# Conclusions

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Thank you!



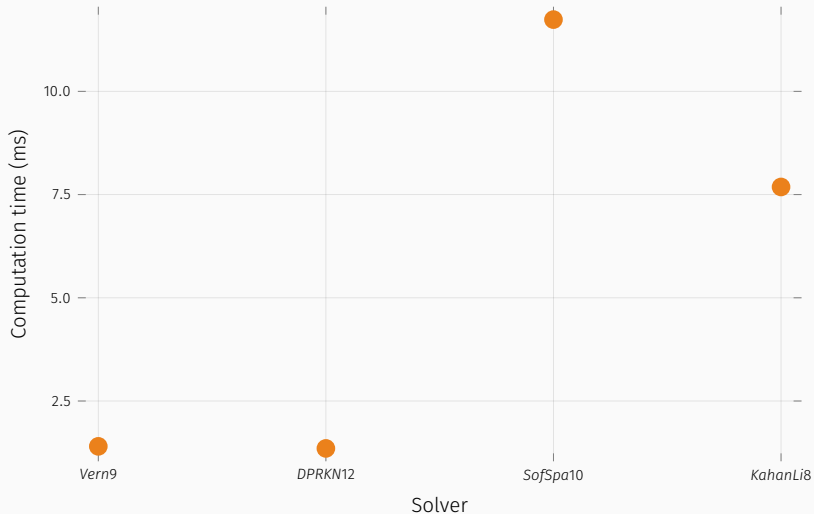


**Figure 21:** The nucleus and the corresponding trajectory in the phase space for a regular trajectory with  $B = 0.5, E = 0.3$

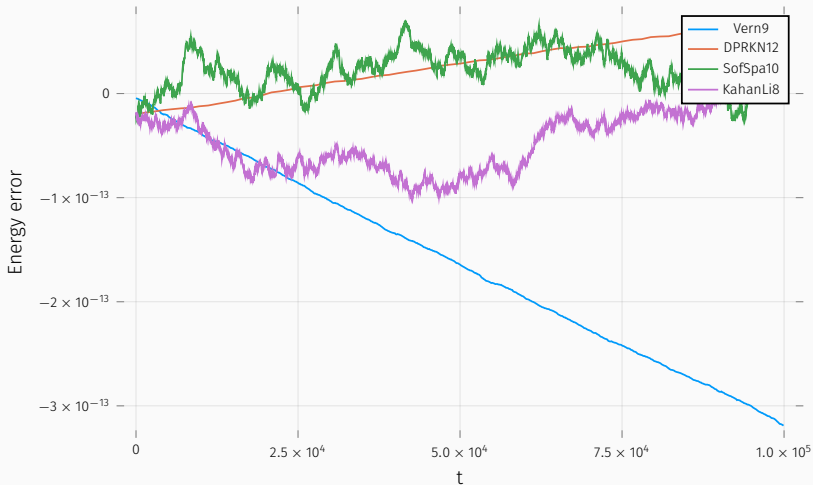




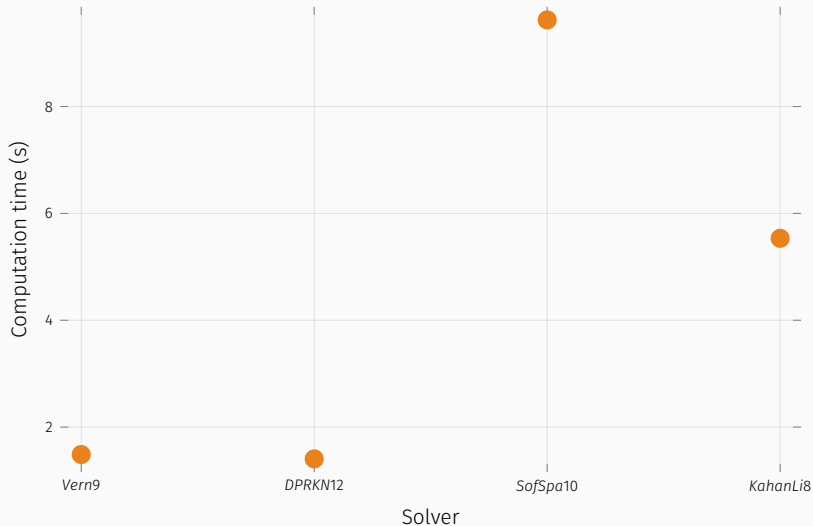
**Figure 22:** Energy error benchmark for short integration time with rescaling



**Figure 23:** Computational time benchmark for short integration time with rescaling



**Figure 24:** Energy error benchmark for long integration time with rescaling



**Figure 25:** Computational time benchmark for long integration time with rescaling