

# Lab 5: Monte Carlo

SI1136 Simulation and Modeling

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## 5.1

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## 5.1 Project

Use the Metropolis method to calculate

$$\langle x \rangle = \frac{\int_0^\infty x e^{-x} dx}{\int_0^\infty e^{-x} dx} . \quad (1)$$

Try different, also sufficiently large values of the parameter delta. Study  $\sigma/N$  and compare with the actual difference to the exact answer.<sup>1</sup>

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<sup>1</sup>Note:  $N_0$  is not critical when  $x_0 = 0$ .

## 5.1 Background

Exact value of the integral (1).

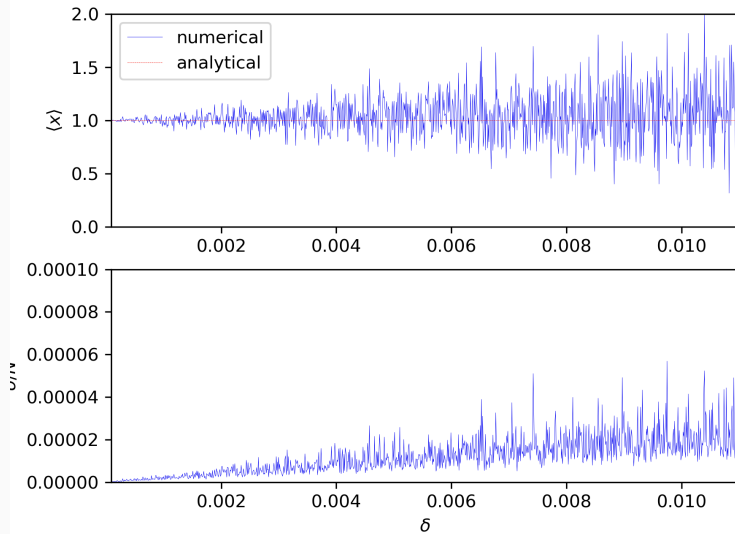
$$\langle x \rangle = \frac{\int_0^\infty x e^{-x} dx}{\int_0^\infty e^{-x} dx} = \frac{\overbrace{[-x e^{-x}]_0^\infty}^0 + \int_0^\infty e^{-x} dx}{\int_0^\infty e^{-x} dx} = 1 .$$

## 5.1 Method i

10 000 points are generated for  $\delta = 0.0001, 0.0002, \dots, 0.1$  after the first 100 points are thrown. Starting parameter  $x_0 = 1$  since this is  $x_i$  should converge to.

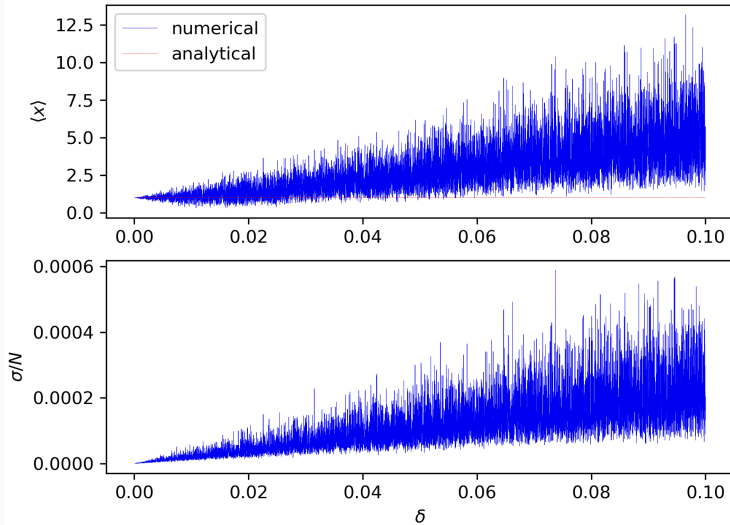
## 5.1 Results i

Calculating  $\langle x \rangle = \frac{\int_0^\infty x e^{-x} dx}{\int_0^\infty e^{-x} dx}$  using the metropolis method.



## 5.1 Results ii

Calculating  $\langle x \rangle = \frac{\int_0^\infty x e^{-x} dx}{\int_0^\infty e^{-x} dx}$  using the metropolis method.





## 5.1 Results iii

**small  $\delta$ :** For small  $\delta < 0.01$ ,  $\langle x \rangle$  varies around the true value  $\langle x \rangle = 1$ .

**large  $\delta$ :** For large  $\delta > 0.01$ ,  $\langle x \rangle$  increases linearly and fluctuates.

**$\sigma/N$ :**  $\sigma/N$  increases as  $\delta$  increases (which is expected).  $\sigma/N$  is several ( $\approx 5$ ) orders of magnitudes smaller than the actual difference to the exact answer. Since  $N = 10\,000$ ,  $\sigma \sim$  the actual difference to the exact answer.

## 5.2

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## 5.2 Project

Implement the Lennard-Jones potential with  $\sigma = 1$  and  $\varepsilon = 1$  in the template. Try to find a large MC step size that gives reasonable acceptance. How does the step size affect the convergence of the energy?

## 5.2 Background

### Lennard-Jones potential.

The Lennard-Jones potential approximates the interaction between a pair of neutral atoms or molecules to the potential

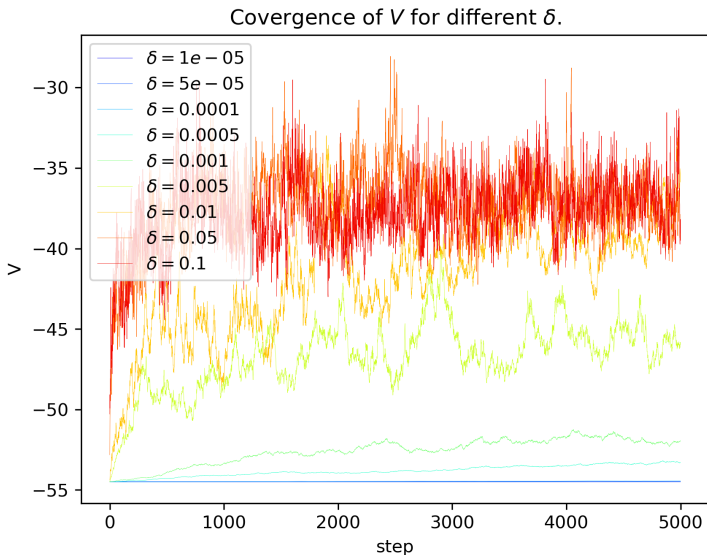
$$V_{LJ}(r) = 4\epsilon \left[ \left( \frac{\sigma}{r} \right)^{12} - \left( \frac{\sigma}{r} \right)^6 \right] .$$

## 5.2 Method

A 5 000 step Monte-Carlo simulation is executed for each  $\delta$  in

```
deltaArray = [0.00001, 0.00005, 0.0001, 0.0005,  
              0.001, 0.005, 0.01, 0.05, 0.1] .
```

## 5.2 Results i



## 5.2 Results ii

**small  $\delta$ :** Poor convergence. Small variance.

**large  $\delta$ :** Good convergence. Large variance.

**best  $\delta$ :**  $0.05 < \delta < 0.1$ .

## 5.3

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## 5.3 Project

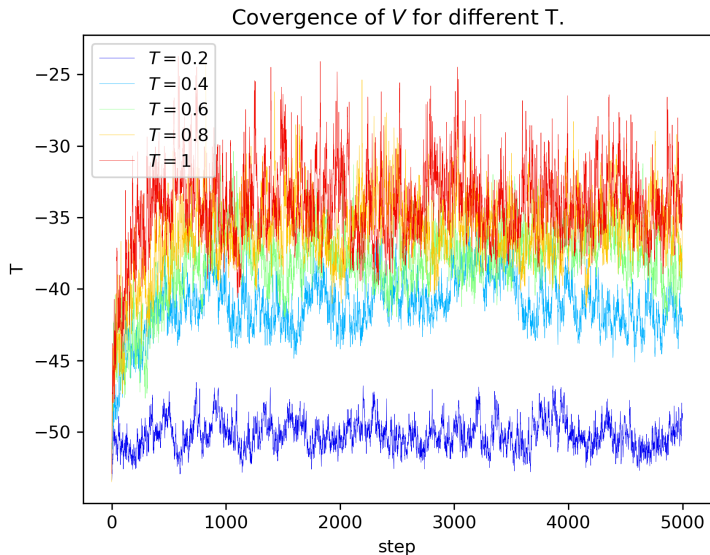
Run LJ simulations at  $T = 1$  and 0.2. What differences in collective behavior do you observe between 1 and 0.2 at long times? A physics question: can you explain what you see?

## 5.3 Method

A 5 000 step Monte-Carlo simulation is executed with  $\delta = 0.05$  for each  $T$  in

$$T_{\text{array}} = [0.2, 0.4, 0.6, 0.8, 1].$$

## 5.3 Results i



## 5.3 Results ii

**convergence:** Convergence is similar for all  $T$ .

**small  $T$ :** Small variance. Converges to small  $V$ .

**large  $T$ :** Large variance. Converges to large  $V$ .

**physical interpretation:**  $T$  is determined by the average kinetic energy of the particles. Large  $T \implies$  higher movement  $\implies$  more chaotic system.  
Furthermore, large  $T \implies$  greater energy  $\implies$  greater  $V$ .

## 5.4

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## 5.4 Project

Use MC to calculate the average energy and heat capacity at several temperatures between  $T = 0.2$  and 1. Can you explain the behavior of the heat capacity by looking at the sampled configurations?

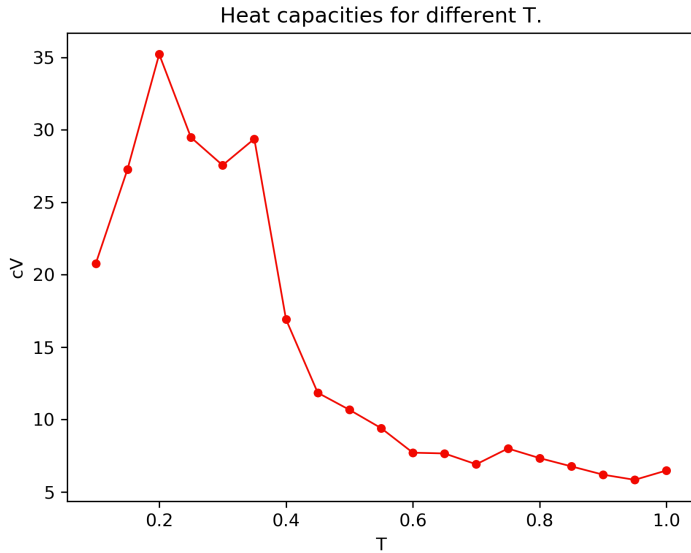
## 5.4 Method i

The heat capacity is determined for  $T$  in

$$TArray = [0.1, 0.15, 0.2, \dots, 1]$$

by averaging over  $2\,000 \leq \text{step} \leq 10\,000$ .

## 5.4 Results i





## 5.4 Results ii

The heat capacity decreases as the temperature increases.

# Appendix

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Available at:

<https://github.com/axelstr/SI1336-Simulations-and-Modeling/tree/master/1%20Monte%20Carlo>