Milestone 4

1.1 Force-Derivation from the Lenard Jones Potential

$$\overrightarrow{f_k} = \sum_{i} \frac{\partial V}{\partial r_{ik}} r_{ik} \tag{1.1}$$

$$V(r) = 4\epsilon \left[\left(\frac{\sigma}{r} \right)^{12} - \left(\frac{\sigma}{r} \right)^{6} \right]$$
 (1.2)

$$\frac{\partial V}{\partial r_{ik}} = 4\epsilon \left(\frac{6\sigma^6}{r_{ik}^6} - \frac{12\sigma^{12}}{r_{ik}^{13}} \right) \tag{1.3}$$

```
[4]: import sympy as sp import warnings warnings ('ignore') sp.init_printing() eps = sp.Symbol("e") sig = sp.Symbol("s") rad = sp.Symbol("r") energyRad = 4 * eps * ((sig/rad)**12 - (sig/rad)**6) energyRad.diff(rad)

[4]: 4e\left(\frac{6s^6}{r^7} - \frac{12s^{12}}{r^{13}}\right)
```

1.2 Different Time Steps

1.3 Simulation Snapshots

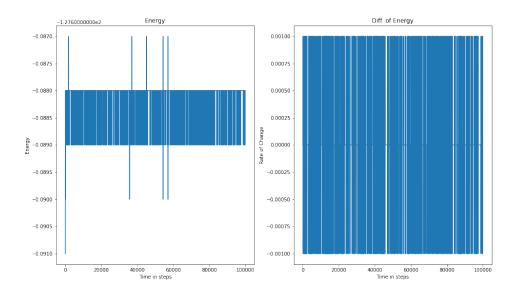


Figure 1.1: Simulation with a time step of 0.001

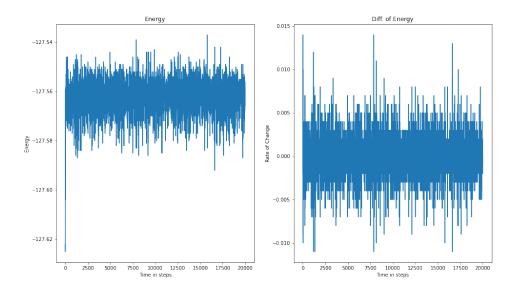


Figure 1.2: Simulation with a time step of 0.005

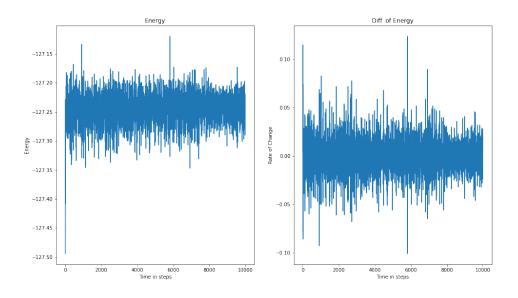


Figure 1.3: Simulation with a time step of 0.01

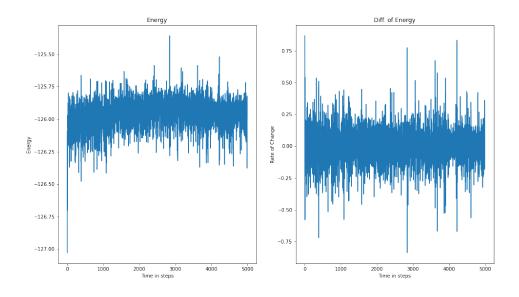


Figure 1.4: Simulation with a time step of 0.02

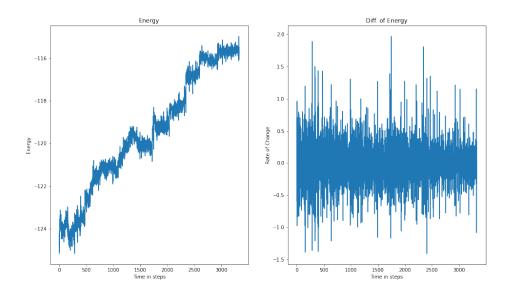


Figure 1.5: Simulation with a time step of 0.03

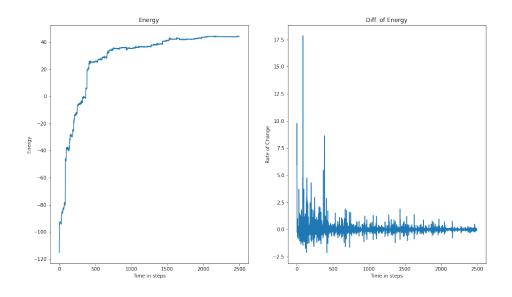


Figure 1.6: Simulation with a time step of 0.04



Figure 1.7: Simulation

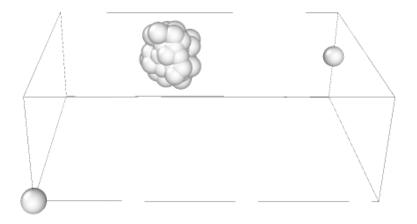


Figure 1.8: Simulation

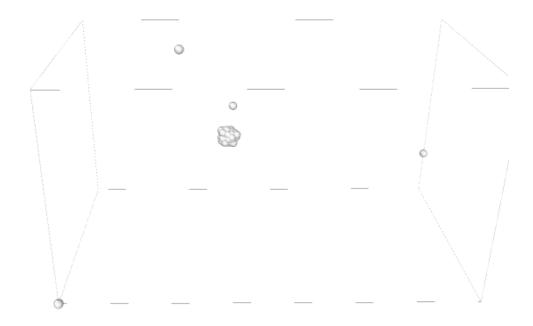


Figure 1.9: Simulation

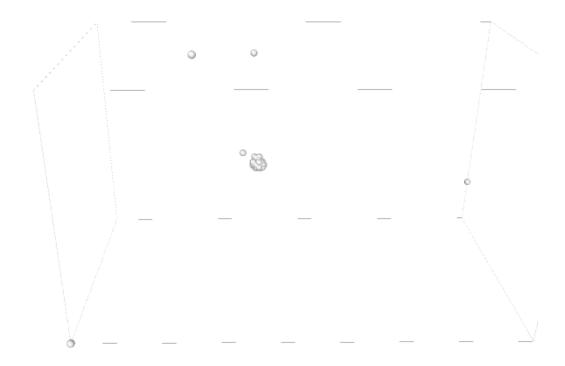


Figure 1.10: Simulation

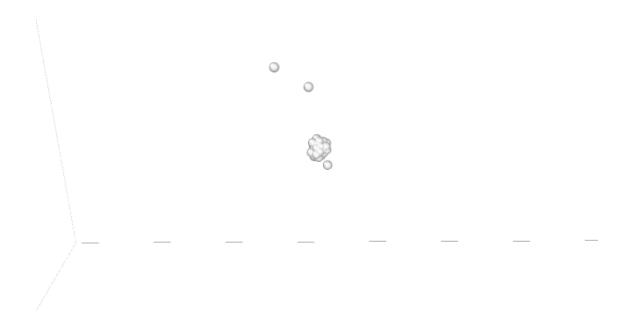


Figure 1.11: Simulation

Milestone 5

2.1 Berendsen Thermostat Teststrategy

$$\lambda = \sqrt{1 + \left(\frac{T_0}{T} - 1\right) \frac{\Delta t}{\tau}} \tag{2.1}$$

$$T(t) = T_0 + (T_1 - T_0)e^{-t/\tau}$$
(2.2)

2.2 Simulation Time

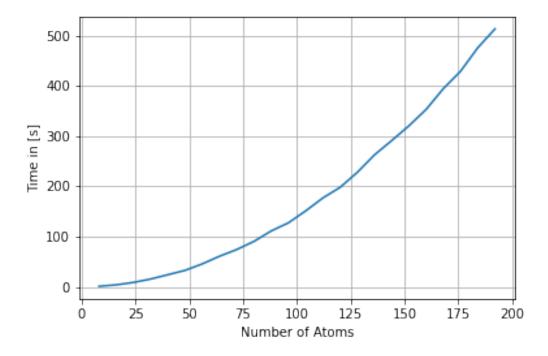


Figure 2.1: Simulation time from 8 to 192 Atoms

Milestone 6

The Images are outdated!!!!

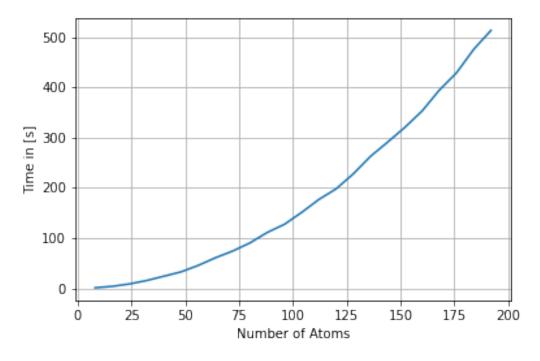


Figure 3.1: Simulationtime without Neighborlist

New images that are correct.

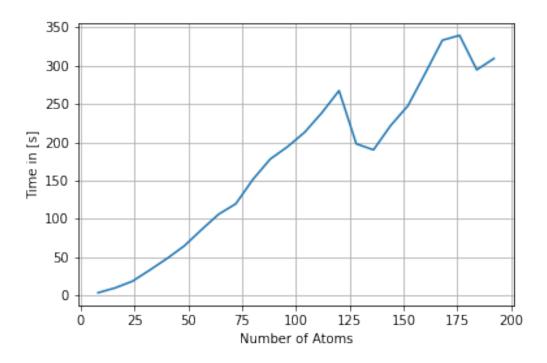


Figure 3.2: Simulationtime with the old Neighborlist

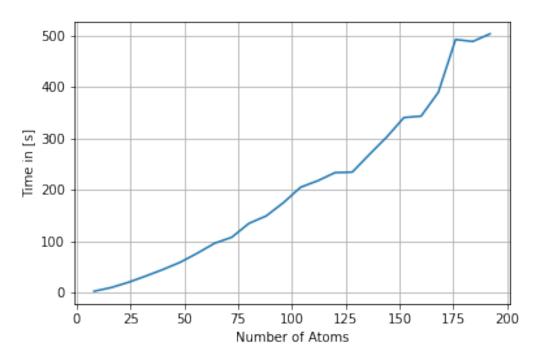


Figure 3.3: Simulation time with the new Neighborlist

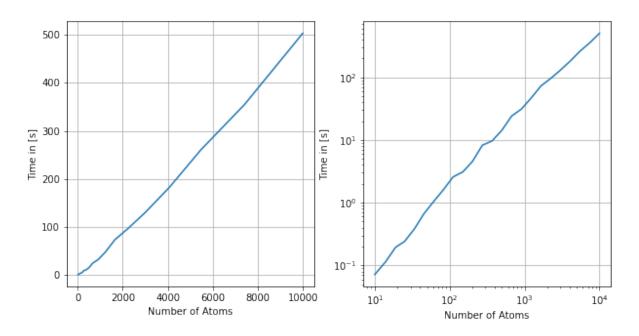


Figure 3.4: Simulationtime with the new Neighborlist

Milestone 7

