

# Markov chain Monte Carlo Simulation

## An Example with Argon

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# Outline

| 1

- 1 Introduction
- 2 Simulation system setup
- 3 Results and conclusion

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| 2

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- For equilibrium distribution, the average of a physical quantity  $A$  in equilibrium is

$$\langle A \rangle = \int d\mathbf{x} \rho(\mathbf{x}) A(\mathbf{x})$$

Here,  $\rho(\mathbf{x}) = e^{-\beta H(\mathbf{x})}/Z$  is the canonical ensemble distribution function, and  $Z = \int d\mathbf{x} e^{-\beta H(\mathbf{x})}$  is the canonical partition function.

- When equilibrium, the probability of the system being in configuration  $\mathbf{x}_a$  do not change with time. Therefore, we have

$$\frac{dP_a}{dt} = \sum_b [w_{b \rightarrow a} P_b - w_{a \rightarrow b} P_a] = 0$$

Here,  $w_{b \rightarrow a}$  represents the transition probability from configuration  $\mathbf{x}_b$  to  $\mathbf{x}_a$ .

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Here,  $w_{b \rightarrow a}$  represents the transition probability from configuration  $\mathbf{x}_b$  to  $\mathbf{x}_a$ .

- ▶ From the above relation, we can deduce that if **detailed balance** is satisfied, we can generate a sequence that follows the Boltzmann distribution:

$$\frac{w_{a \rightarrow b}}{w_{b \rightarrow a}} = \frac{P_b}{P_a} = e^{-\beta(E_b - E_a)} = e^{\beta \Delta E_{ab}}$$

- ▶ Metropolis transition probability:

$$w_{a \rightarrow b} = \begin{cases} e^{\beta \Delta E_{ab}} & (E_b > E_a) \\ 1 & (E_b \leq E_a) \end{cases}$$

- ▶ Metropolis–Hastings algorithm:

1. Initial configuration  $\mathbf{x}_i$  with energy  $E_i$ ;
2. Random move to a new configuration  $\mathbf{x}_{i'}$  with energy  $E_{i'}$ ;
3. If  $E_{i'} \leq E_i$  or  $e^{\beta(E_{i'} - E_i)} < R$ ,  $R$  is random number in  $[0, 1)$ , accept the move. The next configuration will be  $\mathbf{x}_{i+1} = \mathbf{x}_{i'}$ ;
4. Return to step 2 and repeat until  $N$  configurations have been accumulated.

- ▶ Starting from any initial configuration, the system will eventually reach equilibrium and follow the Boltzmann distribution.

# Outline

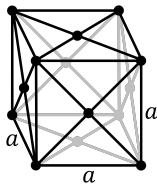
| 6

- 1 Introduction
- 2 Simulation system setup
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# Argon

| 7

- ▶ Argon is a chemical element; it has symbol Ar and atomic number 18. It is in group 18 of the periodic table and is a noble gas.
- ▶ Melting point: 83.81 K, boiling point: 87.302 K, triple point: 83.8058 K
- ▶ Liquid density (at b.p.):  $0.02103 \text{ Å}^{-3}$
- ▶ Lattice: FCC ( $a=5.4691 \text{ Å}$ ), density:  $0.02445 \text{ Å}^{-3}$  (at triple point)





- ▶ Mie potential is an interaction potential describing the interactions between particles on the atomic level. It is mostly used for describing intermolecular interactions, but at times also for modeling intramolecular interaction.

$$U(r) = C\varepsilon \left[ \left( \frac{\sigma}{r} \right)^n - \left( \frac{\sigma}{r} \right)^m \right]$$

with

$$C = \frac{n}{n-m} \left( \frac{n}{m} \right)^{\frac{m}{n-m}}$$

- ▶ The Lennard-Jones potential corresponds to the case where  $n=12$  and  $m=6$ .
- ▶ For Ar,  $\sigma=3.404 \text{ \AA}$ ,  $\varepsilon=117.84 k_B$ ,  $n=12.085$ ,  $m=6.0$ .

# Outline

| 9

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- ▶ Pair correlation function  $g(r)$  in a system of particles describes how density varies as a function of distance from a reference particle.

$$g(\mathbf{r}_1, \mathbf{r}_2) = \frac{\rho(\mathbf{r}_1, \mathbf{r}_2)}{\rho(\mathbf{r})^2}$$

In fluid,  $\rho(\mathbf{r}) = \rho$ ,  $g(\mathbf{r}_1, \mathbf{r}_2) = g(|\mathbf{r}_1 - \mathbf{r}_2|) \equiv g(r)$ .  $g(r)$  can also be referred to as the radial distribution function. When  $r \rightarrow +\infty$ ,  $g(r) \rightarrow 1$ .

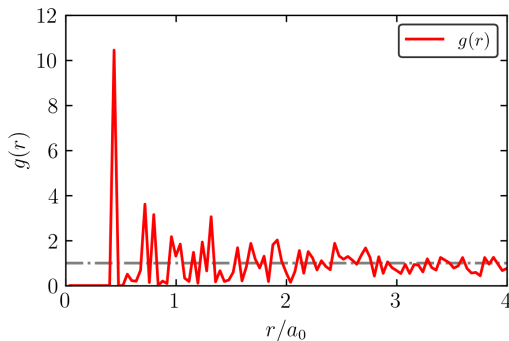
- ▶ In practice, we choose a series of radius grids  $\{r_i\}$ , which have the same difference  $dr$ , and the pair correlation function can be calculated using:

$$g(r_i) = \frac{1}{\rho} \frac{n(r_i)}{4\pi r^2 dr}$$

## Liquid case

| 11

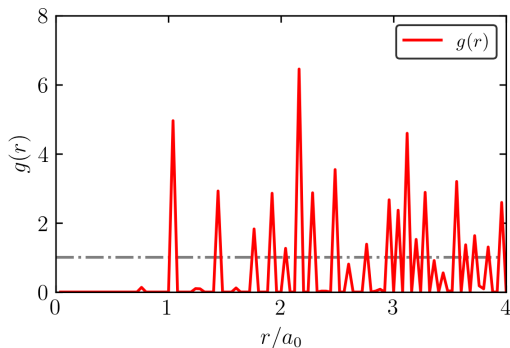
- Density :  $0.02103 \text{ \AA}^{-3}$ , Temperature: 84K
- nmol: 108, box:  $3 \times 3 \times 3$  unit cell,  $a_0$  is box size, ncycle: 1000, ngrid: 100.



## Solid case

| 12

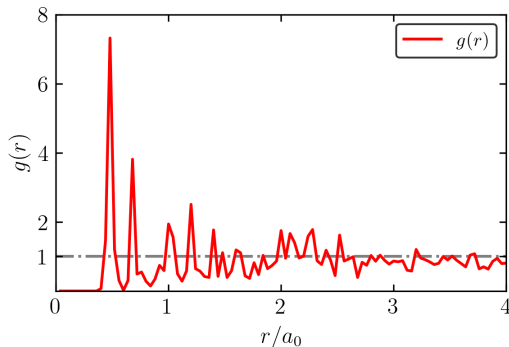
- Density :  $0.2103 \text{ \AA}^{-3}$ , Temperature: 10K
- nmol: 108, box:  $3 \times 3 \times 3$  unit cell,  $a_0$  is box size, ncycle: 1000, ngrid: 100.



## Why higher density

| 13

- Density :  $0.02445 \text{ \AA}^{-3}$ , Temperature: 10K
- nmol: 108, box:  $3 \times 3 \times 3$  unit cell,  $a_0$  is box size, ncycle: 1000, ngrid: 100.



- ▶ Mie potential is used to describe the potential of a fluid, and is not applicable for solids. In this case, only by applying sufficiently large pressure to the system can some solid-like properties be exhibited.
- ▶ The radius of the first peak in the three graphs, when converted to Å, is approximately the same (1.1501 Å, 1.0677 Å, 1.0938 Å), which may suggest the radius of argon ( $1.06 \pm 0.1$  Å).