

Markov-chain Monte Carlo: A modern primer

Lecture 2: Surprises

Part 2.1: Reversible and non-reversible hard-sphere algorithms

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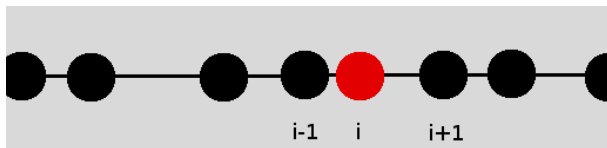
A Set of Lectures
University of Kent
Canterbury, Great Britain

14-17 November 2022

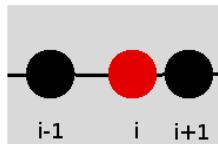
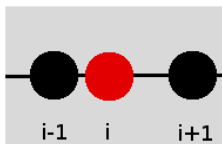
References

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Local Metropolis algorithm 1/5



- Spheres on a $1d$ interval (with or without pbc).
- $\pi(x_1, \dots, x_N)$ non-trivial (see Krauth (2006)).
- Local Metropolis: $x_j \rightarrow x_j \pm \epsilon$ (with pbc).
- $\Omega^{\text{Met}} = \{x_1 < x_2 < \dots < x_N\}$, $\mathbf{x} = \{x_1, \dots, x_N\}$.
- $\mathcal{L} = \mathbb{R} \times \{1, \dots, N\}$ (or some symmetric interval)
- Local Metropolis: Sample element of \mathcal{L} , apply to \mathbf{x}_t

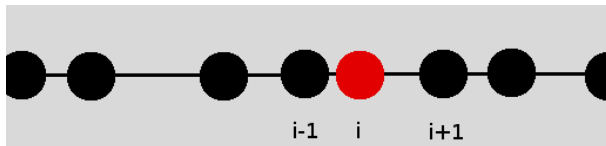


- $\Omega^{\text{Met}} = \{x_1 < x_2 < \dots < x_N\}$, $\mathbf{x} = \{x_1, \dots, x_N\}$.
- $\mathcal{L}^{\text{Met}} = \mathbb{R} \times \{1, \dots, N\}$ (or some symmetric interval)
- Local Metropolis: Sample element of \mathcal{L} , apply to \mathbf{x}_t
- Detailed balance:

$$\pi_a p(a \rightarrow b) = \pi_b p(b \rightarrow a)$$

- Probability distribution on \mathbb{R} symmetric
- Almost no condition on sampling of $\mathcal{N} = \{1, \dots, N\}$.

Local Metropolis algorithm 3/5

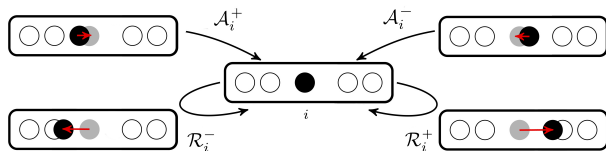


- Reversible Metropolis: $x_i \rightarrow x_i \pm \epsilon$ (reject if overlap)
- Global balance:

$$\mathcal{F}_a^{\text{Met}} = \frac{1}{2N} \sum_i \int_0^\infty d\epsilon \underbrace{(\mathcal{A}_i^+ + \mathcal{R}_i^+ + \mathcal{A}_i^- + \mathcal{R}_i^-)}_{= 2 \text{ for any } \epsilon} = 1.$$

- NB: $\mathcal{A}_i^+(\epsilon) + \mathcal{R}_i^-(\epsilon) = 1$ also $\mathcal{A}_i^-(\epsilon) + \mathcal{R}_i^+(\epsilon) = 1$.

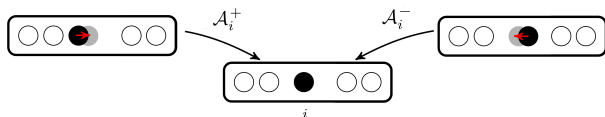
Local Metropolis algorithm 4/5



- Reversible Metropolis: $x_i \rightarrow x_i \pm \epsilon$ (reject if overlap)
- Global balance:

$$\mathcal{F}_a^{\text{rev}} = \frac{1}{2N} \sum_i \int_0^\infty d\epsilon \underbrace{(\mathcal{A}_i^+ + \mathcal{R}_i^+ + \mathcal{A}_i^- + \mathcal{R}_i^-)}_{= 2 \text{ for any } \epsilon} = 1.$$

- NB: $\mathcal{A}_i^+(\epsilon) + \mathcal{R}_i^-(\epsilon) = 1$; $\mathcal{A}_i^-(\epsilon) + \mathcal{R}_i^+(\epsilon) = 1$.
- Good exercise.



Variant:

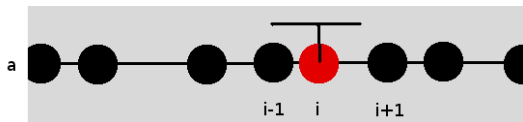
- Heat bath: $x_j \rightarrow \text{ran}(x_{j-1} + 2\sigma, x_{j+1} - 2\sigma)$
- Mixing time $\mathcal{O}(N^3 \log N)$ with fixed boundary conditions (Randall, Winkler 2005).
- Mixing time $\mathcal{O}(N^3) \dots \mathcal{O}(N^3 \log N)$ with periodic boundary conditions (Randall, Winkler 2005).
- Numerically: $\mathcal{O}(N^3 \log N)$ for Metropolis and heat bath.

Symmetric simple exclusion process 1/2

- $\Omega^{\text{Met}} = \{x_1 < x_2 < \dots < x_N\}$, $\mathbf{x} = \{x_1, \dots, x_N\}$.
- $\mathcal{L}^{\text{Met}} = \mathbb{R} \times \{1, \dots, N\}$ (or some symmetric interval)
- $\Omega^{\text{SSEP}} = \{x_1 < x_2 < \dots < x_N\}$, $\mathbf{x} = \{x_1, \dots, x_N\} \in \mathbb{N}^N$.
- $\mathcal{L}^{\text{SSEP}} = \{-, +\} \times \{1, \dots, N\}$

- Mixing time of SSEP $\mathcal{O}(N^3 \log N)$ (Lacoin)
- Many studies

Sequential Metropolis algorithm

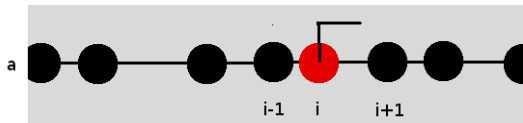


- $\Omega^{\text{Met}} = \{x_1 < x_2 < \dots < x_N\}$, $\mathbf{x} = \{x_1, \dots, x_N\}$.
- $\mathcal{L}^{\text{Met}} = \mathbb{R} \times \{1, \dots, N\}$
- $\hat{\Omega}^{\text{Seq}} = \Omega^{\text{Met}} \times \mathcal{N}$ (particle lifting).
- Sequential Metropolis: Update 0, then 1, then 2, ...
- Global balance:

$$\mathcal{F}_a^{\text{seq}} = \frac{1}{2} \int_0^\infty d\epsilon (\mathcal{A}_i^+ + \mathcal{R}_i^+ + \mathcal{A}_i^- + \mathcal{R}_i^-) = 1.$$

- Many sequences of indices are OK.
- Thm: Any sequential version of a reversible Markov chain satisfies global balance (conditions apply).

Forward Metropolis algorithm 1/2

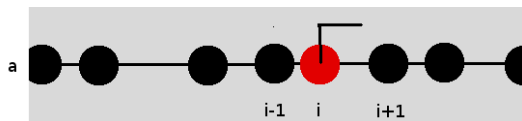


- $\Omega^{\text{Met}} = \{x_1 < x_2 < \dots < x_N\}$, $\mathbf{x} = \{x_1, \dots, x_N\}$.
- $\mathcal{L}^{\text{Met}} = \mathbb{R} \times \mathcal{N} = \mathbb{R}^+ \times \{+, -\} \times \mathcal{N}$
- $\hat{\Omega}^{\text{Forw}} = \Omega^{\text{Met}} \times \{+, -\}$ (displacement lifting).
- $\mathcal{L}^{\text{Forw}} = \mathbb{R}^+ \times \mathcal{N}$
- Forward Metropolis: $x_i \rightarrow x_i + \epsilon$ (NB: random i)
- Sample-space halving...
-

$$\mathcal{F}_a^{\text{forw}} = \frac{1}{N} \sum_i \underbrace{(\mathcal{A}_i^+ + \mathcal{R}_{i-1}^+)}_{=1 \text{ for any } \epsilon} = 1,$$

- Lattice version: Totally asymmetric simple exclusion process (TASEP).
- NB: Forward sequential Metropolis is wrong.

Forward Metropolis algorithm 1/2



- $\Omega^{\text{Met}} = \{x_1 < x_2 < \dots < x_N\}$, $\mathbf{x} = \{x_1, \dots, x_N\}$.
- $\mathcal{L}^{\text{Met}} = \mathbb{R} \times \mathcal{N} = \mathbb{R}^+ \times \{+, -\} \times \mathcal{N}$
- $\hat{\Omega}^{\text{Forw}} = \Omega \times \{+, -\}$ (displacement lifting).
- Forward Metropolis: $x_i \rightarrow x_i + \epsilon$ (NB: random i)
- Sample-space halving...

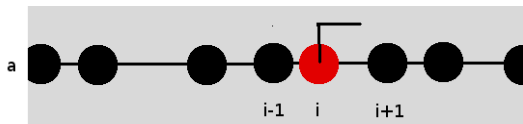
$$\mathcal{F}_a^{\text{forw}} = \frac{1}{N} \sum_i \underbrace{(\mathcal{A}_i^+ + \mathcal{R}_{i-1}^+)}_{=1 \text{ for any } \epsilon} = 1,$$

- NB: Forward sequential Metropolis is wrong (see TD).

Totally asymmetric simple exclusion process

- $\Omega^{\text{SSEP}} = \{x_1 < x_2 < \dots < x_N\} \in \mathbb{N}^N$
- $\mathcal{L}^{\text{SSEP}} = \{+, -\} \times \mathcal{N}$
- $\hat{\Omega}^{\text{TASEP}} = \Omega^{\text{SSEP}} \times \{+, (-)\}$ (displacement lifting).
- $\mathcal{L}^{\text{TASEP}} = \mathcal{N}$
- Sample-space halving...
- TASEP mixing times are known analytically.

Lifted Forward Metropolis algorithm



- $\Omega^{\text{Met}} = \{x_1 < x_2 < \dots < x_N\}$, $\mathbf{x} = \{x_1, \dots, x_N\}$.
- $\mathcal{L}^{\text{Met}} = \mathbb{R} \times \mathcal{N} = \mathbb{R}^+ \times \{+, -\} \times \mathcal{N}$
- $\hat{\Omega}^{\text{Lift-Forw}} = \Omega \times \{+, -\} \times \mathcal{N}$ (displacement + particle lifting).
- $\mathcal{L}^{\text{Lift-Forw}} = \mathbb{R}^+$
- Move i forward if accepted
- Otherwise: lifting move $i \rightarrow i + 1$
- Sample-space halving again.
-

$$\mathcal{F}_{(a,i)}^{\text{lift}} = \mathcal{A}_i^+ + \mathcal{R}_{i-1}^+ = 1.$$

- Analogous to the lifted Metropolis algorithm on the path graph.

Mixing time (operational) 1/2

- Half-system distance (with $\delta_i = x_i - x_{i-1} - d$ with d diameter, L system size, $L_{\text{free}} = L - Nd$):

$$u = \underbrace{\delta_{i+N/2} + \dots + \delta_{i+1}}_{N/2 \text{ terms}}.$$

- Equilibrium distribution of half-system distance:

$$\pi(u) = \frac{\Gamma(N-1)}{L_{\text{free}}^{N-1}} \frac{(L_{\text{free}}u - u^2)^{N/2-1}}{(N/2-1)!^2}.$$

- Variance of half-system distance:

$$\text{Var } u = \frac{N^2 \ell_{\text{free}}^2}{4(N+1)}.$$

- Compact initial condition:

$$\text{Var}_{\text{compact}}(u) = \langle u^2 \rangle_{\text{compact}} - \langle u \rangle_{\text{compact}}^2 = \frac{N^2 \ell_{\text{free}}^2}{4}.$$

Mixing time (operational) 2/2

