Optimization and Simulation Markov Chain Monte Carlo Methods

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Outline

Motivation

- Introduction to Markov chains
- 3 Stationary distributions
- 4 Metropolis-Hastings
- 6 Gibbs sampling
- 6 Simulated annealing

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The knapsack problem

- Patricia prepares a hike in the mountain.
- She has a knapsack with capacity Wkg.
- She considers carrying a list of *n* items.
- Each item has a utility u_i and a weight w_i .
- What items should she take to maximize the total utility, while fitting in the knapsack?



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Knapsack problem



Simulation

- Let \mathcal{X} be the set of all possible configurations (2^n) .
- Define a probability distribution:

$$P(x) = \frac{U(x)}{\sum_{y \in \mathcal{X}} U(y)}$$

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• Question: how to draw from this discrete random variable?

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- 3 Stationary distributions
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- 6 Gibbs sampling
- Simulated annealing

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Markov Chains



Andrey Markov, 1856–1922, Russian mathematician.

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Markov Chains: glossary

Stochastic process

 X_t , t = 0, 1, ..., i, collection of r.v. with same support, or states space $\{1, ..., i, ..., J\}$.

Markov process: (short memory)

$$\Pr(X_t = i | X_0, \dots, X_{t-1}) = \Pr(X_t = i | X_{t-1})$$

Homogeneous Markov process

$$\Pr(X_t = j | X_{t-1} = i) = \Pr(X_{t+k} = j | X_{t-1+k} = i) = P_{ij} \ \forall t \ge 1, k \ge 0.$$

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Markov Chains

Transition matrix

$$P \in \mathbb{R}^{J \times J}$$
.

Properties:

$$\sum_{j=1}^{J} P_{ij} = 1, \ i = 1, \dots, J, \ P_{ij} \ge 0, \ \forall i, j,$$

Ergodicity

- If state *j* can be reached from state *i* with non zero probability, and *i* from *j*, we say that *i* communicates with *j*.
- Two states that communicate belong to the same *class*.
- A Markov chain is *irreducible* or *ergodic* if it contains only one class.
- With an ergodic chain, it is possible to go to every state from any state.

Markov Chains

Aperiodic

- P_{ij}^t is the probability that the process reaches state *j* from *i* after *t* steps.
- Consider all t such that $P_{ii}^t > 0$. The largest common divisor d is called the *period* of state i.
- A state with period 1 is *aperiodic*.
- If $P_{ii} > 0$, state *i* is aperiodic.
- The period is the same for all states in the same class.
- Therefore, if the chain is irreducible, if one state is aperiodic, they all are.

A periodic chain



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Another periodic chain



An aperiodic chain



Aperiodic chain

An equivalent definition

An irreducible Markov chain is said to be aperiodic if for some $t \ge 0$ and some state *i*, we have

$$\Pr(X_t = i | X_0 = i) > 0$$

and

$$\Pr(X_{t+1} = i | X_0 = i) > 0$$

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Intuition

Oscillation

$$P = \left(\begin{array}{cc} 0 & 1 \\ 1 & 0 \end{array}\right)$$

The chain will not "converge" to something stable.

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Markov Chains

Stationary probabilities

$$\Pr(j) = \sum_{i=1}^{J} \Pr(j|i) \Pr(i)$$

• Stationary probabilities: unique solution of the system

$$\pi_{j} = \sum_{i=1}^{J} P_{ij}\pi_{i}, \quad \forall j = 1, \dots, J.$$

$$\sum_{j=1}^{J} \pi_{j} = 1.$$
(1)

• Solution exists for any irreducible chain.

Example

- A machine can be in 4 states with respect to wear
 - perfect condition,
 - partially damaged,
 - seriously damaged,
 - completely useless.
- The degradation process can be modeled by an irreducible aperiodic homogeneous Markov process, with the following transition matrix:

$$P = \left(egin{array}{ccccc} 0.95 & 0.04 & 0.01 & 0.0 \ 0.0 & 0.90 & 0.05 & 0.05 \ 0.0 & 0.0 & 0.80 & 0.20 \ 1.0 & 0.0 & 0.0 & 0.0 \end{array}
ight)$$

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Example

Stationary distribution: $\left(\frac{5}{8}, \frac{1}{4}, \frac{3}{32}, \frac{1}{32}\right)$

$$\left(\frac{5}{8}, \frac{1}{4}, \frac{3}{32}, \frac{1}{32}\right) \left(\begin{array}{cccc} 0.95 & 0.04 & 0.01 & 0.0\\ 0.0 & 0.90 & 0.05 & 0.05\\ 0.0 & 0.0 & 0.80 & 0.20\\ 1.0 & 0.0 & 0.0 & 0.0 \end{array}\right) = \left(\frac{5}{8}, \frac{1}{4}, \frac{3}{32}, \frac{1}{32}\right)$$

- Machine in perfect condition 5 days out of 8, in average.
- Repair occurs in average every 32 days

From now on: Markov process = irreducible aperiodic homogeneous Markov process

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Markov Chains

Detailed balance equations

Consider the following system of equations:

$$x_i P_{ij} = x_j P_{ji}, \quad i \neq j, \quad \sum_{i=1}^J x_i = 1$$
 (2)

We sum over *i*:

$$\sum_{i=1}^{J} x_i P_{ij} = x_j \sum_{i=1}^{J} P_{ji} = x_j.$$

If (2) has a solution, it is also a solution of (1). As π is the unique solution of (1) then $x = \pi$.

$$\pi_i P_{ij} = \pi_j P_{ji}, \quad i \neq j$$

The chain is said time reversible

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Stationary distributions

Property of irreducible aperiodic Marlov chains

$$\pi_j = \lim_{t\to\infty} \Pr(X_t = j) \ j = 1, \dots, J.$$

Ergodicity

- Let f be any function on the state space.
- Then, with probability 1,

$$\lim_{T\to\infty}\frac{1}{T}\sum_{t=1}^T f(X_t) = \sum_{j=1}^J \pi_j f(j).$$

• Computing the expectation of a function of the stationary states is the same as to take the average of the values along a trajectory of the process.

Example: T = 100



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Example: T = 1000



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Example: T = 10000



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A periodic example

It does not work for periodic chains

$$P = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}$$
$$\Pr(X_t = 1) = \begin{cases} 1 & \text{if } t \text{ is odd} \\ 0 & \text{if } t \text{ is even} \end{cases}$$
$$\lim_{t \to \infty} \Pr(X_t = 1) \text{ does not exist}$$

Stationary distribution

$$\pi = \left(\begin{array}{c} 0.5\\ 0.5 \end{array}\right)$$

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Simulation

Motivation

- Sample from very large discrete sets (e.g. the knapsack).
- Full enumeration of the set is infeasible.

Procedure

• We want to simulate a r.v. X with pmf

$$\Pr(X=j)=p_j.$$

- We generate a Markov process with limiting probabilities p_j (how?)
- We simulate the evolution of the process.

$$p_j = \pi_j = \lim_{t \to \infty} \Pr(X_t = j) \ j = 1, \dots, J.$$

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Simulation

Assume that we are interested in simulating

$$\mathsf{E}[f(X)] = \sum_{j=1}^{J} f(j) p_j.$$

We use ergodicity to estimate it with

$$\frac{1}{T}\sum_{t=1}^{T}f(X_t).$$

Drop early states (see above example)

Better estimate:

$$\frac{1}{T}\sum_{t=1+k}^{T+k}f(X_t).$$

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Nicholas Metropolis 1915 – 1999

W. Keith Hastings 1930 –

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Context

- Let b_j , $j = 1, \ldots, J$ be positive numbers.
- Let $B = \sum_{i} b_{j}$. If J is huge, B cannot be computed.

• Let
$$\pi_j = b_j/B$$
.

• We want to simulate a r.v. with pmf π_j .

Explore the set

- Consider a Markov process on $\{1, \ldots, J\}$ with transition probability Q.
- Designed to explore the space $\{1, \ldots, J\}$ efficiently
- Not too fast (and miss important points to sample)
- Not too slowly (and take forever to reach important points)

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Define another Markov process

- Based on the exact same states $\{1, \ldots, J\}$ as the previous ones
- Assume the process is in state *i*, that is $X_t = i$.
- Simulate the (candidate) next state *j* according to *Q*.
- Define

$$X_{t+1} = \begin{cases} j & \text{with probability } \alpha_{ij} \\ i & \text{with probability } 1 - \alpha_{ij} \end{cases}$$

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Transition probability P

$$\begin{array}{rcl} P_{ij} &=& Q_{ij}\alpha_{ij} & \text{if } i \neq j \\ P_{ii} &=& Q_{ii}\alpha_{ii} + \sum_{\ell \neq i} Q_{i\ell}(1 - \alpha_{i\ell}) & \text{otherwise} \end{array}$$

Must verify the property

$$1 = \sum_{j} P_{ij} = P_{ii} + \sum_{j \neq i} P_{ij}$$

= $Q_{ii} \alpha_{ii} + \sum_{\ell \neq i} Q_{i\ell} (1 - \alpha_{i\ell}) + \sum_{j \neq i} Q_{ij} \alpha_{ij}$
= $Q_{ii} \alpha_{ii} + \sum_{\ell \neq i} Q_{i\ell} - \sum_{\ell \neq i} Q_{i\ell} \alpha_{i\ell} + \sum_{j \neq i} Q_{ij} \alpha_{ij}$
= $Q_{ii} \alpha_{ii} + \sum_{\ell \neq i} Q_{i\ell}$

As $\sum_{i} Q_{ij} = 1$, we have $\alpha_{ii} = 1$.

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Time reversibility

$$\pi_i P_{ij} = \pi_j P_{ji}, \quad i \neq j$$

that is

$$\pi_i Q_{ij} \alpha_{ij} = \pi_j Q_{ji} \alpha_{ji}, \quad i \neq j$$

It is satisfied if

$$lpha_{ij} = rac{\pi_j Q_{ji}}{\pi_i Q_{ij}}$$
 and $lpha_{ji} = 1$

or

$$rac{\pi_i Q_{ij}}{\pi_j Q_{ji}} = lpha_{ji} ext{ and } lpha_{ij} = 1$$

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As α_{ij} is a probability

$$\alpha_{ij} = \min\left(\frac{\pi_j Q_{ji}}{\pi_i Q_{ij}}, 1\right)$$

Simplification

Remember: $\pi_j = b_j/B$. Therefore

$$\alpha_{ij} = \min\left(\frac{b_j B Q_{ji}}{b_i B Q_{ij}}, 1\right) = \min\left(\frac{b_j Q_{ji}}{b_i Q_{ij}}, 1\right)$$

The normalization constant B does not play a role in the computation of $\alpha_{ij}.$

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In summary

- Given Q and b_j
- $\bullet~{\rm defining}~\alpha$ as above
- creates a Markov process characterized by P
- with stationary distribution π .

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Algorithm

- Choose a Markov process characterized by Q.
- 2 Initialize the chain with a state *i*: t = 0, $X_0 = i$.
- Simulate the (candidate) next state *j* based on *Q*.
- Let r be a draw from U[0, 1[.

③ Compare
$$r$$
 with $lpha_{ij} = \min\left(rac{b_j \, \mathcal{Q}_{ji}}{b_i \, \mathcal{Q}_{ij}}, 1
ight)$. If

$$r < \frac{b_j Q_{ji}}{b_i Q_{ij}}$$

then
$$X_{t+1} = j$$
, else $X_{t+1} = i$.

Increase t by one.

Example

$$b = (20,8,3,1)$$

$$\pi = (\frac{5}{8}, \frac{1}{4}, \frac{3}{32}, \frac{1}{32})$$

$$Q = \begin{pmatrix} \frac{1}{4} & \frac{1}{4} & \frac{1}{4} & \frac{1}{4} \\ \frac{1}{4} & \frac{1}{4} & \frac{1}{4} & \frac{1}{4} \end{pmatrix}$$

Run MH for 10000 iterations. Collect statistics after 1000.

- Accept: [2488, 1532, 801, 283]
- Reject: [0, 952, 1705, 2239]
- Simulated: [0.627, 0.250, 0.095, 0.028]
- Target: [0.625, 0.250, 0.09375, 0.03125]

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Simulated annealing

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Gibbs sampling

Motivation

- Draw from multivariate distributions.
- Main difficulty: deal with correlations.

Metropolis-Hastings

- Let X = (X¹, X²,..., Xⁿ) be a random vector with pmf (or pdf) p(x).
- Assume we can draw from the marginals:

$$\Pr(X^{i}|X^{j} = x^{j}, j \neq i), i = 1, ..., n.$$

- Markov process. Assume current state is x.
 - Draw randomly (equal probability) a coordinate *i*.
 - Draw r from the *i*th marginal.
 - New state: $y = (x^1, ..., x^{i-1}, r, x^{i+1}, ..., x^n)$.

Gibbs sampling

Transition probability

$$Q_{xy} = \frac{1}{n} \operatorname{Pr}(X^{i} = r | X^{j} = x^{j}, j \neq i) = \frac{p(y)}{n \operatorname{Pr}(X^{j} = x^{j}, j \neq i)}$$

- The denominator is independent of X^{i} .
- So Q_{xy} is proportional to p(y).

Metropolis-Hastings

$$\alpha_{xy} = \min\left(\frac{p(y)Q_{yx}}{p(x)Q_{xy}}, 1\right) = \min\left(\frac{p(y)p(x)}{p(x)p(y)}, 1\right) = 1$$

The candidate state is always accepted.

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Example: bivariate normal distribution

$$\left(\begin{array}{c} X\\ Y \end{array}\right) \sim N\left(\left(\begin{array}{c} \mu_X\\ \mu_Y \end{array}\right), \left(\begin{array}{cc} \sigma_X^2 & \rho\sigma_X\sigma_Y\\ \rho\sigma_X\sigma_Y & \sigma_Y^2 \end{array}\right)\right)$$

Marginal distribution:

$$Y|(X = x) \sim N\left(\mu_Y + \frac{\sigma_Y}{\sigma_X}\rho(x - \mu_X), (1 - \rho^2)\sigma_Y^2\right)$$

Apply Gibbs sampling to draw from:

$$N\left(\left(\begin{array}{c}0\\0\end{array}\right), \left(\begin{array}{c}1&0.9\\0.9&1\end{array}\right)\right)$$

Note: just for illustration. Should use Cholesky factor.

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Gibbs sampling

Example: pdf



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Example: draws from Gibbs sampling



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Simulated annealing

Combinatorial optimization

 $\min_{x\in\mathcal{F}}f(x)$

where the feasible set \mathcal{F} is a large finite set of vectors.

Set of optimal solutions

$$\mathcal{X}^* = \{x \in \mathcal{F} | f(x) \leq f(y), \ \forall y \in \mathcal{F}\} \text{ and } f(x^*) = f^*, \ \forall x^* \in \mathcal{X}^*$$

Probability mass function on \mathcal{F}

$$p_{\lambda}(x) = rac{e^{-\lambda f(x)}}{\sum_{y \in \mathcal{F}} e^{-\lambda f(y)}}, \ \lambda > 0.$$

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Simulated annealing

$$p_{\lambda}(x) = rac{e^{-\lambda f(x)}}{\sum_{y \in \mathcal{F}} e^{-\lambda f(y)}}$$

Equivalently

$$p_{\lambda}(x) = \frac{e^{\lambda(f^* - f(x))}}{\sum_{y \in \mathcal{F}} e^{\lambda(f^* - f(y))}}$$

• As $f^* - f(x) \leq 0$, when $\lambda \to \infty$, we have

$$\lim_{\lambda o \infty} p_\lambda(x) = rac{\delta(x \in \mathcal{X}^*)}{|\mathcal{X}^*|},$$

where

$$\delta(x \in \mathcal{X}^*) = \left\{ egin{array}{cc} 1 & ext{if } x \in \mathcal{X}^* \\ 0 & ext{otherwise.} \end{array}
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Example

$$egin{aligned} \mathcal{F} &= \{1,2,3\} \; f(\mathcal{F}) = \{0,1,0\} \ p_\lambda(1) &= rac{1}{2+e^{-\lambda}} \ p_\lambda(2) &= rac{e^{-\lambda}}{2+e^{-\lambda}} \ p_\lambda(3) &= rac{1}{2+e^{-\lambda}} \end{aligned}$$

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Example



Simulated annealing

- If λ is large,
- we generate a Markov chain with stationary distribution $p_{\lambda}(x)$.
- The mass is concentrated on optimal solutions.
- As the normalizing constant is not needed, only $e^{\lambda(f^*-f(x))}$ is used.
- Construction of the Markov process through the concept of neighborhood.
- A *neighbor* y of x is obtained by simple modifications of x.
- The Markov process will proceed from neighbors to neighbors.
- The neighborhood structure must be designed such that the chain is irreducible, that is the whole space \mathcal{F} must be covered.
- It must be designed also such that the size of the neighborhood is reasonably small.

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Neighborhood

Metropolis-Hastings

- Denote N(x) the set of neighbors of x.
- Define a Markov process where the next state is a randomly drawn neighbor.
- Transition probability:

$$Q_{xy} = rac{1}{|\mathcal{N}(x)|}$$

Metropolis Hastings:

$$\alpha_{xy} = \min\left(\frac{p(y)Q_{yx}}{p(x)Q_{xy}}, 1\right) = \min\left(\frac{e^{-\lambda f(y)}|N(x)|}{e^{-\lambda f(x)}|N(y)|}, 1\right)$$

Neighborhood

Notes

• The neighborhood structure can always be arranged so that each vector has the same number of neighbors. In this case,

$$\alpha_{xy} = \min\left(\frac{e^{-\lambda f(y)}}{e^{-\lambda f(x)}}, 1\right)$$

- If y is better than x, the next state is automatically accepted.
- Otherwise, it is accepted with a probability that depends on λ.
- If λ is high, the probability is small.
- When λ is small, it is easy to escape from local optima.

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Heuristic

lssue

- The number of iterations needed to reach a stationary state and draw an optimal solution may exceed the number of feasible solutions in the set.
- The acceptance probability is very small.
- Therefore, a complete enumeration works better.
- The method is used as a heuristic.