# Introduction to Markov Chain Monte Carlo

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# Outline

Motivation

Examples

Summary

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- ► Today, we will study MCMC and show why and how it works.
- ► MCMC is very powerful in fitting very complicated models such as overly parameterized models.
- When usual MLE and Newton-Raphson methods do not work properly, MCMC will be the best choice.

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▶ A Markov Chain refers to a sequence of random variables  $(X_0, ..., X_n, ...)$  generated by a Markov process.

▶ A particular chain is defined most critically by its transition probabilities (or transition kernel),  $P(i,j) = P(i \rightarrow j)$ , which is the single step move of the process from state i to state j,

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- ▶ We start the chain by specifying a starting vector  $\pi(0)$ . Often all the elements of  $\pi(0)$  are zero except for a single element of 1.
- As the chain processes, the probability values get spread out over the possible states space.

► The probability that the chain has state value  $s_i$  at time or (step) t + 1 is given by the **Chapman** – **Kolomogrov** equation,

$$\pi_{i}(t+1) = P(X_{t+1} = s_{i})$$

$$= \sum_{k} P(x_{t+1} = s_{i} | X_{t} = s_{k}) P(X_{t} = s_{k})$$

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- ► Successive iteration of the **Chapman** − **Kolomogrov** describes the evolution of the chain.
- ▶ Define the probability transition matrix **P** as the matrix whose (i,j)-th element is P(i,j), the probability of moving from state i to state j,  $P(i \rightarrow j)$ . Which implies that

$$\sum_{i} P(i,j) = \sum_{i} P(i \to j) = 1$$

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▶ Defining the *n*-step transition probability  $P_{i,j}^{(n)}$ ,

$$p_{i,j}^{(n)} = P(X_{t+n} = s_j | X_t = s_i)$$

then  $p_{i,i}^{(n)}$  is just the (i,j)-th element of  $\mathbf{P}^n$ .

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- ▶ A chain is said to be **aperiodic** when the number of steps required to move between two states (say *x* and *y*) is not required to be multiple of some integer.
- ► The conditions for a <u>stationary</u> distribution is that the chain is **irreducible** and **aperiodic**.

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 $\blacktriangleright$   $\pi^*$  is the left eigenvector associated with the eigenvalue  $\lambda = 1$  of **P**.

► A sufficient condition for a <u>unique</u> stationary distribution is that the **detailed balance** equation holds (for all *i* and *j*),

$$P(j \to k)\pi_j^* = P(k \to j)\pi_k^*$$

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► The discrete-state Markov chain can be generalized to a continuous state Markov process by having a probability kernel P(x, y) that satisfies

$$\int P(x,y)dy = 1$$

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▶ At equilibrium, that the stationary distribution satisfies,

$$\pi^*(y) = \int \pi^*(x) P(x, y) dx$$

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- ► This is usuall the case in Bayesian statistics, for example,  $f(\theta) = \prod_{i=1}^{n} f_i(\theta), f_i(\theta) > 0.$

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- ► This is usuall the case in Bayesian statistics, for example,  $f(\theta) = \prod_{i=1}^{n} f_i(\theta), f_i(\theta) > 0.$
- ▶ In my next talk, we will study the slice sampler proposed by Neal (2013) to simulate those types of postierior distributions.

## The Metropolis Algorithm

- 1 Start with any initial value  $\theta^{(0)}$  satisfying  $f(\theta^{(0)}) > 0$ . Set k = 0.
- 2 Using the current values  $\theta^{(K)}$ , generate a value  $\theta^*$  from some jumping (candidate, or proposal) distribution,  $q(\theta^{(K)}, \theta^*)$ , which is the probability of returning a values of  $\theta^*$  given a previous value of  $\theta^{(K)}$ . Here the jump distribution is symmetric such that  $q(\theta^{(k)}, \theta^*) = q(\theta^*, \theta^{(k)})$ .
- 3 Calculate the ratio of the density at the candidate  $\theta^*$  and current  $\theta^{(k)}$  points,

$$\alpha = \frac{p(\theta^*)}{p(\theta^{(k-1)})} = \frac{f(\theta^*)}{f(\theta^{(k-1)})}$$

- 4 If  $\alpha \ge 1$  then set  $\theta^{(k+1)} = \theta^*$  and go to step 2. If  $\alpha < 1$  then with probability  $\alpha$  accept the candidate point, otherwise reject  $\theta^*$  and set  $\theta^{(k+1)} = \theta^{(k)}$  then go to step 2.
- 4' Generate a uniform distribution u = U(0, 1). if  $u < \min(\alpha, 1)$  then set  $\theta^{(k+1)} = \theta^*$ ; Otherwise set  $\theta^{(k+1)} = \theta^{(k)}$  and go step 2.

### The Metropolis-Hastings Algorithm (Cont'd)

▶ Hastings (1970) generalized the Metropolis algorithm by using an arbitrary transition probability function  $q(\theta_1, \theta_2) = Pr(\theta_1 \rightarrow \theta_2)$ , and setting the acceptance probability for a candidate point as

$$\alpha = \frac{f(\theta^*)q(\theta^*|\theta^{(k-1)})}{f(\theta^{(k-1)}q(\theta^{(k-1)}|\theta^*)}$$

► Chib and Greenberg (1995) said that finding a jumping function appears to be a search for the proverbial needle in a haystack.

## Example 1 – fit a Student -t distribution

Suppose that we have a data set  $\mathbf{y} = \{y_i, i = 1, ..., n\}$ , which is generated from a Student-t distribution  $y \sim t(v)$  with pdf  $f(y) = \frac{v^{v/2}\Gamma((v+1)/2)}{\Gamma(v/2)\Gamma(1/2)}(v+y^2)^{-(v+1)/2}.$ 

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▶ The likelihood of the data based on the model is

$$f(\mathbf{y}|\nu) = \prod_{i=1}^{n} f(y_i|\nu)$$

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 $\triangleright$  The posterior distribution of v is

$$f(v|\mathbf{y}) \propto f(v) \prod_{i=1}^{n} \frac{v^{v/2} \Gamma((v+1)/2)}{\Gamma(v/2)} (v+y_i^2)^{-(v+1)/2}$$

where f(v) is the prior distribution of v.

#### Example 1 – fit a Student -t distribution (Cont'd)

Since this full conditional is an unknown distribution, we use a random-walk Metropolis-Hastings algorithm, in which the proposal density is a standard Gaussian density.

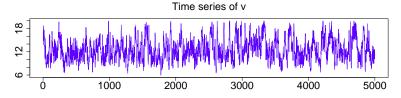


Figure: The training data was drawn from t(10) with sample size =1000. The estimate of v is v = 12.57 with standard deviation sd = 2.73.

► The MH algorithm was iterated 10,000 times. After the first 5000 iterations were discarded as burn in, parameter was estimated by sample mean.

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- ▶ Define  $X \in \mathbb{R}^m$  be the associated vector of covariates.
- ▶ Supposed that we obtain a sample  $\{y_i, X_i\}_{i=1}^n$ , where n is the sample size.
- ► The relationship between *y* and *X* can be modeled by a generalized linear model (GLM),

$$f(y_i|X_i,\theta) = \beta_0 + X_i^T \beta + \epsilon_i, \quad i = 1,...,n$$

where  $\epsilon_i$  are *iid* white noises with mean 0 and variance  $\sigma^2$ .

## Logistic regression (LR) model (Cont'd)

▶ In logistic regression, we use a logistic function p(.), defined below,

$$p(y|X,\theta) = \frac{e^{\beta_0 + X^T \beta}}{1 + e^{\beta_0 + X^T \beta}}.$$

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▶ After a bit of manipulation, we obtain the following

$$\frac{p(y|X,\theta)}{1-p(y|X,\theta)} = e^{\beta_0 + X^T \beta}, \theta^T = (\beta_0, \beta_1, ..., \beta_m)$$

The quantity of  $p(y|X,\theta)/(1-p(y|X,\theta))$  is called the odds, which can take any values in  $(0,\infty)$ .

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► The usual logistic regression model is,

$$\log\left(\frac{p(y|X,\theta)}{1-p(y|X,\theta)}\right) = \beta_0 + X^T \beta.$$

#### Markov Chain Monte Carlo (MCMC) estimation

▶ The likelihood function of the model is

$$L(y|X,\theta) = \prod_{i=1}^{n} p(X_i;\theta)^{y_i} (1 - p(X_i;\theta))^{1 - y_i}$$

$$= \prod_{i=1}^{n} \left( \frac{e^{\beta_0 + X_i^T \beta}}{1 + e^{\beta_0 + X_i^T \beta}} \right)^{y_i} \left( 1 - \frac{e^{\beta_{10} + X_i^T \beta}}{1 + e^{\beta_0 + X_i^T \beta}} \right)^{1 - y_i}$$

Note: All the parameters in the model are treated as random variables with prior distributions.

# Metropolis-Hastings (MH) sampler

0 Let  $\gamma$  be a parameter in  $\theta$ , and define  $\theta_{-\gamma}$  as all parameters in  $\theta$  except  $\gamma$ .

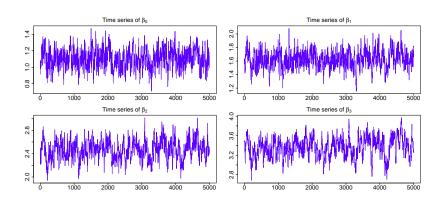
## Metropolis-Hastings (MH) sampler

- 0 Let  $\gamma$  be a parameter in  $\theta$ , and define  $\theta_{-\gamma}$  as all parameters in  $\theta$  except  $\gamma$ .
- 1. Initialize  $\gamma^{(0)} \sim q(.)$ , where q(.) is a candidate distribution, which is a normal distribution,  $N(\mu, \sigma^2)$ .
- 2. For k=1, ..., K, repeat:
  - (a) Sample  $\gamma^* \sim q(\gamma|\gamma^{(k-1)})$  such that  $\gamma^* \sim N(\gamma^{(k-1)}, \sigma^2)$ .
  - (b) Calculate acceptance probability:

$$\begin{split} \alpha &= \min \bigg\{ 1, \frac{L(y|x, \theta_{-\gamma}, \gamma^{(*)}) q(\gamma^{(k-1)}|\gamma^{(*)})}{L(y|x, \theta_{-\gamma}, \gamma^{(k-1)}) q(\gamma^{(*)}|\gamma^{(k-1)})} \bigg\} \\ &= \min \bigg\{ 1, \frac{L(y|x, \theta_{-\gamma}, \gamma^{(*)})}{L(y|x, \theta_{-\gamma}, \gamma^{(k-1)})} \bigg\}, \underbrace{\text{as } q(.) \text{ is symmetric.}}. \end{split}$$

- (c) Sample  $u \sim \text{Uniform } (0, 1)$ .
- (d) If  $u < \alpha$  then accept the proposal  $\gamma^{(k)} = \gamma^*$ ; Otherwise, reject the proposal and set  $\gamma^{(k)} = \gamma^{(k-1)}$ .

## Example 2 – fit a logistic regression model – sampled time series



The MH algorithm was iterated 10,000 times. After the first 5000 iterations were discarded, parameters were estimated by sample means.

Table: Fit a logistic regression model.

Parameter	True	Estimate	SD
$eta_0$	1	1.10	0.12
$eta_1$	1.5	1.62	0.13
$eta_2$	2.2	2.45	0.17
$\beta_3$	3	3.36	0.20

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- ► Two examples were given to show how MCMC works.
- ▶ The usual logistic regression model can be extended to,

$$\log\left(\frac{p(y|X,\theta)}{1-p(y|X,\theta)}\right) = \beta_0 + X^T \beta + \epsilon,$$

where  $\epsilon$  is white noises with mean 0 and variance  $\sigma^2$ .

Motivation Examples Summary

# Thank you!

#### References

- [1] Robert, C. and G., Casella. 2013. Monte Carlo Statistical Methods: Edition 2. London: Chapman and Hall.
- [2] Gilks, W. R., Richardson, S., and Spiegelhalter, D. J. 1995. Markov Chain Monte Carlo in Practice. London: Chapman and Hall.
- [3] Neal, R. N. 2003. Slice sampling. *The Annals of Statistics* 31: 705-767.
- [4] Chib, S. and Greenberg E. 1995. Understanding the Metropolis-Hastings Algorithm. *The American Statistician* Vol. 49, No. 4: 327-335.