

# **Advanced Statistical Computing**

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# Markov Chain Monte Carlo

- The goal is to generate a sequence of random samples from an arbitrary probability density function (usually high dimensional),
- The sequence of random samples form a Markov chain,
- The purpose is simulation (Monte Carlo).

# What is Monte Carlo



# What is Monte Carlo?

- Rely on repeated sampling to study the results of a experiment or study the properties of certain procedure.
  - Often used in complex and uncertain scenarios
  - Difficult to formulate, high correlation.
  - Cheap
  - Take advantage of faster computers
- History
  - John von Neumann, Stanislaw Ulam, Nicholas Metropolis

# Applications of Monte Carlo

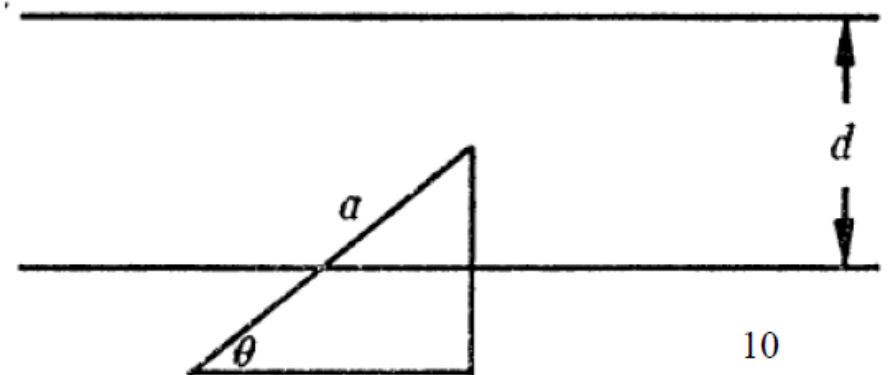
- Optimization
- Numerical integration
- Generate random samples

# Buffon's needle

Georges-Louis Leclerc, Comte de Buffon  
(1707-1788)



Given a needle of length  $a$  and an infinite grid of parallel lines with common distance  $d$  between them, what is the probability  $P(E)$  that a needle, tossed at the grid randomly, will cross one of the parallel lines?





# Buffon's needle

- Assume  $a < d$

$$P(E) = \int_0^\pi \frac{a \sin \theta d\theta}{\pi d} = (a/\pi d) \int_0^\pi \sin \theta d\theta = 2a/\pi d.$$

<http://web.student.tuwien.ac.at/~e9527412/buffon.html>

# Motivation

- Generate *iid* r.v. from high-dimensional arbitrary distributions is extremely difficult.
- Drop the “*independent*” requirement.
- How about also drop the “*identically distributed*” requirement?



# Markov chain

- Assume a finite state, discrete Markov chain with  $N$  different states.

- Random process  $X_n$ ,  $n = 0, 1, 2, \dots$

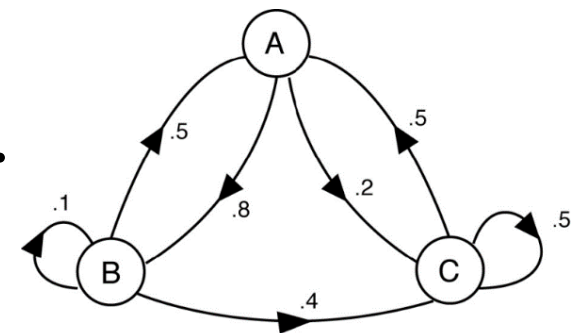
$$x_n \in S = \{1, 2, \dots, N\}$$

- Markov property,

$$P(X_{n+1} = x \mid X_1 = x_1, X_2 = x_2, \dots, X_n = x_n) = P(X_{n+1} = x \mid X_n = x_n)$$

- Time-homogeneous
- Order

– Future state depends on the past  $m$  states.



Markov graph of transition probabilities between states A, B and C

# Key parameters

- Transition matrix

$$P(X_n = j \mid X_{n-1} = i) = p(i, j),$$

$$P = \{p(i, j)\}.$$

- Initial probability distribution  $\pi^{(0)}$

$$\pi^{(n)}(i) = P(x_n = i).$$

- Stationary distribution (invariant/equilibrium)

$$\pi = \pi P.$$

# Reducibility

- A state  $j$  is accessible from state  $i$  (written  $i \rightarrow j$ ) if  $P(X_n = j \mid X_0 = i) = p_{ij}^{(n)} > 0$ .
- A Markov chain is *irreducible* if it is possible to get to any state from any state.

# Recurrence

- A state  $i$  is *transient* if given that we start in state  $i$ , there is a non-zero probability that we will never return to  $i$ . State  $i$  is *recurrent* if it is not transient.

# Ergodicity

- A state  $i$  is *ergodic* if it is aperiodic and positive recurrent. If all states in an irreducible Markov chain are ergodic, the chain is *ergodic*.

# Reversible Markov chains

- Consider an ergodic Markov chain that converges to an invariant distribution  $\pi$ . A Markov chain is *reversible* if for all  $x, y \in S$ ,

$$\pi(x)p(x, y) = \pi(y)p(y, x).$$

which is known as the detailed balance equation.

- An ergodic chain in equilibrium and satisfying the detailed balance condition has  $\pi$  as its unique stationary distribution.

# Markov Chain Monte Carlo

- The goal is to generate sequence of random samples from an arbitrary probability density function (usually high dimensional),
- The sequence of random samples form a Markov chain,

in Markov chain,  $P \rightarrow \pi$

in MCMC,  $\pi \rightarrow P$ .



# Examples

$$P(X | E, \beta, \sigma^2) \propto \prod_{k=1}^K \prod_{E(i)=k} \prod_{j=1}^M \left( (\sigma_{kj}^2)^{-\frac{1}{2}} e^{-\frac{1}{2\sigma_{kj}^2} (x_{ij} - \beta_{kj})^2} \right)$$

$$P(X | E) = \prod_{k=1}^K \prod_{j=1}^M \iint \prod_{E(i)=k} P(x_{ij} | \beta_{kj}, \sigma_{kj}^2) p(\beta_{kj} | \beta_0, \sigma_{kj}^2) p(\sigma_{kj}^2) d\beta_{kj} d\sigma_{kj}^2$$

$$= \prod_{k=1}^K \prod_{j=1}^M \left[ \frac{b^a}{\Gamma(a)} \frac{(2\pi)^{-\frac{n_k}{2}}}{\sqrt{n_k+1}} \frac{\Gamma\left(\frac{n_k}{2} + a\right)}{\left( b + \frac{1}{2} \left( \sum_{E(i)=k} x_{ij}^2 + \beta_0^2 - \frac{\left( \sum_{E(i)=k} x_{ij} + \beta_0 \right)^2}{n_k + 1} \right) \right)^{\frac{n_k}{2} + a}} \right]$$

# Bayesian Inference

Genotype  $Y = (y_1, \dots, y_n)$

Haplotype  $Z = (z_1, \dots, z_n)$

Frequency  $\Theta = (\theta_1, \dots, \theta_m)$

Prior  $\theta \sim \text{Dirichlet}(\beta)$

$$P(Y, Z, \Theta) = \prod_{i=1}^n \theta_{z_{i1}} \theta_{z_{i2}} \prod_{g=1}^m \theta_g^{\beta_g - 1}$$

# History

- Metropolis, N., Rosenbluth, A. W., Rosenbluth, M. N., Teller, A. H. and Teller, E. (1953).

Equation of state calculations by fast computing machines. *Journal of Chemical Physics*, **21**, 1087–1092.

# Metropolis algorithm

- Direct sampling from the target distribution is difficult,
- Generating candidate draws from a proposal distribution,
- These draws then “corrected” so that asymptotically, they can be viewed as random samples from the desired target distribution.

# Pseudo code

- Initialize  $X_0$ ,
- Repeat
  - Sample  $Y \sim q(x,.)$ ,
  - Sample  $U \sim \text{Uniform}(0,1)$ ,
  - If  $U \leq \alpha(X,Y)$ , set  $X_i = y$ ,
  - Otherwise  $X_i = x$ .

# An informal derivation

- Find  $\alpha(X,Y)$ :
- Joint density of current Markov chain state and the proposal is  $g(x,y) = q(x,y)\pi(x)$
- Suppose  $q$  satisfies detail balance
$$q(x,y) \pi(x) = q(y,x)\pi(y)$$
- If  $q(x,y) \pi(x) > q(y,x)\pi(y)$ , introduce
$$\alpha(x,y) < 1 \text{ and } \alpha(y,x) = 1$$
hence  $\alpha(x,y) = \frac{\pi(y)q(y,x)}{\pi(x)q(x,y)}$ .
- If  $q(x,y) \pi(x) > q(y,x)\pi(y)$ , ...
- The probability of acceptance is  $\alpha(x,y) = \min\left(1, \frac{\pi(y)q(y,x)}{\pi(x)q(x,y)}\right)$ .

# Metropolis-Hastings Algorithm

- Start with any  $X^{(0)}=x_0$ , and a “*proposal chain*”  $T(x,y)$
- Suppose  $X^{(t)}=x_t$ . At time  $t+1$ ,
  - **Draw**  $y \sim T(x_t, y)$  (i.e., propose a move for the next step)
  - Compute “*goodness ratio*”

$$r = \frac{\pi(y)T(y, x_t)}{\pi(x_t)T(x_t, y)}$$

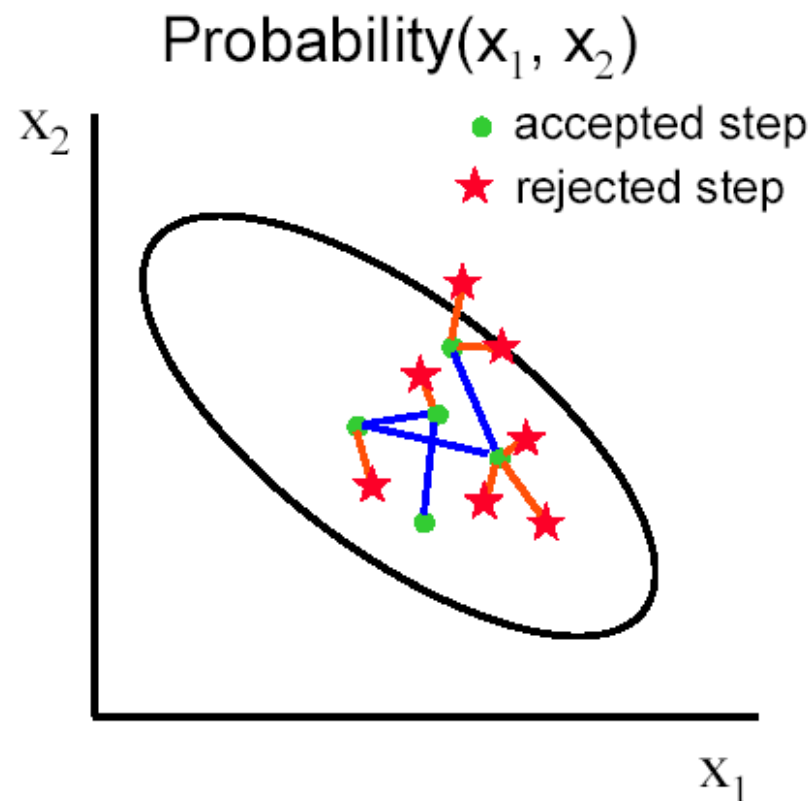
- **Acceptance/Rejection decision:** Let

$$X^{(t+1)} = \begin{cases} y, & \text{with } p = \min\{1, r\} \\ x_t, & \text{with } 1 - p \end{cases}$$



# Remarks

- Relies only on calculation of target pdf up to a normalizing constant.



# Remarks

- How to choose a good proposal function is crucial.
- Sometimes tuning is needed.
  - Rule of thumb: 30% acceptance rate
- Convergence is slow for high dimensional cases.

# Illustration of Metropolis-Hastings

- Suppose we try to sample from a bi-variate normal distributions.

$$N\left(\begin{pmatrix} 0 \\ 0 \end{pmatrix}, \begin{pmatrix} 1 & \rho \\ \rho & 1 \end{pmatrix}\right)$$

- Start from  $(0, 0)$

- Proposed move at each step is a two dimensional random walk

$$x_{t+1} = x_t + s \cos \theta$$

$$y_{t+1} = y_t + s \sin \theta$$

*with*

$$s \sim U(0,1)$$

$$\theta \sim U(0,2\pi)$$

# Illustration of Metropolis-Hastings

- At each step, calculate  $r = \frac{\pi(x_{t+1}, y_{t+1})}{\pi(x_t, y_t)}$

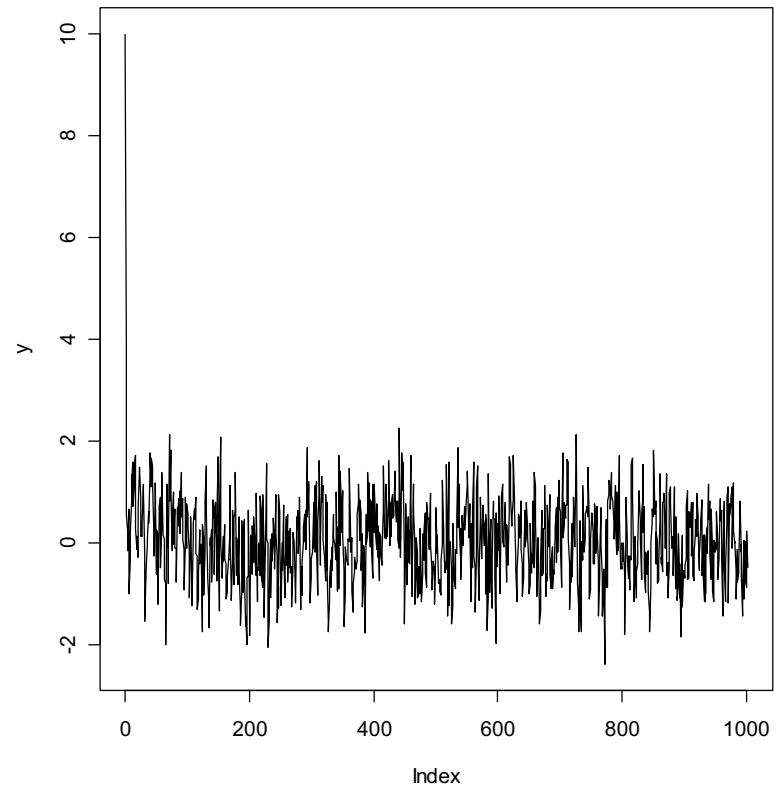
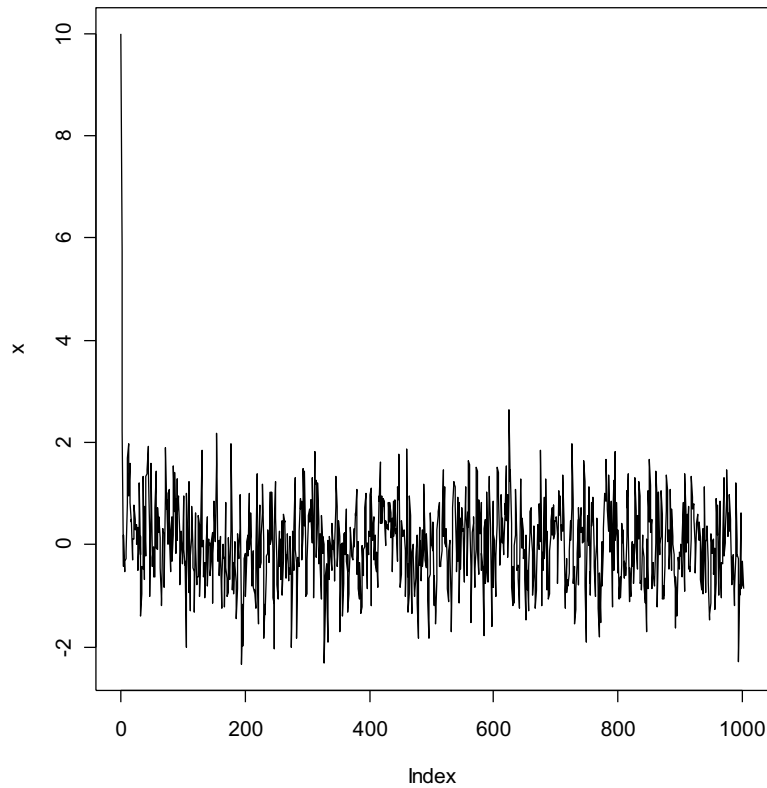
$$T((x_t, y_t), (x_{t+1}, y_{t+1})) = T((x_{t+1}, y_{t+1}), (x_t, y_t)) = 1/\pi^2$$

since

$$\begin{aligned} r &= \frac{\frac{1}{2\pi\sqrt{1-\rho^2}} \exp\left(-\frac{1}{2(1-\rho^2)}(x_{t+1}^2 - 2\rho x_{t+1}y_{t+1} + y_{t+1}^2)\right)}{\frac{1}{2\pi\sqrt{1-\rho^2}} \exp\left(-\frac{1}{2(1-\rho^2)}(x_t^2 - 2\rho x_t y_t + y_t^2)\right)} \\ &= \exp\left(-\frac{1}{2(1-\rho^2)}\left((x_{t+1}^2 - 2\rho x_{t+1}y_{t+1} + y_{t+1}^2) - (x_t^2 - 2\rho x_t y_t + y_t^2)\right)\right) \end{aligned}$$

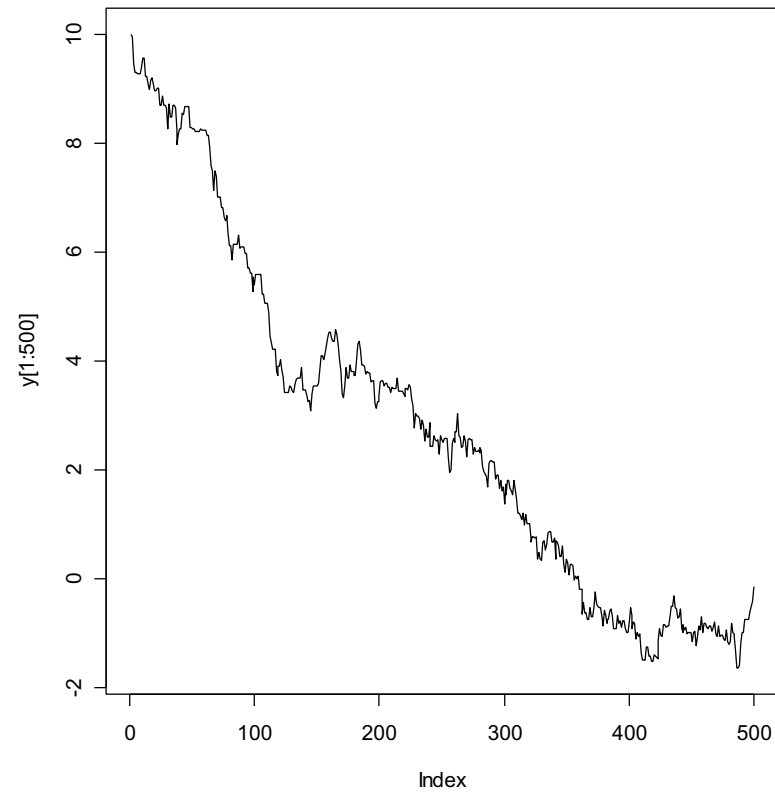
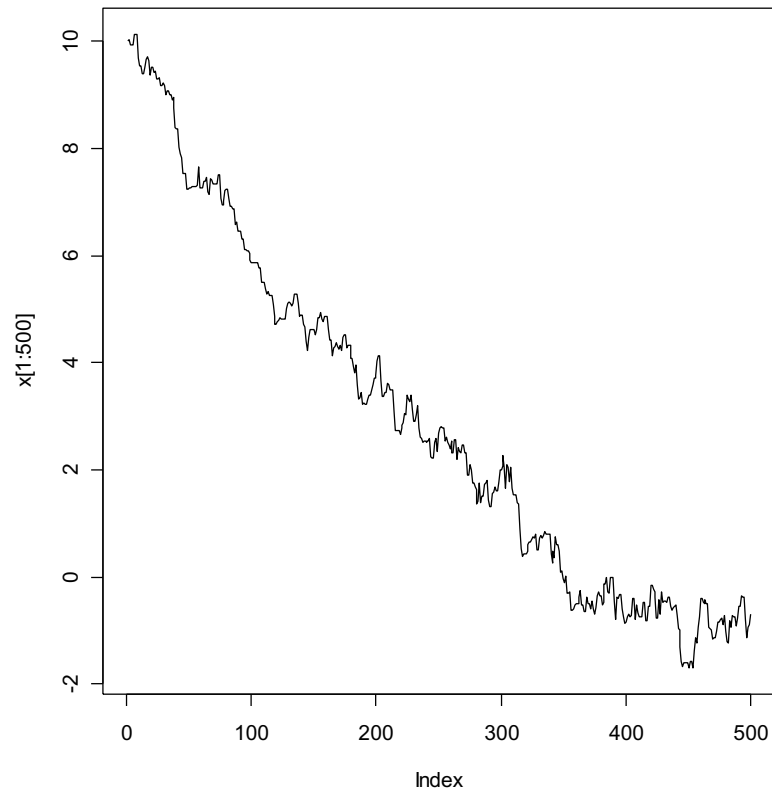
# Convergence

- Trace plot      Good



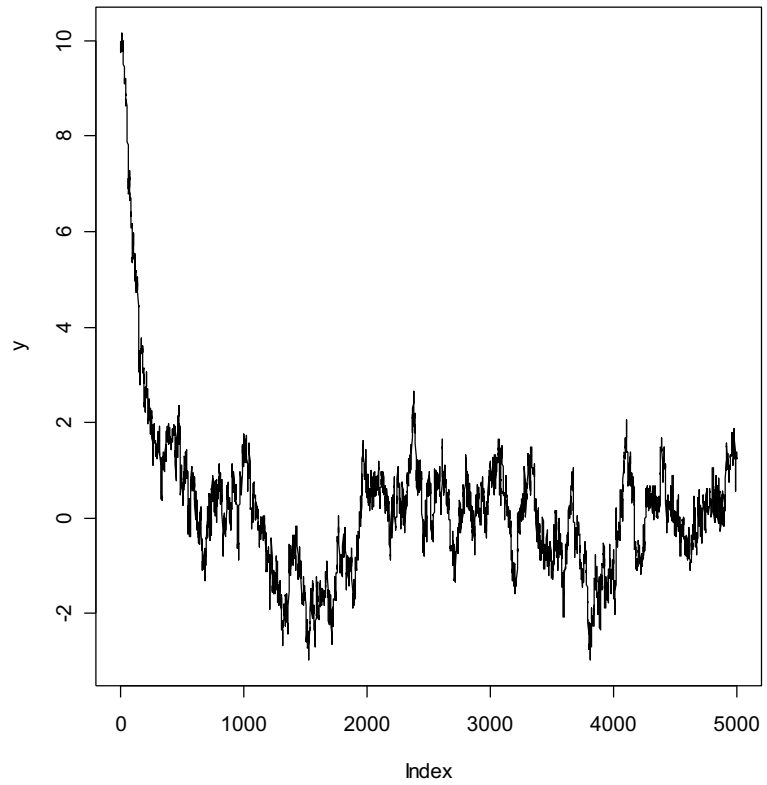
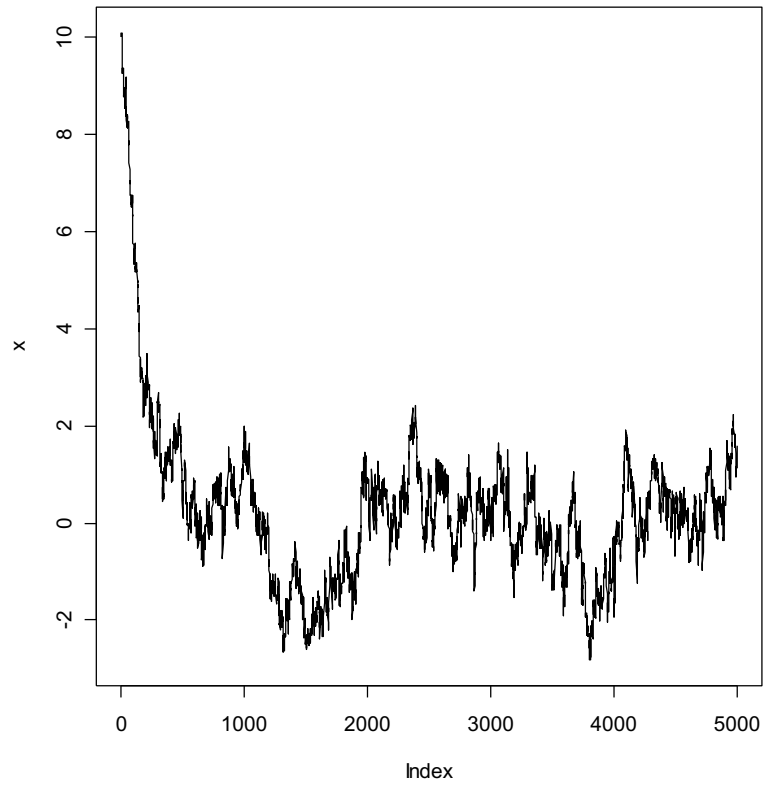
# Convergence

- Trace plot      Bad



# Convergence

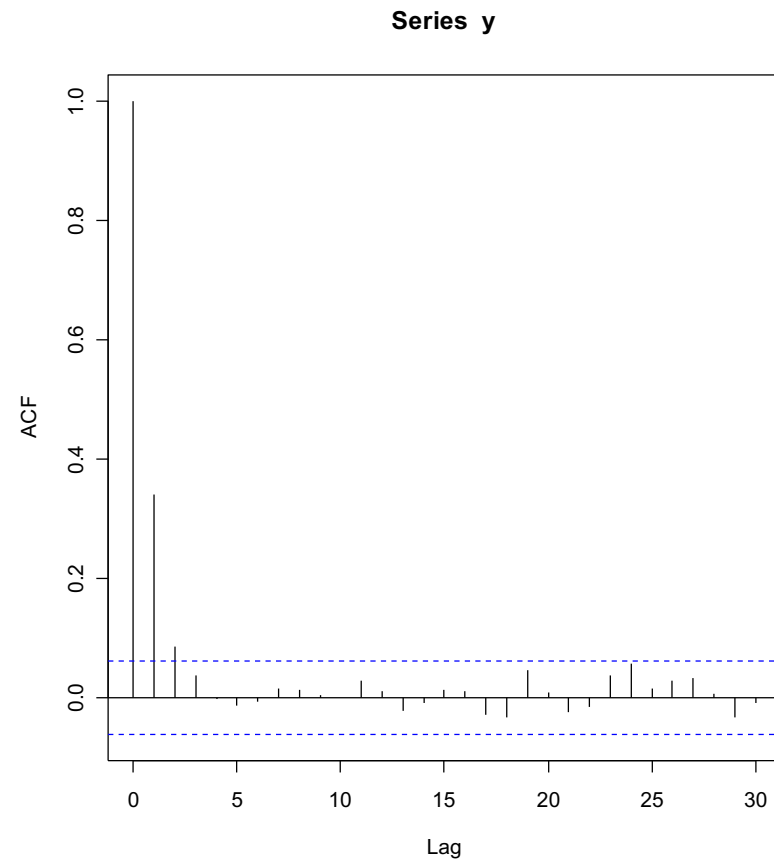
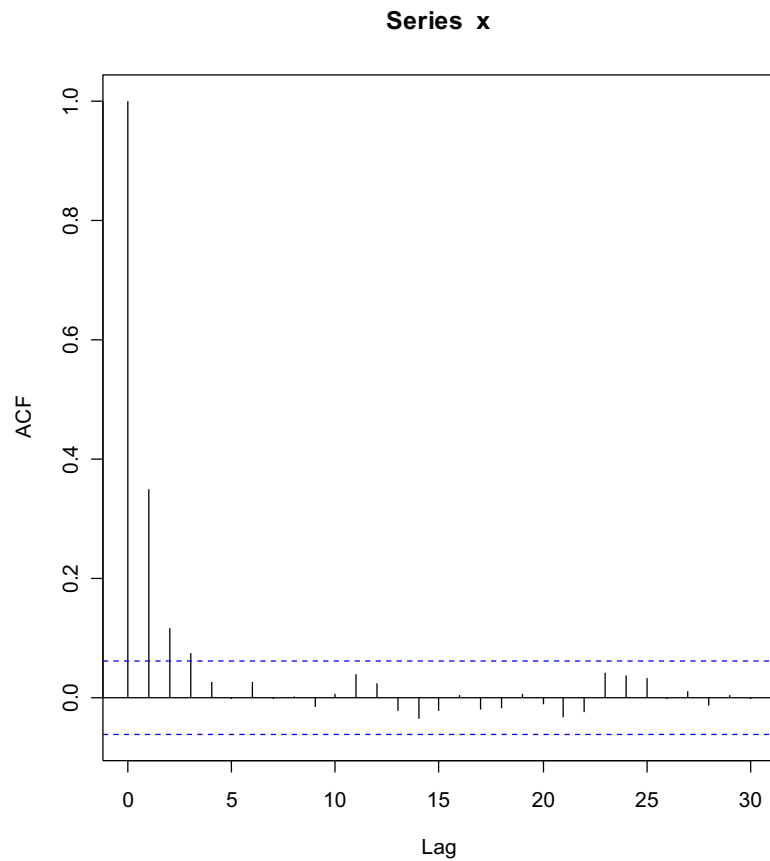
- Trace plot





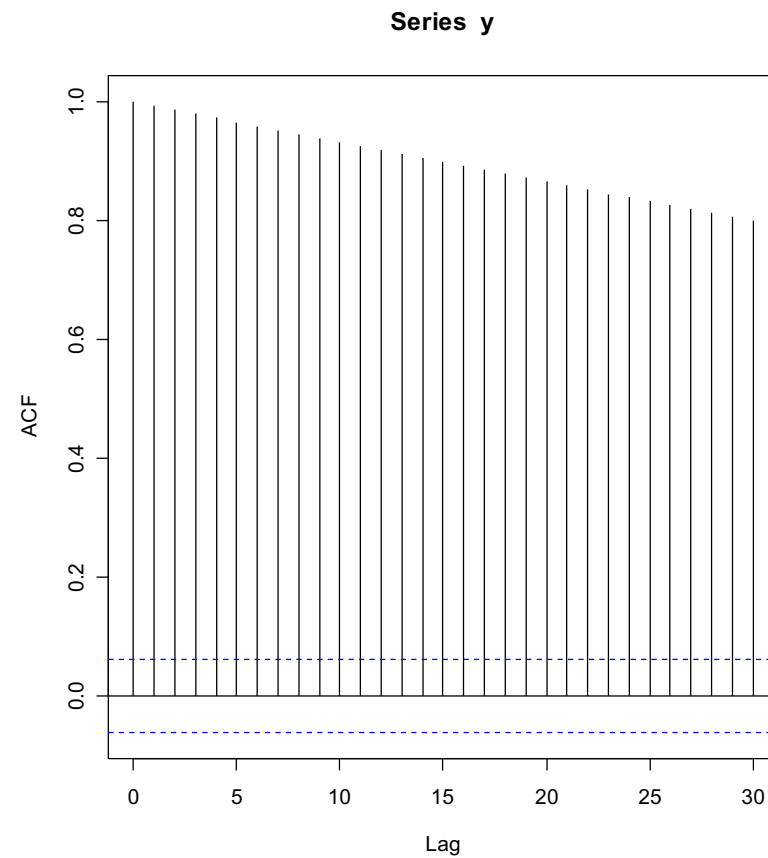
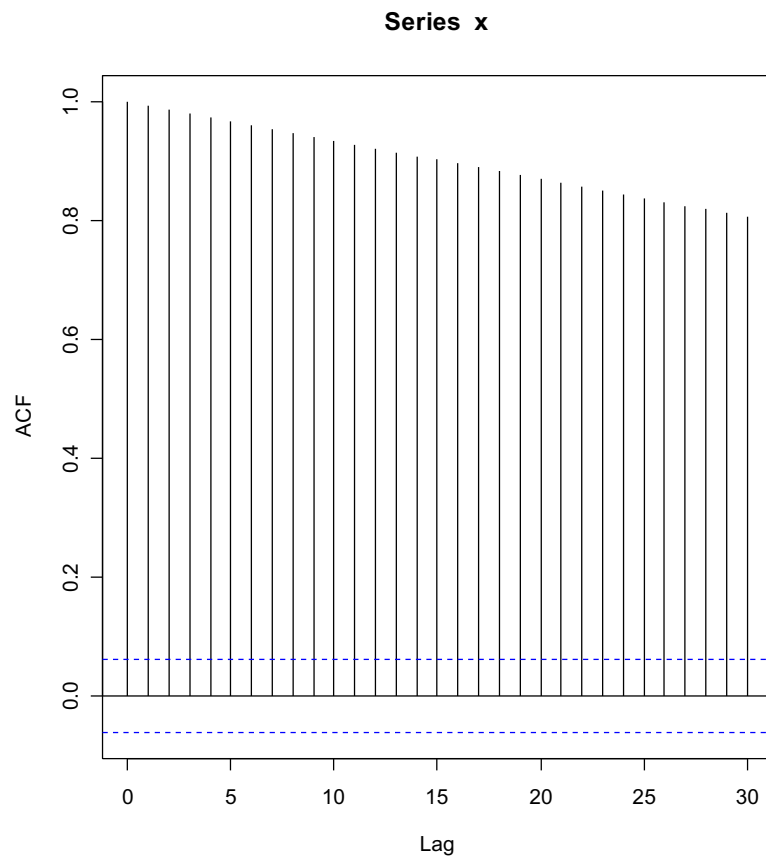
# Convergence

- Autocorrelation plot    Good



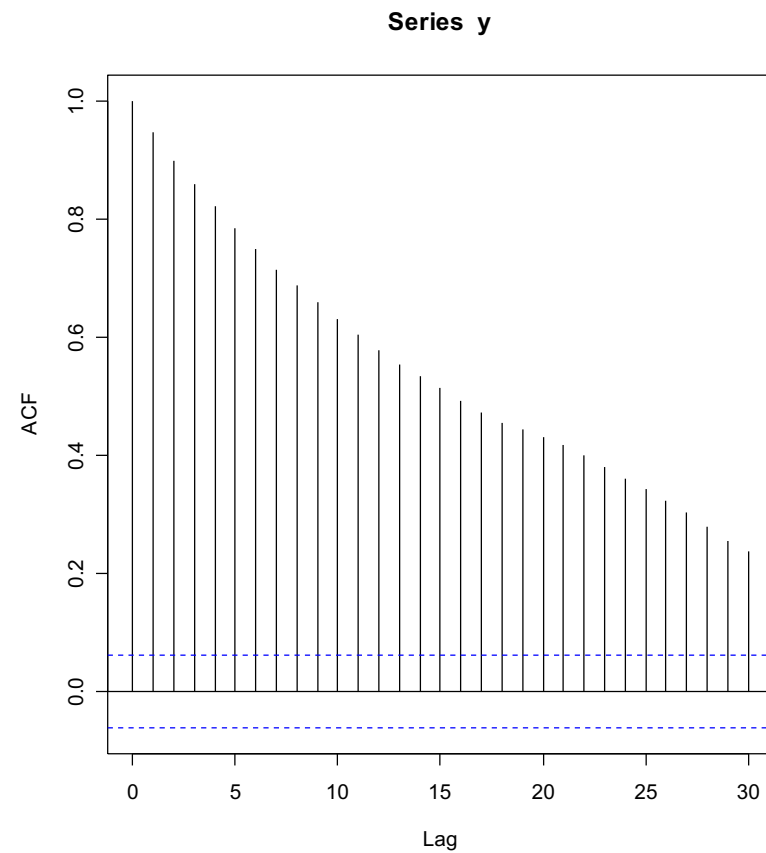
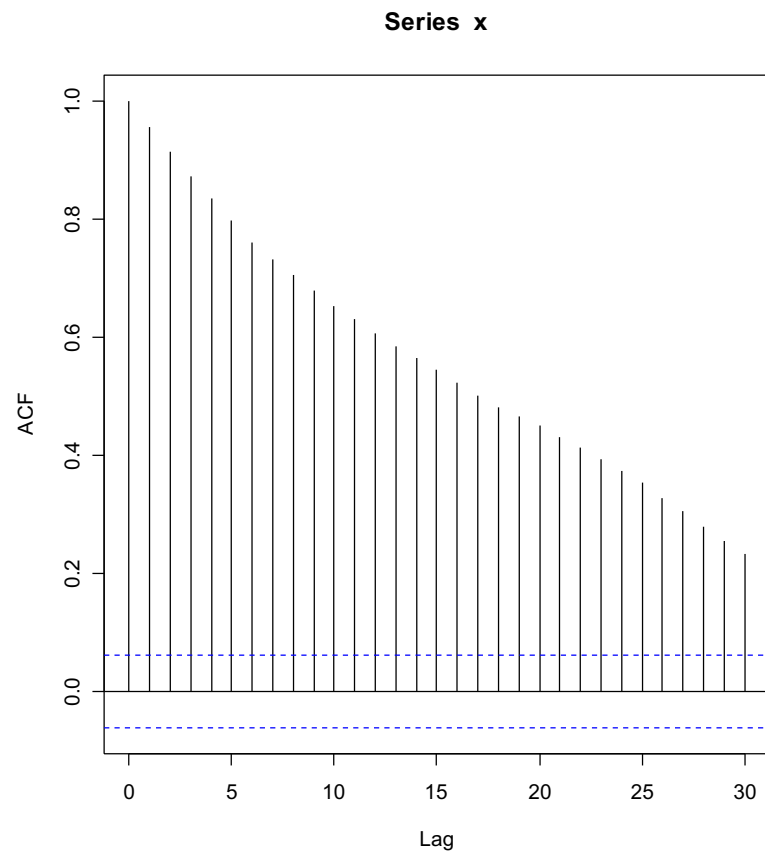
# Convergence

- Autocorrelation plot    Bad     $s = 0.5$



# Convergence

- Autocorrelation plot    Okay     $s = 3.0$



# References

- Metropolis et al. 1953,
- Hastings 1973,
- Tutorial paper:  
Chib and Greenberg (1995). Understanding the Metropolis--Hastings Algorithm. *The American Statistician* **49**, 327-335.

# Gibbs Sampler

- **Purpose:** Draw random samples from a joint distribution (high dimensional)

$$x = (x_1, x_2, \dots, x_n) \text{ Target } \pi(x)$$

- **Method:** Iterative conditional sampling

$$\forall i, \text{ draw } x_i \sim \pi(x_i \mid x_{[-i]})$$

# Illustration of Gibbs Sampler

- Suppose the target distribution is:

$$(X, Y) \sim N\left(\begin{pmatrix} 0 \\ 0 \end{pmatrix}, \begin{pmatrix} 1 & \rho \\ \rho & 1 \end{pmatrix}\right)$$

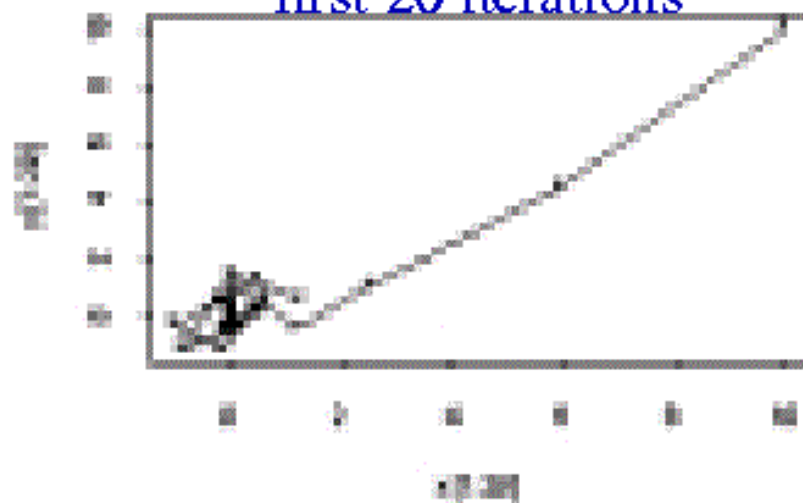
- Gibbs sampler:

$$[X|Y = y] \sim N(\rho y, 1 - \rho^2)$$

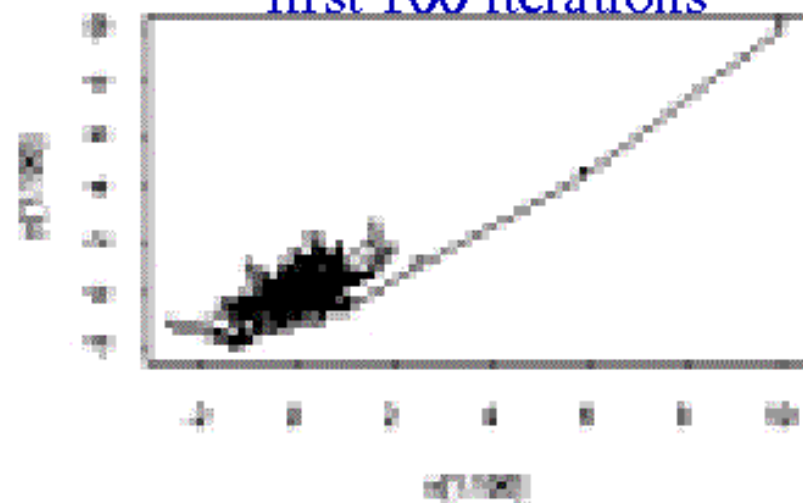
$$[Y|X = x] \sim N(\rho x, 1 - \rho^2)$$

Start from, say,  $(X, Y) = (10, 10)$ , we can take a look at the trajectories. We took  $\rho = 0.6$ .

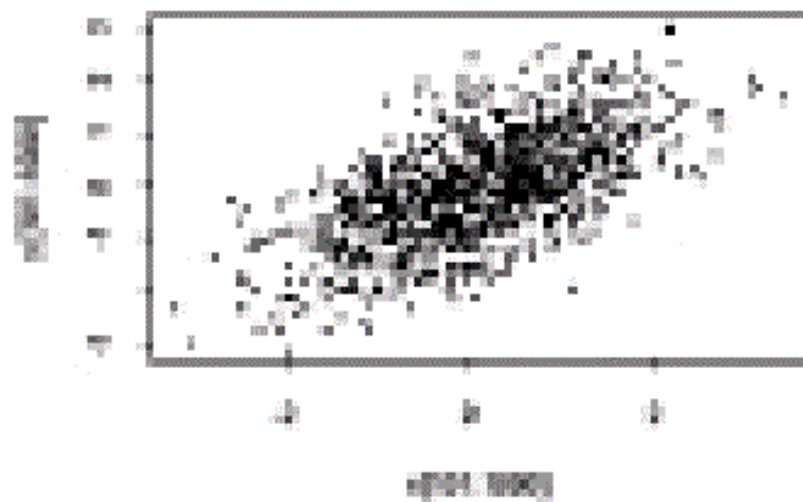
first 20 iterations



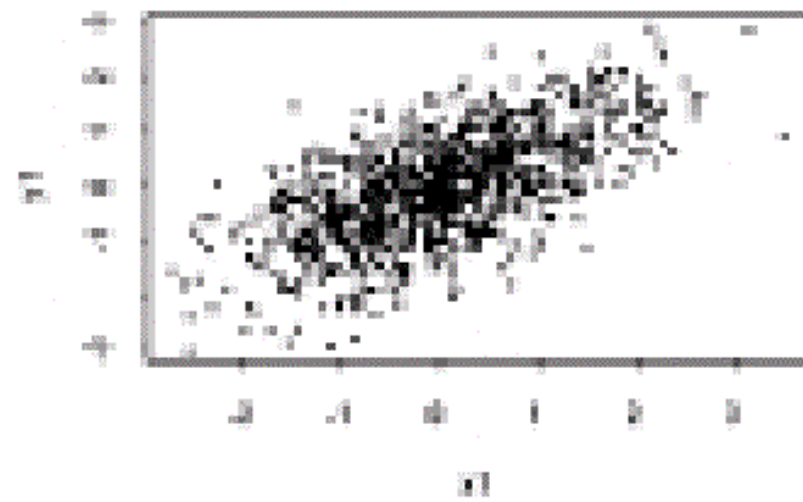
first 100 iterations



101-1000 iterations



900 iid samples





# References

- Geman and Geman 1984,
- Gelfand and Smith 1990,
- Tutorial paper:  
Casella and George (1992). Explaining the Gibbs sampler. *The American Statistician*, **46**, 167-174.

# Remarks

- Gibbs Sampler is a special case of Metropolis-Hastings
- Compare to EM algorithm, Gibbs sampler and Metropolis-Hastings are stochastic procedures
- Verify convergence of the sequence
- Require Burn in
- Use multiple chains

# References

- Geman and Geman 1984,
- Gelfand and Smith 1990,
- Tutorial paper:  
Casella and George (1992). Explaining the Gibbs sampler. *The American Statistician*, **46**, 167-174.