Probabilistic Inference (CO-493)

Imperial College London

Sampling

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February 12, 2019

Learning Material

- ▶ Bishop: Pattern Recognition and Machine Learning, Chapter 11
- ► MacKay: Information Theory, Inference and Learning Algorithms, Chapter 29
 http://www.inference.org.uk/itprnn/book.html
- ► Iain Murray's MCMC Tutorial: http://videolectures.net/mlss09uk_murray_mcmc/

Monte Carlo Methods—Motivation

- Monte Carlo methods are computational techniques that make use of random numbers
- ► Two typical problems:
 - 1. **Problem 1:** Generate samples $\{x^{(s)}\}$ from a given probability distribution p(x), e.g., for simulation (generative models) or representations of distributions

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➤ Examples: Means/variances of distributions, marginal likelihood, predictions in a Bayesian model

Complication: Integral cannot be evaluated analytically

Approximate Integration

- Numerical integration (low-dimensional problems)
- Bayesian quadrature, e.g., O'Hagan (1987, 1991); Rasmussen & Ghahramani (2003)
- Laplace approximation
- ► Variational inference, e.g., Jordan et al. (1999), Blei et al. (2017)
- ► Expectation Propagation, Opper & Winther (2001); Minka (2001)
- Monte-Carlo Methods, e.g., Gilks et al. (1996), Robert & Casella (2013), Bishop (2006)

Problem 2: Monte Carlo Estimation

Computing expectations via statistical sampling:

$$\mathbb{E}[f(\mathbf{x})] = \int f(\mathbf{x})p(\mathbf{x})d\mathbf{x}$$

$$\approx \frac{1}{S} \sum_{s=1}^{S} f(\mathbf{x}^{(s)}), \quad \mathbf{x}^{(s)} \sim p(\mathbf{x})$$

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► Making predictions (e.g., Bayesian regression with inputs *x* and targets *y*)

$$p(y|x) = \int p(y|\theta, x) \underbrace{p(\theta)}_{\text{Parameter distribution}} d\theta$$

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Key problem: Generating samples from p(x) or p(θ)
 Need to solve Problem 1

Properties of Monte Carlo Sampling

$$\mathbb{E}[f(\mathbf{x})] = \int f(\mathbf{x})p(\mathbf{x})d\mathbf{x}$$

$$\approx \frac{1}{S} \sum_{s=1}^{S} f(\mathbf{x}^{(s)}), \quad \mathbf{x}^{(s)} \sim p(\mathbf{x})$$

Estimator is asymptotically consistent, i.e.,

$$\lim_{S \to \infty} \frac{1}{S} \sum_{s=1}^{S} f(\mathbf{x}^{(s)}) = \mathbb{E}[f(\mathbf{x})] + \epsilon$$

- ▶ Error ϵ is normal (Gaussian) and its variance shrinks $\propto 1/S$, independent of the dimensionality
- Estimator is unbiased

Monte Carlo Estimation

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How do we get these samples?

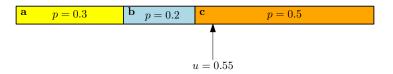
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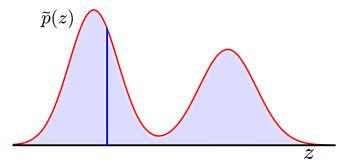
- ▶ How do we get these samples?
- ▶ Need to solve Problem 1
 - Sampling from simple distributions
 - Sampling from complicated distributions

Sampling Discrete Values



- $u \sim \mathcal{U}[0,1]$, where \mathcal{U} is the uniform distribution
- $u = 0.55 \Rightarrow x = c$

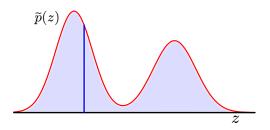
Continuous Variables



More complicated.

Geometric intuition: sample uniformly from the area under the curve

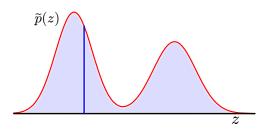
Rejection Sampling: Setting



Assume:

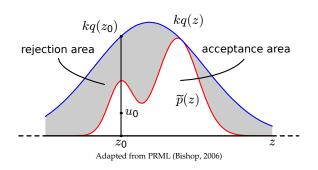
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- ► Evaluating $\tilde{p}(z) = Zp(z)$ is easy (and Z may be unknown)

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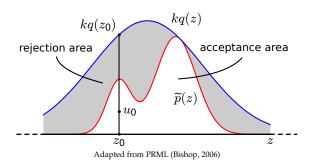
- Assume:
 - Sampling from p(z) is difficult
 - Evaluating $\tilde{p}(z) = Zp(z)$ is easy (and Z may be unknown)
- Find a simpler distribution (proposal distribution) q(z) from which we can easily draw samples (e.g., Gaussian, Laplace)
- ► Find an upper bound $kq(z) \ge \tilde{p}(z)$

Rejection Sampling: Algorithm



- 1. Generate $z_0 \sim q(z)$
- 2. Generate $u_0 \sim \mathcal{U}[0, kq(z_0)]$
- 3. If $u_0 > \tilde{p}(z_0)$, reject the sample. Otherwise, retain z_0

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- Accepted pairs (z, u) are uniformly distributed under the curve of $\tilde{p}(z)$
- Marginal probability density of the *z*-coordiantes of accepted points must be proportional to $\tilde{p}(z)$

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• Samples are independent samples from p(z)

Sampling in High Dimensions

Example:

- $p(x) = \mathcal{N}(\mathbf{0}, \sigma_p^2 \mathbf{I}), \quad q(x) = \mathcal{N}(\mathbf{0}, \sigma_q^2 \mathbf{I}) \text{ where } \sigma_q = 1.01\sigma_p$
- ▶ What is the value of *k* if $x \in \mathbb{R}^{1000}$?

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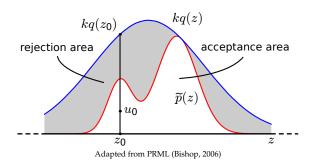
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- ▶ What is the value of *k* if $x \in \mathbb{R}^{1000}$?
- $q(0) = 1/(2\pi\sigma_q^2)^{500}$ For $kq \ge p$ we need to set

$$k \geqslant \frac{p(0)}{q(0)} = \frac{(\sigma_q^2)^{500}}{(\sigma_p^2)^{500}} = \exp\left(1000\ln\frac{\sigma_q}{\sigma_p}\right) = \exp(1000\ln 1.01) \approx 20,000$$

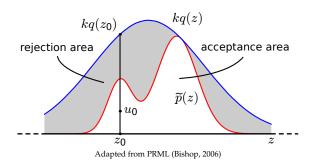
- Acceptance rate is the ratio of the volume under p to the volume under kq. In our example: 1/k = 1/20,000.
- In high dimensions the factor k is probably huge
 Low acceptance rate
- ▶ Finding *k* is tricky

Shortcomings



► Finding the upper bound *k* is tricky

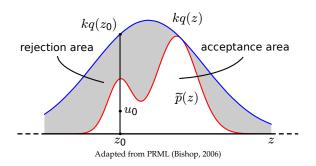
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Shortcomings



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- ► In high dimensions the factor *k* is probably huge
- ► Low acceptance rate/high rejection rate of samples

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Key idea: Do not throw away all rejected samples, but give them lower weight by rewriting the integral as an expectation under a simpler distribution q (proposal distribution):

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If we choose q in a way that we can easily sample from it, we can approximate this last expectation by Monte Carlo:

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- ► Does not scale to interesting (high-dimensional) problems
- ▶ Different approach to sample from complicated (high-dimensional) distributions

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Markov Chain Monte Carlo

Objective

Generate samples from an unknown target distribution.

Target distribution: the distribution we are interested in (e.g., posterior)

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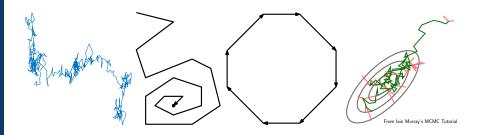
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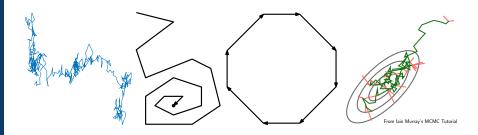
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- Example: $T(x^{(t+1)}|x^{(t)}) = \mathcal{N}(x^{(t+1)}|x^{(t)}, \sigma^2 I)$
- ► Samples $x^{(1)}, ..., x^{(t)}$ form a Markov chain
- Samples $x^{(1)}, \dots, x^{(t)}$ are no longer independent, but unbiased We can still average them

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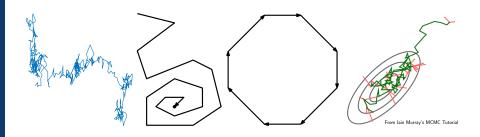
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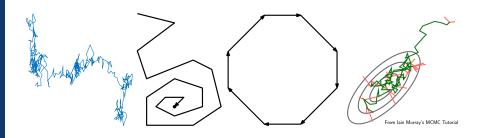
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- ► Converge to an absorbing state
- ► Converge to a (deterministic) limit cycle
- Converge to an equilibrium distribution p^* : Markov chain remains in a region, bouncing around in a random way

- Remember objective: Explore/sample parameters that may have generated our data (generate samples from posterior)
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- ▶ Design the Markov chain such that the equilibrium distribution is the desired distribution p(x)
- Generate a Markov chain that converges to that equilibrium distribution (independent of start state)
- ► Although successive samples are dependent we can effectively generate independent samples by running the Markov chain long enough: Discard most of the samples, retain only every *M*th sample

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▶ Use ergodic and stationary Markov chains to generate samples from the equilibrium distribution

Invariance and Detailed Balance

► Invariance: Each step leaves the distribution p^* invariant (we stay in p^*):

$$p^*(x') = \sum_{x} T(x'|x)p^*(x)$$
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Sufficient condition for p* being invariant:
 Detailed balance:

$$p^*(\mathbf{x})T(\mathbf{x}'|\mathbf{x}) = p^*(\mathbf{x}')T(\mathbf{x}|\mathbf{x}')$$

Also ensures that the Markov chain is reversible

Metropolis-Hastings

- ► Assume that $\tilde{p} = Zp$ can be evaluated easily (in practice: $\log \tilde{p}$)
- ► Proposal density $q(x'|x^{(t)})$ depends on last sample $x^{(t)}$. Example: Gaussian with mean $x^{(t)}$: $q(x'|x^{(t)}) = \mathcal{N}(x^{(t)}, \Sigma)$

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Metropolis-Hastings Algorithm

- 1. Generate proposal $x' \sim q(x'|x^{(t)})$
- 2. If

$$\frac{q(\mathbf{x}^{(t)}|\mathbf{x}')\tilde{p}(\mathbf{x}')}{q(\mathbf{x}'|\mathbf{x}^{(t)})\tilde{p}(\mathbf{x}^{(t)})} \ge u, \qquad u \sim U[0,1]$$

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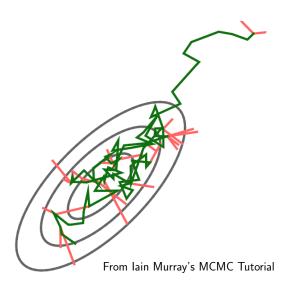
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- $p(x^{(t)}) \xrightarrow{t \to \infty} p^*(x)$ Converge to equilibrium distribution
- ► If proposal distribution is symmetric: Metropolis Algorithm (Metropolis et al., 1953); Otherwise Metropolis-Hastings

Example



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Step-Size Demo

- ► Explore $p(x) = \mathcal{N}(x | 0, 1)$ for different step sizes σ .
- We can only evaluate $\log \tilde{p}(x) = -x^2/2$
- ▶ Proposal distribution *q*: Gaussian $\mathcal{N}(x^{(t+1)} | x^{(t)}, \sigma^2)$ centered at the current state for various step sizes σ
- ► Expect to explore the space between −2,2 with high probability

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- ► Theoretical results: in 1D 44%, in higher dimensions about 25% acceptance rate for good mixing properties
- ► Tune the step size

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- Unlike rejection sampling, the previous sample is used to reset the chain (if a sample was discarded)
- ▶ If q > 0, we will end up in the equilibrium distribution: $p(x^{(t)}) \stackrel{t \to \infty}{\longrightarrow} p^*(x)$

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 - ➤ Adaptive rejection sampling generates independent samples
- Unlike rejection sampling, the previous sample is used to reset the chain (if a sample was discarded)
- ► If q > 0, we will end up in the equilibrium distribution: $p(\mathbf{x}^{(t)}) \stackrel{t \to \infty}{\longrightarrow} p^*(\mathbf{x})$
- ► Explore the state space by random walk
 - ➤ May take a while in high dimensions
- No further catastrophic problems in high dimensions

Gibbs Sampling (Geman & Geman, 1984)

- Assumption: $p(x) = p(x_1, ..., x_n)$ is too complicated to draw samples from directly, but its conditionals $p(x_i|x_{\setminus i})$ are tractable to work with
- ► Any distribution "with a name" (Gaussian, Laplace, Bernoulli, Gamma, Wishart, ...) is easy to sample from (standard libraries)

Algorithm

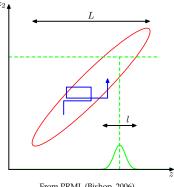
Assuming *n* parameters x_1, \ldots, x_n , Gibbs sampling samples individual variables conditioned on all others:

1.
$$x_1^{(t+1)} \sim p(x_1|x_2^{(t)}, \dots, x_n^{(t)})$$

2.
$$x_2^{(t+1)} \sim p(x_2|x_1^{(t+1)}, x_3^{(t)}, \dots, x_n^{(t)})$$

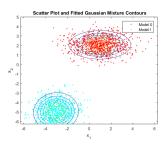
3. :

4.
$$x_n^{(t+1)} \sim p(x_n|x_1^{(t+1)},\ldots,x_{n-1}^{(t+1)})$$



From PRML (Bishop, 2006)

Gibbs Sampling: Ergodicity



- p(x) is invariant
- Ergodicity: Sufficient to show that all conditionals are greater than 0.
 - Then any point in *x*-space can be reached from any other point (potentially with low probability) in a finite number of steps involving one update of each of the component variables.

Finding the Conditionals

- 1. Write down the (log-) joint distribution $p(x_1, ..., x_n)$
- 2. For each x_i
 - 2.1 Throw away all terms that do not depend on the current sampling variable
 - 2.2 Pretend this is the density for your variable of interest and all other variables are fixed. What distribution does the log-density remind you of?
 - 2.3 That is your conditional sampling density $p(x_i|x_{\setminus i})$

Example

► Model:

$$y_i \sim \mathcal{N}(\mu, \tau^{-1})$$
, $\mu \sim \mathcal{N}(\mu \mid 0, 1)$, $\tau \sim \text{Gamma}(\tau \mid 2, 1)$
 $\text{Gamma}(\tau \mid 2, 1) = \frac{1}{\Gamma(2)} \tau \exp(-\tau)$

▶ **Objective:** Generate samples from the parameter posterior $p(\mu, \tau | y_1, ..., y_N)$ given N observations $y_1, ..., y_N$

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- ▶ Then

$$p(y, \mu, \tau) = \prod_{i=1}^{N} p(y_i | \mu, \tau) p(\mu) p(\tau)$$
$$\propto \tau^{N/2} \exp(-\frac{\tau}{2} \sum_{i} (y_i - \mu)^2) \exp(-\frac{\tau}{2} \mu^2) \tau \exp(-\tau)$$

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$$\begin{split} p(\boldsymbol{y}, \mu, \tau) &= \prod_{i=1}^{N} p(y_i | \mu, \tau) p(\mu) p(\tau) \\ &\propto \tau^{N/2} \exp(-\frac{\tau}{2} \sum_{i} (y_i - \mu)^2) \exp(-\frac{1}{2} \mu^2) \tau \exp(-\tau) \\ p(\mu | \tau, \boldsymbol{y}) &= \mathcal{N}\left(\frac{\tau \sum_{i} y_i}{1 + N \tau}, (1 + N \tau)^{-1}\right) \\ p(\tau | \mu, \boldsymbol{y}) &= \operatorname{Gamma}(2 + \frac{N}{2}, 1 + \frac{1}{2} \sum_{i} (y_i - \mu)^2) \end{split}$$

- Gibbs is Metropolis-Hastings with acceptance probability 1:
 Sequence of proposal distributions q is defined in terms of conditional distributions of the joint p(x)
 - **▶** Converge to equilibrium distribution: $p^{(t)}(x) \stackrel{t \to \infty}{\longrightarrow} p(x)$
 - >> Exploration by random walk behavior can be slow

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- Statistical software derives the conditionals of the model, and it works out how to do the updates: STAN¹, WinBUGS², JAGS³

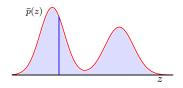
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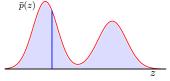
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Flavors of Gibbs Sampling

- Collapsed Gibbs sampler: Analytically integrate out some parameters and sample the rest.
 - Tends to be much more efficient with smaller variance (see Rao-Blackwellization in the state estimation literature)
- Block-Gibbs sampler: Sample groups of variables at a time instead of single-site updating



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- ► Introduce additional variable u, define joint $\hat{p}(x, u)$:

$$\hat{p}(x,u) = \begin{cases} 1/Z_p & \text{if } 0 \leq u \leq \tilde{p}(x) \\ 0 & \text{otherwise} \end{cases}$$
, $Z_p = \int \tilde{p}(x)dx$

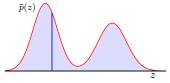


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$$\int \hat{p}(x,u)du = \int_0^{\tilde{p}(x)} 1/Z_p du = \tilde{p}(x)/Z_p = p(x)$$



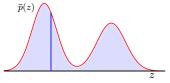
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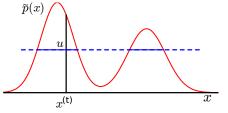
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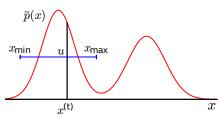
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- **▶** Obtain samples from unknown p(x) by sampling from $\hat{p}(x, u)$ and then ignore u values
- ► Gibbs sampling: Update one variable at a time

Slice Sampling Algorithm

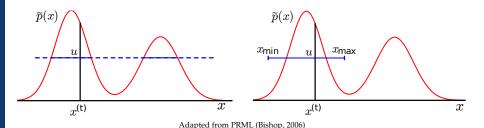




Adapted from PRML (Bishop, 2006)

- ► Repeat the following steps:
 - 1. Draw $u|x^{(t)} \sim \mathcal{U}[0, \tilde{p}(x)]$
 - 2. Draw $x^{(t+1)}|u \sim \mathcal{U}[\{x : \tilde{p}(x) > u\}]$ \Longrightarrow slice

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- ▶ In practice, we sample $x^{(t+1)}|u$ uniformly from an interval $[x_{\min}, x_{\max}]$ around $x^{(t)}$.
- ► The interval is found adaptively (see Neal (2003) for details)

Relation to other Sampling Methods

Similar to:

- ▶ Metropolis: Just need to be able to evaluate $\tilde{p}(x)$ More robust to the choice of parameters (e.g., step size is automatically adapted)
- ► Gibbs: 1-dimensional transitions in state space
 No longer required that we can easily sample from 1-D
 conditionals
- Rejection: Asymptotically draw samples from the volume under the curve described by p

 No upper-bounding of p required

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Properties

- Slice sampling can be applied to multivariate distributions by repeatedly sampling each variable/dimension in turn (similar to Gibbs sampling).
 - See (Neal, 2003; Murray et al., 2010) for more details
- ▶ This requires to compute a function that is proportional to $p(x_i|x_{\setminus i})$ for all variables x_i .

Properties

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 - See (Neal, 2003; Murray et al., 2010) for more details
- ► This requires to compute a function that is proportional to $p(x_i|x_{\setminus i})$ for all variables x_i .
- ▶ No rejections
- Adaptive step sizes
- Easy to implement
- Broadly applicable

 Samples from the Markov chain before the equilibrium distribution is reached should be discarded (burn-in phase)

40

- Samples from the Markov chain before the equilibrium distribution is reached should be discarded (burn-in phase)
- ► MCMC generates dependent samples
 - >> Introduces additional variance in the Monte-Carlo estimator

$$\frac{1}{S} \sum_{s=1}^{S} f(x^{(s)}), \quad x^{(s)} \sim p(x)$$

due to correlation of samples

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- ► Autocorrelation is an indicator for choosing *K*

MCMC Diagnostics: Trace Plots

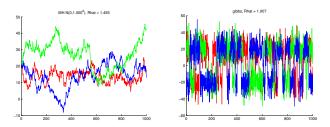


Figure from Murphy (2012)

- Mixing time: Amount of time it takes the Markov chain to converge to the stationary distribution and forget its initial state.
- ► Trace plots: Run multiple chains from very different starting points, plot the samples of the variables of interest. If the chain has mixed, the trace plots should converge to the same distribution.

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Summary

- Solving integrals, computing expectations
- Monte Carlo methods use random numbers
- Rejection and importance sampling do not work well in high dimensions
- MCMC generates a Markov chain of dependent samples that allow us to generate samples from the target distribution
- ► Metropolis Hastings, Gibbs, Slice sampling

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