

Markov chain Monte Carlo

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Outline

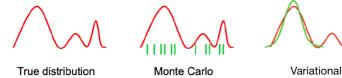


- Introduction
- Rejection sampling
- Importance sampling
- Metropolis-Hastings
- Hamiltonian Monte Carlo
- Gibbs sampling

Previously, in MBML...



- Representation (Weeks 1-4)
- Modelling toolbox (Weeks 5-8, +12)
- Inference (Weeks 9-11)
 - Exact inference
 - Analytical (Bayes' rule)
 - Variable elimination
 - Belief propagation
 - Approximate inference
 - Stochastic methods Markov chain Monte Carlo (MCMC)
 - Deterministic methods Variational inference (VI)



Approximate inference



- Suppose we wish to compute the posterior distribution of the latent variables ${\bf z}$ given some observed data ${\bf x}$: $p({\bf z}|{\bf x})$
- So far, we discussed various exact algorithms for posterior inference
- But, for many problems of interest exact posterior inference is intractable
 - Cannot determine the posterior distribution analytically
 - Cannot even compute expectations with respect to the posterior
- Reasons for intractability:
 - Dimensionality of the latent space is too high to work with directly
 - Posterior distribution has a highly complex form for which expectations are not analytically tractable
- We must resort to approximate inference methods!

Monte Carlo methods



• In most situations, the posterior distribution $p(\mathbf{z}|\mathbf{x})$ is required only for evaluating **expectations** of some function $f(\mathbf{z})$ (e.g. to make predictions)

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

- (Note: we are using **z** to denote the variables whose posterior we wish to infer!)
- For example, the mean of z is given by

$$\mathbb{E}_{p(\mathbf{z}|\mathbf{x})}[\mathbf{z}] = \int \mathbf{z} \, p(\mathbf{z}|\mathbf{x}) \, d\mathbf{z}$$

- The idea behind sampling methods is to obtain a set of samples $\mathbf{z}^{(s)}$, for $s \in \{1, \dots, S\}$, drawn independently from the distribution $p(\mathbf{z}|\mathbf{x})$
- Allows to approximate expectations as finite sums!

$$\mathbb{E}_{p(\mathbf{z}|\mathbf{x})}[f(\mathbf{z})] \approx \frac{1}{S} \sum_{s=1}^{S} f(\mathbf{z}^{(s)})$$

Example: making predictions



- Let **z** denote the parameters of our model (e.g. Bayesian logistic regression)
- ullet Suppose we wish to make a prediction y_* for a new observation ${f x}_*$

$$\begin{split} \mathbb{E}_{p(\mathbf{z}|\mathbf{x})}[y_*|\mathbf{x}_*] &= \int p(y_*|\mathbf{x}_*, \mathbf{z}, \mathbf{x}) \, p(\mathbf{z}|\mathbf{x}) \, d\mathbf{z} \\ &\approx \frac{1}{S} \sum_{s=1}^S p(y_*|\mathbf{x}_*, \mathbf{z}^{(s)}, \mathbf{x}) \end{split}$$

where $\mathbf{z}^{(s)} \sim p(\mathbf{z}|\mathbf{x})$

Notice that

$$\frac{1}{S} \sum_{s=1}^{S} p(y_* | \mathbf{x}_*, \mathbf{z}^{(s)}, \mathbf{x}) \neq p\left(y_* \middle| \mathbf{x}_*, \frac{1}{S} \sum_{s=1}^{S} \mathbf{z}^{(s)}, \mathbf{x}\right)$$

- Even if the posterior $p(\mathbf{z}|\mathbf{x})$ is intractable to compute, we can still make predictions as long as we can obtain samples from the posterior!
- But if $p(\mathbf{z}|\mathbf{x})$ is intractable, how can we sample from it?

Properties of Monte Carlo



Monte Carlo estimator:

$$\begin{split} \mathbb{E}_{p(\mathbf{z}|\mathbf{x})}[f(\mathbf{z})] &= \int f(\mathbf{z}) \, p(\mathbf{z}|\mathbf{x}) \, d\mathbf{z} \\ &\approx \hat{f} \triangleq \frac{1}{S} \sum_{s=1}^{S} f(\mathbf{z}^{(s)}), \quad \mathbf{z}^{(s)} \sim p(\mathbf{z}|\mathbf{x}) \end{split}$$

Estimator is unbiased:

$$\mathbb{E}_{p(\{\mathbf{z}^{(s)}\}|\mathbf{x})}[\hat{f}] = \frac{1}{S} \sum_{s=1}^{S} \mathbb{E}_{p(\mathbf{z}|\mathbf{x})}[f(\mathbf{z})] = \mathbb{E}_{p(\mathbf{z}|\mathbf{x})}[f(\mathbf{z})]$$

• Variance shrinks proportionally to 1/S:

$$\mathbb{V}_{p(\{\mathbf{z}^{(s)}\}|\mathbf{x})}[\hat{f}] = \frac{1}{S^2} \sum_{s=1}^{S} \mathbb{V}_{p(\mathbf{z}|\mathbf{x})}[f(\mathbf{z})] = \frac{1}{S} \mathbb{V}_{p(\mathbf{z}|\mathbf{x})}[f(\mathbf{z})]$$

Sampling from distributions



- Draw mass to the left of point: $u \sim \mathsf{Uniform}(0,1)$
- ullet Use inverse CDF to obtain sample $y(u)=h^{-1}(u)$

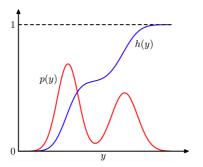


Figure: from PRML book, Bishop (2006)

 \bullet However, keep in mind that we can't always compute and invert $h(\boldsymbol{y})$

Rejection sampling



- ullet Suppose we wish to sample from the unnormalized density $ilde{p}(z) \propto p(z)$
- Construct **proposal distribution** $kq(z) \ge \tilde{p}(z)$

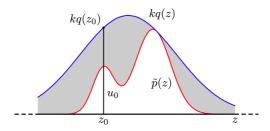


Figure: from PRML book, Bishop (2006)

- Draw sample $z \sim q(z)$
- Draw height $u \sim \mathsf{Uniform}(0, kq(z))$
- **Reject** sample z if $u > \tilde{p}(z)$

Playtime!



- Rejection sampling
 - See "Part 1" of "11 Markov chain Monte Carlo methods.ipynb" notebook
 - Expected duration: 30 minutes



• As previously mentioned, the posterior distribution p(z) is often required only for evaluating **expectations** of some function f(z)

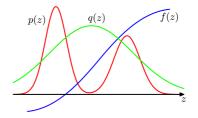


Figure: from PRML book, Bishop (2006)

- Importance sampling allows for just that!
- It does **not** provide a mechanism for drawing samples from the posterior p(z), but it allows to compute expectations with respect to it:

$$\mathbb{E}_{p(z)}[f(z)] \approx \frac{1}{S} \sum_{s=1}^{S} f(z^{(s)}), \quad z^{(s)} \sim p(z)$$



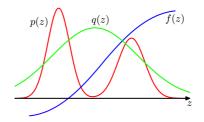


Figure: from PRML book, Bishop (2006)

Rewrite the integral as an expectation under q:

$$\mathbb{E}_{p(z)}[f(z)] = \int f(z) \, p(z) \, dz = \int f(z) \, \frac{p(z)}{q(z)} \, q(z) \, dz$$

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

- The ratio $\frac{p(z^{(s)})}{q(z^{(s)})}$ is the **importance/weights** of each sample!
- But, what if we cannot evaluate p(z) but only $\tilde{p}(z) \propto p(z)$?



- But, what if we cannot evaluate p(z) but only $\tilde{p}(z) \propto p(z)$?
- Let $p(z) = \frac{1}{Z_p} \tilde{p}(z)$ and $q(z) = \frac{1}{Z_q} \tilde{q}(z)$

$$\mathbb{E}_{p(z)}[f(z)] = \int f(z) \, p(z) \, dz \approx \frac{Z_q}{Z_p} \frac{1}{S} \sum_{s=1}^S f(z^{(s)}) \underbrace{\frac{\tilde{p}(z^{(s)})}{\tilde{q}(z^{(s)})}}_{z^{(s)}}, \quad z^{(s)} \sim q(z)$$



- But, what if we cannot evaluate p(z) but only $\tilde{p}(z) \propto p(z)$?
- \bullet Let $p(z)=\frac{1}{Z_p}\tilde{p}(z)$ and $q(z)=\frac{1}{Z_q}\tilde{q}(z)$

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$$\approx \frac{1}{S} \sum_{s=1}^S f(z^{(s)}) \underbrace{\frac{\tilde{p}(z^{(s)})}{\tilde{p}(z^{(s)})}}_{w^{(s)}} = \sum_{s=1}^S f(z^{(s)}) \, w^{(s)}$$

since
$$\frac{Z_p}{Z_q} \approx \frac{1}{S} \sum_s \tilde{r}^{(s)}$$
.

• All we need to do is compute the importance/weight $w^{(s)}$ of each sample s as the normalized ratio $\tilde{r}^{(s)}=\frac{\tilde{p}(z^{(s)})}{\tilde{q}(z^{(s)})}$

Playtime!



- Importance sampling
 - See "Part 2' of "11 Markov chain Monte Carlo methods.ipynb" notebook
 - Expected duration: 30 minutes



- Rejection and importance sampling scale badly with dimensionality
- Also, the whole procedure still feels naive... A 2-dimensional example:

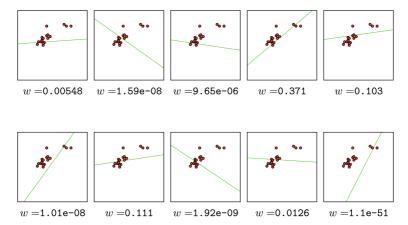


Figure: Samples from posterior over linear regression coefficients $\beta = (\beta_0, \beta_1)^T$ using importance sampling. Figure from Iain Murray (2009)

Metropolis algorithm



- Start with some initial state **z** and iterate the following steps:
- Perturb variables using proposal distribution $q(\mathbf{z}'|\mathbf{z})$, e.g. $q(\mathbf{z}'|\mathbf{z}) = \mathcal{N}(\mathbf{z}'|\mathbf{z}, \sigma^2\mathbf{I})$
- Accept proposal with probability: $\min\left(1, \frac{\tilde{p}(\mathbf{z}')}{\tilde{p}(\mathbf{z})}\right)$
- In other words: always accept if new state \mathbf{z}' has higher probability, otherwise go to new state \mathbf{z}' anyway with probability $\frac{\tilde{p}(\mathbf{z}')}{\tilde{p}(\mathbf{z})}$

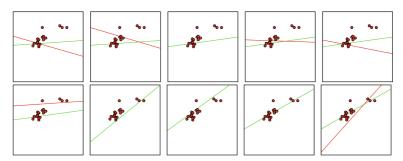


Figure: Samples from posterior over linear regression coefficients $\beta = (\beta_0, \beta_1)^T$ using the Metropolis algorithm. Figure from Iain Murray (2009)

Metropolis algorithm



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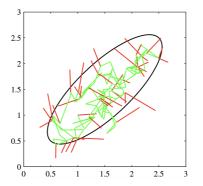


Figure: from PRML book, Bishop (2006)

• Green = accepted steps, Red = rejected steps

Markov chain Monte Carlo



- Construct a **biased** random walk that explores the posterior distribution $\tilde{p}(\mathbf{z})$
- Markov steps at each time step t: $\mathbf{z}^{(t)} \sim p(\mathbf{z}^{(t)}|\mathbf{z}^{(t-1)})$

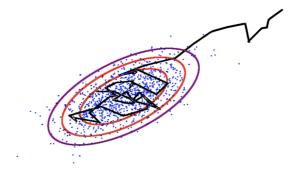


Figure: from Iain Murray (2009)

- MCMC gives approximate (but **correlated!**) samples from $\tilde{p}(\mathbf{z})$
- Can we choose any transition operator $T(\mathbf{z}^{(t)} \leftarrow \mathbf{z}^{(t-1)}) \triangleq p(\mathbf{z}^{(t)}|\mathbf{z}^{(t-1)})$?

Some definitions...



- ullet Consider a Markov chain with transition probabilities: $T_t(\mathbf{z}^{(t)} \leftarrow \mathbf{z}^{(t-1)})$
- ullet A Markov chain is called **homogeneous** if the transition probabilities are the same for all t
- A distribution $p^*(\mathbf{z})$ is said to be **invariant**, or **stationary**, with respect to a Markov chain if each step in the chain leaves that distribution invariant
- For a homogeneous Markov chain with transition probabilities $T(\mathbf{z}' \leftarrow \mathbf{z})$, the distribution $p^*(\mathbf{z})$ is invariant if

$$p^*(\mathbf{z}) = \sum_{\mathbf{z}'} T(\mathbf{z} \leftarrow \mathbf{z}') \, p^*(\mathbf{z}')$$

- Our goal is to use Markov chains to sample from a given distribution
- We can achieve this if we set up a Markov chain such that the desired distribution is invariant!

MCMC properties



• A sufficient (but not necessary) condition for ensuring that the required distribution $p(\mathbf{z})$ is invariant is to choose the transition probabilities to satisfy the property of **detailed balance**, defined by

$$p^*(\mathbf{z}') T(\mathbf{z} \leftarrow \mathbf{z}') = p^*(\mathbf{z}) T(\mathbf{z}' \leftarrow \mathbf{z})$$

- A Markov chain that respects detailed balance is said to be reversible
- Lastly, we must also require that for $t \to \infty$, $p(\mathbf{z}^{(t)})$ converges to the required invariant distribution $p^*(\mathbf{z})$, irrespective of the choice of initial distribution $p(\mathbf{z}^{(0)})$
- This property is called **ergodicity**
- The invariant distribution is then called the equilibrium distribution



- A more general version of the Metropolis algorithm:
 - ullet Propose a move from the current state: $\mathbf{z}' \sim q(\mathbf{z}'|\mathbf{z})$
 - Accept proposal with probability:

$$\min\left(1,\frac{\tilde{p}(\mathbf{z}')\,q(\mathbf{z}|\mathbf{z}')}{\tilde{p}(\mathbf{z})\,q(\mathbf{z}'|\mathbf{z})}\right)$$

- Satisfies detailed balance! (see Bishop (2006) for proof)
- Since the Gaussian is a symmetric function, i.e. $\mathcal{N}(\mathbf{z}'|\mathbf{z},\sigma^2) = \mathcal{N}(\mathbf{z}|\mathbf{z}',\sigma^2)$, when $q(\mathbf{z}'|\mathbf{z}) = \mathcal{N}(\mathbf{z}'|\mathbf{z},\sigma^2\mathbf{I})$ we get back the Metropolis algorithm:

$$\min\left(1, \frac{\tilde{p}(\mathbf{z}') \, q(\mathbf{z} \!\!\!\!\! + \!\!\!\! \mathbf{z}')}{\tilde{p}(\mathbf{z}) \, q(\mathbf{z}' \!\!\!\!\! + \!\!\!\! \mathbf{z})}\right) = \min\left(1, \frac{\tilde{p}(\mathbf{z}')}{\tilde{p}(\mathbf{z})}\right)$$



ullet Be careful how you choose the step size σ^2

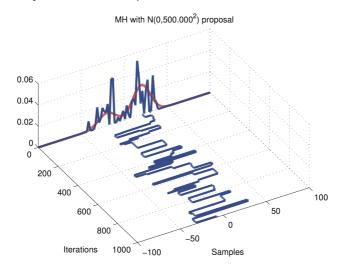


Figure: from ML: A prob. perspective book, Murphy (2011)



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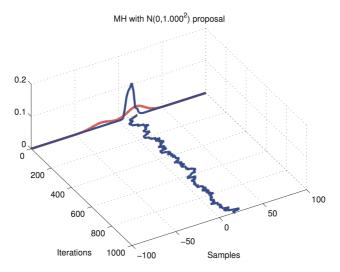


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• Careful how you choose the step size σ^2

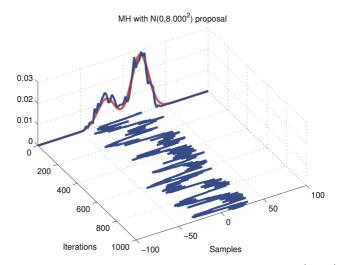


Figure: from ML: A prob. perspective book, Murphy (2011)

MCMC in practice



- ullet The samples are correlated! We should ${f thin}$ only keep every n^{th} sample
- Arbitrary initialization means early samples are bad discard a burn-in period
- A good idea is to run multiple chains (e.g. in parallel multicore)
- How do you know when to stop?
- Make sure to check diagnostics!
 (e.g. "Rhat" and number of effective samples in STAN)

Playtime!

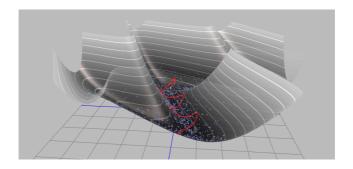


- Metropolis-Hastings
 - See "Part 3' of "11 Markov chain Monte Carlo methods.ipynb" notebook
 - Expected duration: 30 minutes

Hamiltonian Monte Carlo: basic idea



- Include Hamiltonian dynamics
 - Construct a landscape with gravitational potential energy $E(\mathbf{z}) = -\log \, p^*(\mathbf{z})$
 - ullet Introduce velocity v (auxiliary variable) carrying kinetic energy K(v)
- Define joint distribution: $p(\mathbf{z}, v) \propto e^{-E(\mathbf{z})} e^{-K(v)} = e^{-H(\mathbf{z}, v)}$
- ullet Velocity v is independent of position and Gaussian distributed



Hamiltonian Monte Carlo: basic idea



- Procedure:
 - Gibbs sample velocity v (we will see Gibbs sampling next!)
 - Simulate Hamiltonian dynamics then flip sign of velocity
 - Accept new position with probability: $\min[1, e^{H(\mathbf{z}, v) H(\mathbf{z}', v')}]$
- Hamiltonian dynamics are simulated approximately
- Distances between successive generated points are typically large, so we need less iterations to get representative sampling!
- Each iteration is more computationally expensive, but sampling is more efficient
- For a detailed explanation, see: http://arogozhnikov.github.io/2016/12/19/markov_chain_monte_carlo.html

Hamiltonian Monte Carlo: basic idea



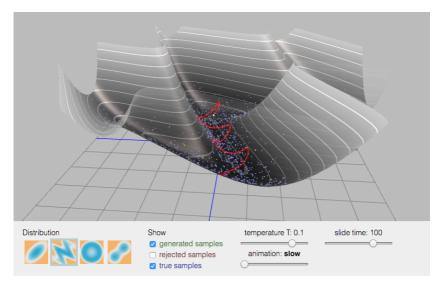
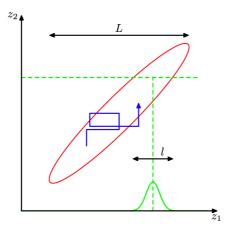


Figure: http://arogozhnikov.github.io/2016/12/19/markov_chain_monte_carlo.html

Gibbs sampling





- Suppose we wish to sample from the posterior $p(\mathbf{z}) = p(z_1, \dots, z_M)$
- Start with initial state $\mathbf{z}^{(0)}$
- Pick each variable z_m in turn (or randomly) and resample it from the conditional:

$$z_m \sim p(z_m|\mathbf{z}_{i\neq m})$$

Gibbs sampling



- Simple and widely applicable MCMC algorithm
- Gibbs sampling has no rejection! (see next slide)
- Must be able to sample from conditional distributions $p(z_m|\mathbf{z}_{i\neq m})$
 - Discrete conditionals with a few possible settings can be explicitly normalized

$$p(z_m|\mathbf{z}_{i\neq m}) = \frac{p(z_m, \mathbf{z}_{i\neq m})}{\sum_{z_j} p(z_j|\mathbf{z}_{i\neq m})}$$

- Continuous conditionals require either:
 - Conjugacy to allow for analytical solutions
 - Or efficient approximate solutions (e.g. standard sampling methods)

Gibbs sampling



- Gibbs sampling satisfies detailed balance (see e.g. Bishop (2006) for proof)
- It is a particular instance of the Metropolis-Hastings algorithm where $q(\mathbf{z}'|\mathbf{z}) = p(z_m'|\mathbf{z}_{i \neq m})$
- Probability of acceptance:

$$\min\left(1, \frac{p(\mathbf{z}')\,q(\mathbf{z}|\mathbf{z}')}{p(\mathbf{z})\,q(\mathbf{z}'|\mathbf{z})}\right) = \min\left(1, \underbrace{\frac{p(z_m'|\mathbf{z}_{i\neq m}')\,p(\mathbf{z}_{i\neq m}')\,p(\mathbf{z}_{i\neq m}')\,p(z_m'|\mathbf{z}_{i\neq m}')}_{=1}}\right) = 1$$

where we used the fact that $p(\mathbf{z}) = p(z_m | \mathbf{z}_{i \neq m}) p(\mathbf{z}_{i \neq m})$ and $\mathbf{z}'_{i \neq m} = \mathbf{z}_{i \neq m}$

- Gibbs sampling has no rejection!
- A practical example (discrete variables): see slides 18-25 from David Sontag (http://people.csail.mit.edu/dsontag/courses/pgm13/slides/lecture9.pdf)

Summary



- We need approximate methods to compute posteriors (sums/integrals)
- Monte Carlo does not explicitly depend on dimension, but simple methods work only in low dimensions
- MCMC methods can make local moves important for higher dimensions!
- General and often easy to implement (simple computations), but hard to diagnose

Playtime!



- Gibbs sampling
 - See "Part 4' of "11 Markov chain Monte Carlo methods.ipynb" notebook
 - Expected duration: 30 minutes