

Statistical Learning for Text Data Analytics

Latent Dirichlet Allocation

Yangqiu Song

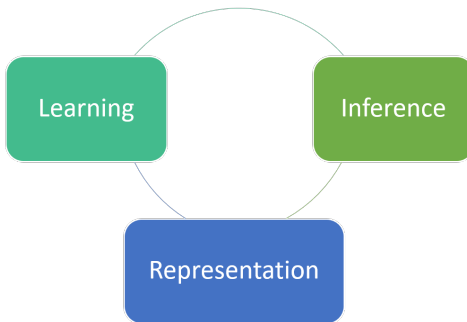
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*Contents are based on materials created by David Mackay (Introduction to Monte Carlo methods, 1998)

Course Topics



- Representation: language models, word embeddings, **topic models**
- Learning: supervised learning, semi-supervised learning, sequence models, deep learning, optimization techniques
- Inference: constraint modeling, joint inference, search algorithms

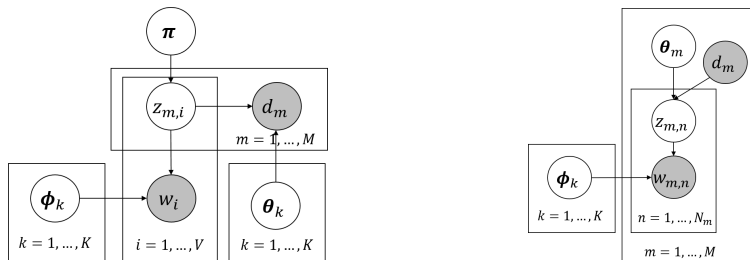
NLP applications: tasks introduced in Lecture 1

Overview

- 1 Language Models: Recap
- 2 Topic Models
- 3 Probabilistic Latent Semantic Analysis (PLSA)
- 4 Latent Dirichlet Allocation (LDA)
 - Motivation: Bayesian Modeling
 - Background of Monte Carlo Methods
 - Importance Sampling
 - Rejection Sampling
 - Metropolis Methods
 - Gibbs Sampling
 - Sampling for EM Algorithm
 - Collapsed Gibbs Sampling for LDA

Alternative Way for PLSA to Generate Texts

$$\begin{aligned}
 P(\mathcal{D}, \mathcal{W}) &= \prod_{m=1}^M \prod_{i=1}^{N_m} \sum_{k=1}^K P(z_{m,i} = k) P(d_m | \theta_k) P(w_i | \phi_k) \\
 &= \prod_{m=1}^M \prod_{i=1}^V \left(\sum_{k=1}^K P(z_{m,i} = k) P(d_m | \theta_k) P(w_i | \phi_k) \right)^{c_{d_m}(w_i)}
 \end{aligned}$$



$$P(\mathcal{D}, \mathcal{W}) = \prod_{m=1}^M \prod_{i=1}^V P(d_m) \left(\sum_{k=1}^K P(z_{m,i} = k | \theta_m) P(w_i | \phi_k) \right)^{c_{d_m}(w_i)}$$

Bayesian Modeling: Topic Models

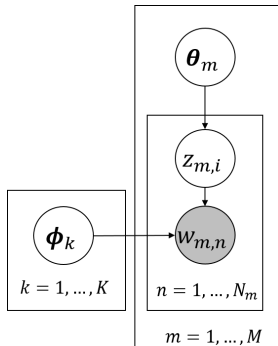


Figure: PLSA

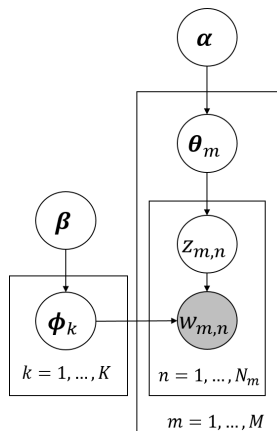


Figure: LDA

Generative Process of Latent Dirichlet Allocation

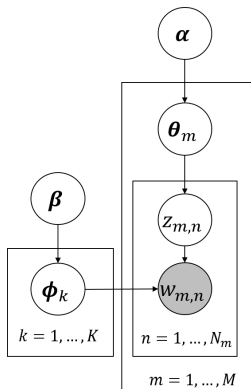


Figure: LDA

- For all clusters/components $k \in [1, K]$:
 - Choose mixture components $\phi_k \sim \text{Dir}(\phi|\beta)$
- For all documents $m \in [1, M]$:
 - Choose $N_m \sim \text{Poisson}(\xi)$
 - Choose mixture probability $\theta_m \sim \text{Dir}(\theta|\alpha)$
 - For all words $n \in [1, N_m]$ in document d_m :
 - Choose a component index
 $z_{m,n} \sim \text{Mult}(z|\theta_m)$
 - Choose a word $w_{m,n} \sim \text{Mult}(w|\phi_{z_{m,n}})$

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Bayesian Inference

- Suppose we have a Bayesian learning problem
 $P(\Theta|X) \propto P(X|\Theta)P(\Theta)$ ($X = \{x_1, \dots, x_N\}$)
- If we want to predict for a new coming data x
- Maximum a posterior (MAP) makes a point estimation
 $\Theta^* = \max_{\Theta} P(\Theta|X)$, and makes a prediction as $P(x|\Theta^*)$
- Full Bayesian uses $P(x|X) = \int_{\Theta} P(x|\Theta)P(\Theta|X)d\Theta$
- In general, we have a lot of following cases need to be estimated:

$$\mathbb{E}_{P(\mathbf{x})}[\phi(\mathbf{x})] = \int_{\mathbf{x}} \phi(\mathbf{x})P(\mathbf{x})d\mathbf{x}$$

- One way to solve this (especially when $P(\mathbf{x})$ is difficult to compute) is using sampling:

$$\hat{\mathbb{E}}_{P(\mathbf{x})}[\phi(\mathbf{x})] = \frac{1}{R} \sum_{\mathbf{x}^{(r)} \sim P(\mathbf{x})} \phi(\mathbf{x}^{(r)})$$

Sampling

$$\Phi = \mathbb{E}_{P(\mathbf{x})}[\phi(\mathbf{x})] = \int_{\mathbf{x}} \phi(\mathbf{x}) P(\mathbf{x}) d\mathbf{x}$$

- We call $P(\mathbf{x})$ the target density
- We assume $\mathbf{x}$ is a $\mathbb{R}^d$ vector with real/discrete components $\mathbf{x}^i$
- We concentrate on the sampling problem, because if we have solved it, then we can solve the expectation problem by

$$\hat{\Phi} = \hat{\mathbb{E}}_{P(\mathbf{x})}[\phi(\mathbf{x})] = \frac{1}{R} \sum_{\mathbf{x}^{(r)} \sim P(\mathbf{x})}^R \phi(\mathbf{x}^{(r)})$$

- The expectation of $\hat{\Phi}$ is Φ
- The variance of $\hat{\Phi}$ will decrease as $\frac{\sigma^2}{R}$ where σ^2 is the variance of Φ :

$$\sigma^2 = \int_{\mathbf{x}} [\phi(\mathbf{x}) - \Phi]^2 P(\mathbf{x}) d\mathbf{x}$$

which means the accuracy of the sampling is independent of the dimensionality of the space sampled

- A few as a dozen independent samples $\mathbf{x}^{(r)}$ suffice to estimate Φ satisfactorily

However, why is sampling from $P(\mathbf{x})$ hard?

- We assume that the density from which we wish to draw samples $P(\mathbf{x})$ can be evaluated, at least to within a multiplicative constant:

$$P(\mathbf{x}) = P^*(\mathbf{x})/Z$$

- If we can evaluate $P^*(\mathbf{x})$, why can we not easily obtain Φ ?
 - We do not know the normalizing constant

$$Z = \int_{\mathbf{x}} d\mathbf{x} P^*(\mathbf{x})$$

- Even if we know Z , drawing samples from $P(\mathbf{x})$ is still challenging, especially in high-dimensional spaces

One Dimensional Sampling Example

- Consider $P^* = \exp\{0.4(x - 0.4)^2 - 0.08x^4\}$, $x \in (-\infty, \infty)$
- To give a simpler problem, we can discretize the variable x and ask for samples from the discrete prob.

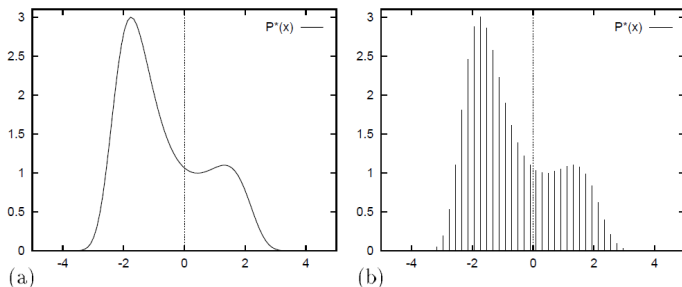
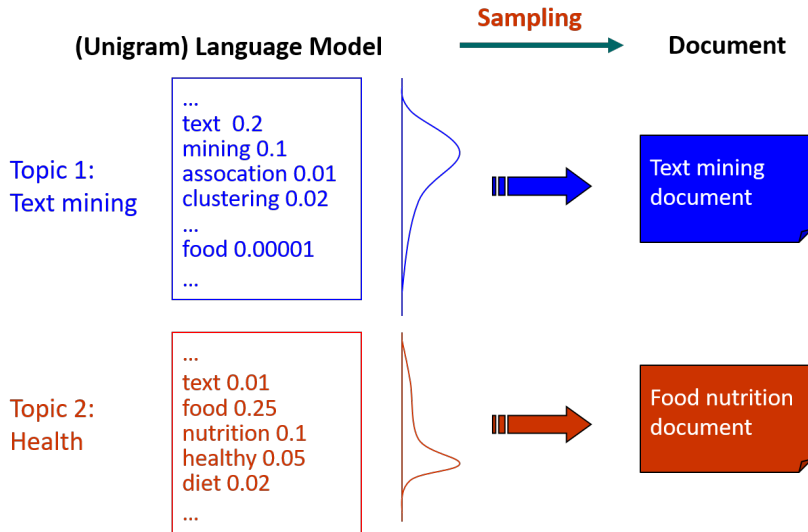


Figure 1. (a) The function $P^*(x) = \exp[0.4(x - 0.4)^2 - 0.08x^4]$. How to draw samples from this density? (b) The function $P^*(x)$ evaluated at a discrete set of uniformly spaced points $\{x_i\}$. How to draw samples from this discrete distribution?

- If we evaluate $p_i^* = P^*(x_i)$ at each point x_i , we can compute $Z = \sum_i p_i^*$ and $p_i = p_i^*/Z$

Recall: Generative View of Text Documents



Generating Text from Language Models

Example

$$P(\text{of}) = 3/66$$

$$P(\text{her}) = 2/66$$

$$P(\text{Alice}) = 2/66$$

$$P(\text{sister}) = 2/66$$

$$P(\text{was}) = 2/66$$

$$P(,) = 4/66$$

$$P(\text{to}) = 2/66$$

$$P(') = 4/66$$

Under a unigram language model:



Alice was beginning to get very tired of sitting by her sister on the bank, and of having nothing to do: once or twice she had peeped into the book her sister was reading, but it had no pictures or conversations in it, 'and what is the use of a book,' thought Alice 'without pictures or conversation?'

Generating Text from Language Models

Example

$P(\text{of}) = 3/66$	$P(\text{her}) = 2/66$
$P(\text{Alice}) = 2/66$	$P(\text{sister}) = 2/66$
$P(\text{was}) = 2/66$	$P(,) = 4/66$
$P(\text{to}) = 2/66$	$P(') = 4/66$

Under a unigram language model:



The same likelihood!

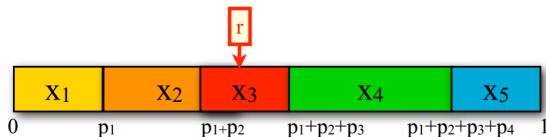
beginning by, very Alice but was and?
reading no tired of to into sitting
sister the, bank, and thought of without
her nothing: having conversations Alice
once do or on she it get the book her had
peeped was conversation it pictures or
sister in, 'what is the use had twice of
a book 'pictures or' to

Recall: Computer Simulation

Sample from a discrete distribution $P(X)$, assuming n outcomes in the event space X

Algorithm 1 Sample from a distribution $P(X)$

- 1: **for** $t = 1$ to T **do**
 - 2: Divide the interval $[0, 1]$ into n intervals according to the probabilities of the outcomes
 - 3: Generate a random number r between 0 and 1
 - 4: Return x_i where r falls into $[\sum_0^{i-1} p_i, \sum_0^i p_i]$
 - 5: **end for**
-



The Cost of Computing Z

- To compute Z , we have to visit every point in the space

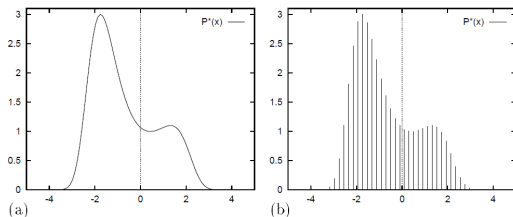


Figure 1. (a) The function $P^*(x) = \exp[0.4(x-0.4)^2 - 0.08x^4]$. How to draw samples from this density? (b) The function $P^*(x)$ evaluated at a discrete set of uniformly spaced points $\{x_i\}$. How to draw samples from this discrete distribution?

- There are 50 uniformly spaced points in one dimension
- If our system had N dimensions, e.g., $K = 2$ clusters (latent variables), $N = 1000$ data examples
 - Then the corresponding number of points would be 2^{1000}

Uniform Sampling

- Having agreed that we cannot visit every location in the space, we might consider trying to solve sampling by uniform sampling:
 - Sample $\mathbf{x}^{(r)}$ uniformly and evaluate $P^*(\mathbf{x}^{(r)})$ to give

$$Z_R = \sum_{r=1}^R P^*(\mathbf{x}^{(r)})$$

- and estimate $\Phi = \mathbb{E}_{P(\mathbf{x})}[\phi(\mathbf{x})] = \int_{\mathbf{x}} \phi(\mathbf{x}) P(\mathbf{x}) d\mathbf{x}$ by

$$\hat{\Phi} = \sum_{r=1}^R \phi(\mathbf{x}^{(r)}) \frac{P^*(\mathbf{x}^{(r)})}{Z_R}$$

- Is there anything wrong with this strategy?

Is there anything wrong with this strategy?

- Let's assume $\phi(\mathbf{x})$ is a benign, smoothly varying function, and concentrated on the nature of $P^*(\mathbf{x})$
- A high dimensional distribution is often concentrated in a small region of the state space known as its typical set T
 - whose volume is given by $|T| \simeq 2^{H(\mathbf{X})}$
 - $H(\mathbf{X})$ is the entropy of the probability distribution
$$H(\mathbf{X}) = \sum_{\mathbf{x}} P(\mathbf{x}) \log_2 \frac{1}{P(\mathbf{x})}$$
- $\Phi = \int_{\mathbf{x}} \phi(\mathbf{x}) P(\mathbf{x}) d\mathbf{x}$ will be principally determined by values in typical set
- If we have N random variables with binary values, the total size of state space is 2^N and the typical set size is 2^H
 - Each sample has a chance $2^H/2^N$ of falling into typical set
 - We need $R_{\min} \simeq O(2^{N-H})$ samples

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Importance Sampling

- Importance sampling is not a method for generating samples from $P(\mathbf{x})$
- It is just a method for estimating the expectations of a function $\phi(\mathbf{x})$
- Let's imagine the target distribution is a one-dimensional density $P(x)$

$$P(x) = \frac{P^*(x)}{Z}$$

but $P(x)$ is too complicated to sample from directly

- We assume $Q(x) = \frac{Q^*(x)}{Z_Q}$ is a simpler density from which we can generate samples
- In importance sampling, we generate R samples $\{x^{(r)}\}_{r=1}^R$ from $Q(x)$
- Then Φ can be estimated by

$$\hat{\Phi} = \sum_{r=1}^R \frac{\sum_r w_r \phi(x^{(r)})}{\sum_r w_r}$$

where $w_r = \frac{P^*(x^{(r)})}{Q^*(x^{(r)})}$

Importance Sampling: A Toy Example

- Φ can be estimated by

$$\hat{\Phi} = \sum_{r=1}^R \frac{\sum_r w_r \phi(x^{(r)})}{\sum_r w_r}$$

where $w_r = \frac{P^*(x^{(r)})}{Q^*(x^{(r)})}$

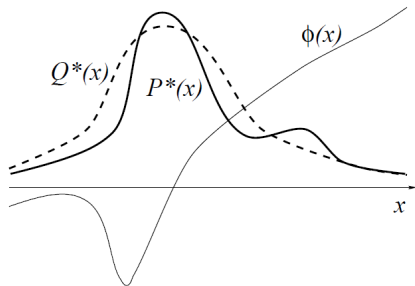


Figure 2. Functions involved in importance sampling. We wish to estimate the expectation of $\phi(x)$ under $P(x) \propto P^*(x)$. We can generate samples from the simpler distribution $Q(x) \propto Q^*(x)$. We can evaluate Q^* and P^* at any point.

Importance Sampling: A Toy Example

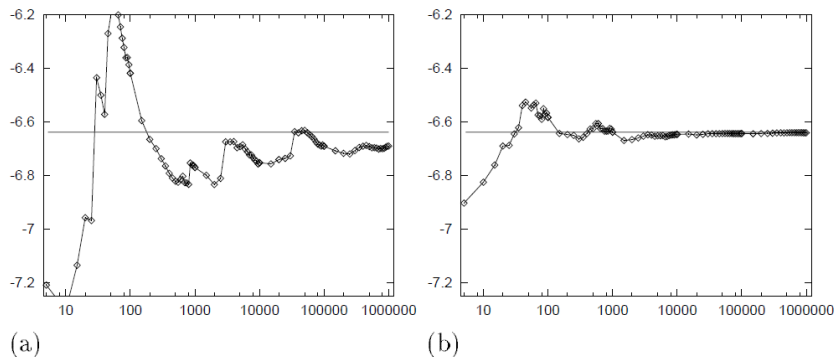


Figure 3. Importance sampling in action: a) using a Gaussian sampler density; b) using a Cauchy sampler density. Horizontal axis shows number of samples on a log scale. Vertical axis shows the estimate $\hat{\Phi}$. The horizontal line indicates the true value of Φ .

Importance Sampling in Many Dimensions

- Importance sampling suffers from two difficulties
 - We clearly need to obtain samples that lie in the typical set
 - Even if we obtain samples in the typical set, the weights associated with these samples are likely to vary by large factors

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Rejection Sampling

- We again assume one dimensional complicated density $P(x) = \frac{P^*(x)}{Z}$
- We assume a simpler **proposal density** $Q(x)$ which we can evaluate and can generate samples from
- We further assume for all x , $cQ^*(x) > P^*(x)$

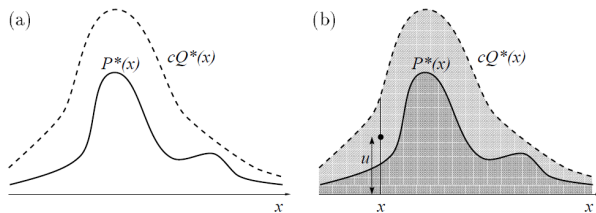


Figure 4. Rejection sampling. a) The functions involved in rejection sampling. We desire samples from $P(x) \propto P^*(x)$. We are able to draw samples from $Q(x) \propto Q^*(x)$, and we know a value c such that $cQ^*(x) > P^*(x)$ for all x . b) A point (x, u) is generated at random in the lightly shaded area under the curve $cQ^*(x)$. If this point also lies below $P^*(x)$ then it is accepted.

Rejection Sampling

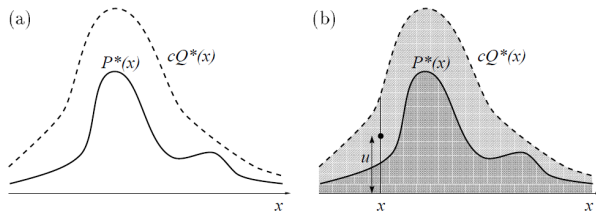


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- We first generate x from $Q(x)$
- We evaluate $cQ^*(x)$ and generate a uniformly distributed variable from the interval $[0, cQ^*(x)]$
- We then evaluate $P^*(x)$ and accept or reject the sample x by comparing u with $P^*(x)$
 - If $u > P^*(x)$ then x is rejected

Rejection Sampling

- Reject sampling will work best when Q is a good proximation of P
- c grows exponentially with the dimensionality N (MacKay (1998))
- While it is a useful method for one-dimensional problems, it is not a practical technique for high-dimensional distributions $P(\mathbf{x})$

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Motivation

- Importance sampling and rejection sampling only work well if the proposal density $Q(x)$ is similar to $P(x)$
- In large and complex problems, it is difficult to create a single density $Q(x)$ that has this property
- The Metropolis algorithm makes use of a proposal density $Q(x', x^{(t)})$ which depends on the current state $x^{(t)}$
 - It is not necessarily for $Q(x', x^{(t)})$ to look at all similar to $P(x)$

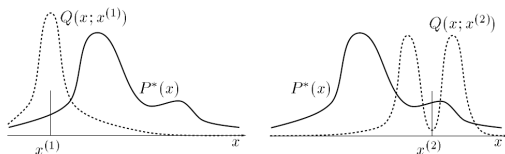


Figure 6. Metropolis method in one dimension. The proposal distribution $Q(x'; x)$ is here shown as having a shape that changes as x changes, though this is not typical of the proposal densities used in practice.

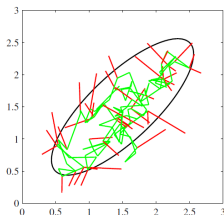
Metropolis Method

- We again assume that we can evaluate $P^*(x)$ for any x
- A tentative new x' is generated from the proposal density $Q(x'|x^{(t)})$
- To decide whether to accept the new state, we compute the quantity:

$$a = \frac{P^*(x')Q(x^{(t)}|x')}{P^*(x^{(t)})Q(x'|x^{(t)})}$$

- If $a \geq 1$, then accept the new x'
 - Otherwise, the new state is accepted with probability a
- If the state is accepted, we set $x^{(t+1)} = x'$
 - If the state is rejected, we set $x^{(t+1)} = x^{(t)}$
- This is different from rejection sampling: in Metropolis method, a rejection causes the current state sent to the generated list another time
- The samples in a Metropolis simulation of T iterations are correlated

$$a = \frac{P^*(x')Q(x^{(t)}|x')}{P^*(x^{(t)})Q(x'|x^{(t)})}$$



Steps that are accepted are shown as green lines, and rejected steps are shown in red.

- When $Q(x^{(t)}, x') = Q(x', x^{(t)})$, it is called Metropolis algorithm (Metropolis et al., 1953)
- When $Q(x^{(t)}, x') \neq Q(x', x^{(t)})$, it is known as Metropolis-Hastings algorithm (Hastings, 1970)
- Metropolis methods are known as Markov chain Monte Carlo methods

Markov Chains

- A first order Markov chain is defined to be a series of random variables $x^{(1)}, \dots, x^{(M)}$ such that the following conditional independence property holds

$$P(x^{(t+1)} | x^{(1)}, \dots, x^{(t)}) = P(x^{(t+1)} | x^{(t)})$$

where we define the transition probability

$$T(x^{(t)}, x^{(t+1)}) = P(x^{(t+1)} | x^{(t)})$$

- A Markov chain is called **homogeneous** if the transition probabilities are the same for all t
- The marginal probability for a particular variable can be expressed in terms of the marginal probability for the previous variable in the chain in the form

$$P(x^{(t+1)}) = \sum_{x^{(t)}} P(x^{(t+1)} | x^{(t)}) P(x^{(t)})$$

Markov Chains (Cont'd)

- A distribution 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 $P(x)$ is invariant if

$$P(x) = \sum_{x'} T(x', x) P(x')$$

- A sufficient (but not necessary) condition for ensuring that the required distribution $P(x)$ is invariant is to choose the transition probabilities to satisfy the property of detailed balance, defined by

$$P(x) T(x, x') = P(x') T(x', x)$$

- It is easy to verify that

$$\sum_{x'} T(x', x) P(x') = \sum_{x'} P(x) T(x, x') = P(x) \sum_{x'} T(x, x') = P(x)$$

Metropolis methods are known as Markov chain Monte Carlo methods

- From the probability to accept a new state:

$$a(x, x^{(t)}) = \min \left(1, \frac{P^*(x') Q(x^{(t)} | x')}{P^*(x^{(t)}) Q(x' | x^{(t)})} \right)$$

- We have the joint probability of two consecutive states as

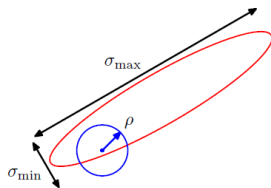
$$\begin{aligned} & P^*(x^{(t)}) \cdot Q(x' | x^{(t)}) a(x, x^{(t)}) \\ = & \min(P^*(x^{(t)}) Q(x' | x^{(t)}), P^*(x') Q(x^{(t)} | x')) \\ = & \min(P^*(x') Q(x^{(t)} | x'), P^*(x^{(t)}) Q(x' | x^{(t)})) \\ = & P^*(x) \cdot Q(x^{(t)} | x') a(x^{(t)}, x') \end{aligned}$$

as required

- Metropolis method actually samples from the required distribution $P(x)$

Sampling Effects

- The specific choice of proposal distribution can have a marked effect on the performance of the algorithm



- The scale ρ of the proposal distribution should be on the order of the smallest standard deviation $\sigma_{\min}$
- The iteration T should be at least $(\sigma_{\max}/\sigma_{\min})^2$ to obtain approximately independent samples

Simulation of Sampling

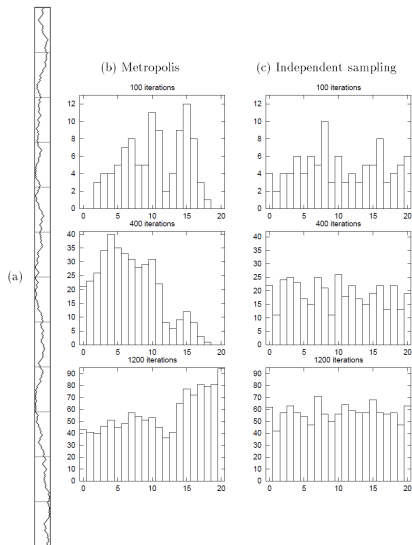


Figure 8. Metropolis method for a toy problem. (a) The state sequence for $t = 1 \dots 600$. Horizontal direction = states from 0 to 20; vertical direction = time from 1 to 600; the cross bars mark time intervals of duration 50. (b) Histogram of occupancy of the states after 100, 400 and 1200 iterations. (c) For comparison, histograms resulting when successive points are drawn *independently* from the target distribution.

$$P(x) = \begin{cases} \frac{1}{21} & x \in \{0, 1, \dots, 20\} \\ 0 & \text{otherwise} \end{cases}$$

$$Q(x'|x) = \begin{cases} \frac{1}{2} & x' = x \pm 1 \\ 0 & \text{otherwise} \end{cases}$$

- Rejection will occur only when the proposal takes the state $x' = -1$ or $x' = 21$
- It takes $\approx T^2 = 100$ (178) iterations to reach 0 or 20
- It takes ≈ 400 (540) iterations to reach both 0 and 20

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Gibbs Sampling

- In the general case of a system with K variables, a single iteration involves sampling one parameter at a time:

- $x_1^{(t+1)} \sim P(x_2^{(t)}, x_3^{(t)}, \dots, x_K^{(t)})$
- $x_2^{(t+1)} \sim P(x_1^{(t+1)}, x_3^{(t)}, \dots, x_K^{(t)})$
- $x_3^{(t+1)} \sim P(x_1^{(t+1)}, x_2^{(t+1)}, \dots, x_K^{(t)})$
- ...
- $x_K^{(t+1)} \sim P(x_1^{(t+1)}, x_2^{(t+1)}, \dots, x_{K-1}^{(t+1)})$

- Denote $\mathbf{x}_{\setminus k}^{(t)} = P(x_1^{(t+1)}, x_2^{(t+1)}, \dots, x_{k-1}^{(t+1)}, x_{k+1}^{(t)}, \dots, x_K^{(t)})$

- Gibbs sampling can be viewed as a Metropolis method

$$\begin{aligned} a_G &= \frac{P^*(\mathbf{x}')Q(\mathbf{x}^{(t)}|\mathbf{x}')}{P^*(\mathbf{x}^{(t)})Q(\mathbf{x}'|\mathbf{x}^{(t)})} = \frac{P(\mathbf{x}')P(x_k^{(t)}|\mathbf{x}'_{\setminus k})}{P(\mathbf{x}^{(t)})P(x'_k|\mathbf{x}_{\setminus k}^{(t)})} \\ &= \frac{P(x'_k|\mathbf{x}'_{\setminus k})P(\mathbf{x}'_{\setminus k})P(x_k^{(t)}|\mathbf{x}'_{\setminus k})}{P(x_k^{(t)}|\mathbf{x}_{\setminus k}^{(t)})P(\mathbf{x}_{\setminus k}^{(t)})P(x'_k|\mathbf{x}_{\setminus k}^{(t)})} \stackrel{\mathbf{x}'_{\setminus k} = \mathbf{x}_{\setminus k}^{(t)}}{=} \frac{P(x'_k|\mathbf{x}'_{\setminus k})P(\mathbf{x}'_{\setminus k})P(x_k^{(t)}|\mathbf{x}'_{\setminus k})}{P(x_k^{(t)}|\mathbf{x}'_{\setminus k})P(\mathbf{x}'_{\setminus k})P(x'_k|\mathbf{x}'_{\setminus k})} = 1 \end{aligned}$$

- The samples are always accepted

Example of Gibbs Sampling

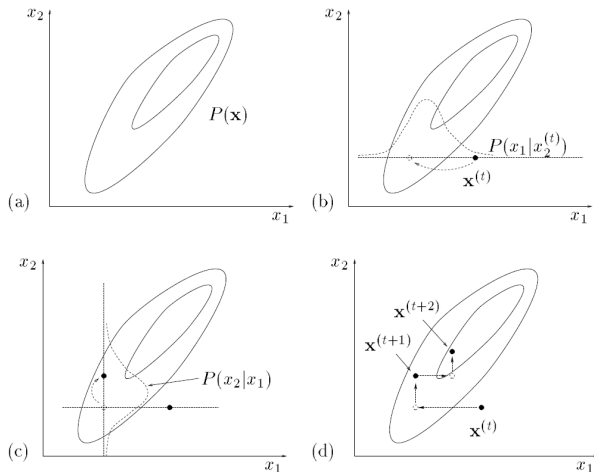


Figure 9. Gibbs sampling. (a) The joint density $P(\mathbf{x})$ from which samples are required. (b) Starting from a state $\mathbf{x}^{(t)}$, x_1 is sampled from the conditional density $P(x_1|x_2^{(t)})$. (c) A sample is then made from the conditional density $P(x_2|x_1)$. (d) A couple of iterations of Gibbs sampling.

MacKay, D. J. C. (1998). Introduction to Monte Carlo methods. In Jordan, M. I., editor, *Learning in Graphical Models*, NATO Science Series, pages 175–204. Kluwer Academic Press.