

# Sampling and Monte Carlo Integration

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# Recap

Learning and inference often involves intractable integrals, e.g.

- ▶ Marginalisation

$$p(\mathbf{x}) = \int_{\mathbf{y}} p(\mathbf{x}, \mathbf{y}) d\mathbf{y}$$

- ▶ Expectations

$$\mathbb{E}[g(\mathbf{x}) \mid \mathbf{y}_o] = \int g(\mathbf{x}) p(\mathbf{x} \mid \mathbf{y}_o) d\mathbf{x}$$

for some function  $g$ .

- ▶ For unobserved variables, likelihood and gradient of the log lik

$$L(\boldsymbol{\theta}) = p(\mathcal{D}; \boldsymbol{\theta}) = \int_{\mathbf{u}} p(\mathbf{u}, \mathcal{D}; \boldsymbol{\theta}) d\mathbf{u},$$
$$\nabla_{\boldsymbol{\theta}} \ell(\boldsymbol{\theta}) = \mathbb{E}_{p(\mathbf{u} \mid \mathcal{D}; \boldsymbol{\theta})} [\nabla_{\boldsymbol{\theta}} \log p(\mathbf{u}, \mathcal{D}; \boldsymbol{\theta})]$$

Notation:  $\mathbb{E}_{p(\mathbf{x})}$  is sometimes used to indicate that the expectation is taken with respect to  $p(\mathbf{x})$ .

# Recap

Learning and inference often involves intractable integrals, e.g.

- ▶ For unnormalised models with intractable partition functions

$$L(\boldsymbol{\theta}) = \frac{\tilde{p}(\mathcal{D}; \boldsymbol{\theta})}{\int_{\mathbf{x}} \tilde{p}(\mathbf{x}; \boldsymbol{\theta}) d\mathbf{x}}$$

$$\nabla_{\boldsymbol{\theta}} \ell(\boldsymbol{\theta}) \propto \mathbf{m}(\mathcal{D}; \boldsymbol{\theta}) - \mathbb{E}_{p(\mathbf{x}; \boldsymbol{\theta})} [\mathbf{m}(\mathbf{x}; \boldsymbol{\theta})]$$

- ▶ Combined case of unnormalised models with intractable partition functions and unobserved variables.
- ▶ Evaluation of intractable integrals can sometimes be avoided by using other learning criteria (e.g. score matching).
- ▶ Here: methods to approximate integrals like those above using sampling.

# Program

1. Monte Carlo integration
2. Sampling

# Program

## 1. Monte Carlo integration

- Approximating expectations by averages
- Importance sampling

## 2. Sampling

# Averages with iid samples

- ▶ (From tutorials): For Gaussians, the sample average is an estimate (MLE) of the mean (expectation)  $\mathbb{E}[x]$

$$\bar{x} = \frac{1}{n} \sum_{i=1}^n x_i \approx \mathbb{E}[x]$$

- ▶ Gaussianity not needed: assume  $x_i$  are iid observations of  $x \sim p(x)$ .

$$\mathbb{E}[x] = \int xp(x)dx \approx \bar{x}_n \qquad \bar{x}_n = \frac{1}{n} \sum_{i=1}^n x_i$$

- ▶ Subscript  $n$  reminds us that we used  $n$  samples to compute the average.
- ▶ Approximating integrals by means of sample averages is called Monte Carlo integration.

# Averages with iid samples

- ▶ Sample average is unbiased

$$\mathbb{E}[\bar{x}_n] = \frac{1}{n} \sum_{i=1}^n \mathbb{E}[x_i] \stackrel{*}{=} \frac{n}{n} \mathbb{E}[x] = \mathbb{E}[x]$$

(\*: “identically distributed” assumption is used, not independence)

- ▶ Variability

$$\mathbb{V}[\bar{x}_n] = \frac{1}{n^2} \mathbb{V}\left[\sum_{i=1}^n x_i\right] \stackrel{*}{=} \frac{1}{n^2} \sum_{i=1}^n \mathbb{V}[x_i] = \frac{1}{n} \mathbb{V}[x]$$

(\*: independence assumption used)

- ▶ Expected squared error decreases as  $1/n$

$$\mathbb{V}[\bar{x}_n] = \mathbb{E}\left[(\bar{x}_n - \mathbb{E}[x])^2\right] = \frac{1}{n} \mathbb{V}[x]$$

# Averages with iid samples

- ▶ Weak law of large numbers:

$$\mathbb{P} (|\bar{x}_n - \mathbb{E}[x]| \geq \epsilon) \leq \frac{\mathbb{V}[x]}{n\epsilon^2}$$

- ▶ As  $n \rightarrow \infty$ , the probability for the sample average to deviate from the expected value goes to zero.
- ▶ We say that sample average converges in probability to the expected value.
- ▶ Speed of convergence depends on the variance  $\mathbb{V}[x]$ .
- ▶ Different “laws of large numbers” exist that make different assumptions.



# Chebyshev's inequality

- ▶ Weak law of large numbers is a direct consequence of Chebyshev's inequality
- ▶ Chebyshev's inequality: Let  $s$  be some random variable with mean  $\mathbb{E}[s]$  and variance  $\mathbb{V}[s]$ .

$$\mathbb{P}(|s - \mathbb{E}[s]| \geq \epsilon) \leq \frac{\mathbb{V}[s]}{\epsilon^2}$$

- ▶ This means that for *all* random variables:
  - ▶ probability to deviate more than three standard deviation from the mean is less than  $1/9 \approx 0.11$   
(set  $\epsilon = 3\sqrt{\mathbb{V}(s)}$ )
  - ▶ Probability to deviate more than 6 standard deviations:  $1/36 \approx 0.03$ .

These are conservative values; for many distributions, the probabilities will be smaller.

- ▶ Chebyshev's inequality follows from Markov's inequality.
- ▶ Markov's inequality: For a random variable  $y \geq 0$

$$\mathbb{P}(y \geq t) \leq \frac{\mathbb{E}[y]}{t} \quad (t > 0)$$

- ▶ Chebyshev's inequality is obtained by setting  $y = |s - \mathbb{E}[s]|$

$$\begin{aligned} \mathbb{P}(|s - \mathbb{E}[s]| \geq t) &= \mathbb{P}\left((s - \mathbb{E}[s])^2 \geq t^2\right) \\ &\leq \frac{\mathbb{E}[(s - \mathbb{E}[s])^2]}{t^2}. \end{aligned}$$

Chebyshev's inequality follows with  $t = \epsilon$ , and because  $\mathbb{E}[(s - \mathbb{E}[s])^2]$  is the variance  $\mathbb{V}[s]$  of  $s$ .

# Proofs (not examinable)

Proof for Markov's inequality: Let  $t$  be an arbitrary positive number and  $y$  a one-dimensional non-negative random variable with pdf  $p$ . We can decompose the expectation of  $y$  using  $t$  as split-point,

$$\mathbb{E}[y] = \int_0^{\infty} up(u)du = \int_0^t up(u)du + \int_t^{\infty} up(u)du.$$

Since  $u \geq t$  in the second term, we obtain the inequality

$$\mathbb{E}[y] \geq \int_0^t up(u)du + \int_t^{\infty} tp(u)du.$$

The second term is  $t$  times the probability that  $y \geq t$ , so that

$$\begin{aligned} \mathbb{E}[y] &\geq \int_0^t up(u)du + t\mathbb{P}(y \geq t) \\ &\geq t\mathbb{P}(y \geq t), \end{aligned}$$

where the second line holds because the first term in the first line is non-negative. This gives Markov's inequality

$$\mathbb{P}(y \geq t) \leq \frac{\mathbb{E}(y)}{t} \quad (t > 0)$$

# Averages with correlated samples

- ▶ When computing the variance of the sample average

$$\mathbb{V}[\bar{x}_n] = \frac{\mathbb{V}[x]}{n}$$

we assumed the samples are identically and independently distributed.

- ▶ The variance shrinks with increasing  $n$  and the average becomes more and more concentrated around  $\mathbb{E}[x]$ .
- ▶ Corresponding results exist for the case of statistically dependent samples  $x_i$ . Known as “ergodic theorems”.
- ▶ Important for the theory of Markov chain Monte Carlo methods (outside the scope of our lecture).

# More general expectations

- ▶ So far, we have considered

$$\mathbb{E}[x] = \int xp(x)dx \approx \frac{1}{n} \sum_{i=1}^n x_i$$

where  $x_i \sim p(x)$

- ▶ This generalises

$$\mathbb{E}[g(\mathbf{x})] = \int g(\mathbf{x})p(\mathbf{x})d\mathbf{x} \approx \frac{1}{n} \sum_{i=1}^n g(\mathbf{x}_i)$$

where  $\mathbf{x}_i \sim p(\mathbf{x})$

- ▶ Variance of the approximation if the  $\mathbf{x}_i$  are iid is  $\frac{1}{n}\mathbb{V}[g(\mathbf{x})]$

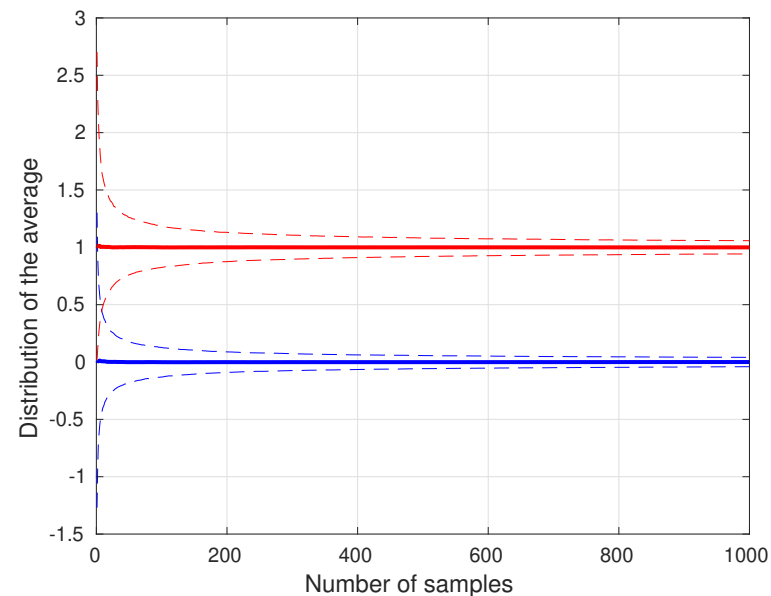
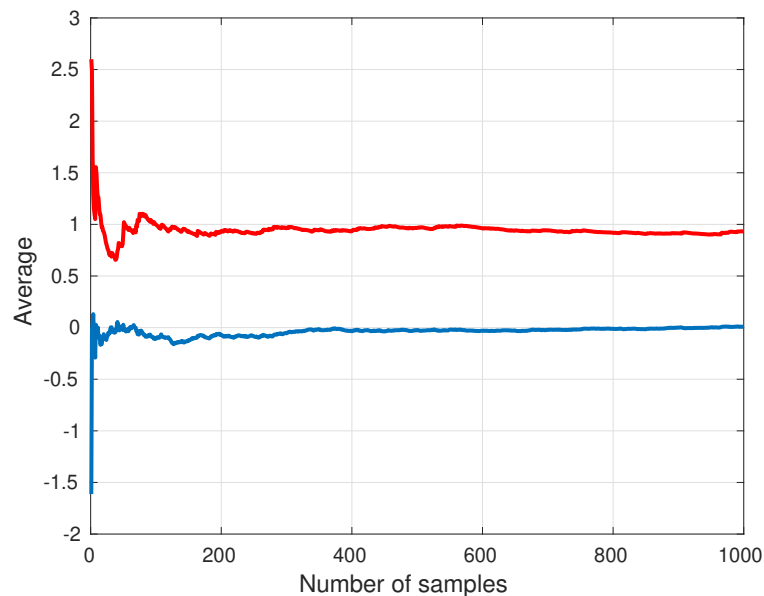
# Example (Based on a slide from Amos Storkey)

$$\mathbb{E}[g(x)] = \int g(x)\mathcal{N}(x; 0, 1)dx \approx \frac{1}{n} \sum_{i=1}^n g(x_i) \quad (x_i \sim \mathcal{N}(x; 0, 1))$$

for  $g(x) = x$  and  $g(x) = x^2$

Left: sample average as a function of  $n$

Right: Variability (0.5 quantile: solid, 0.1 and 0.9 quantiles: dashed)



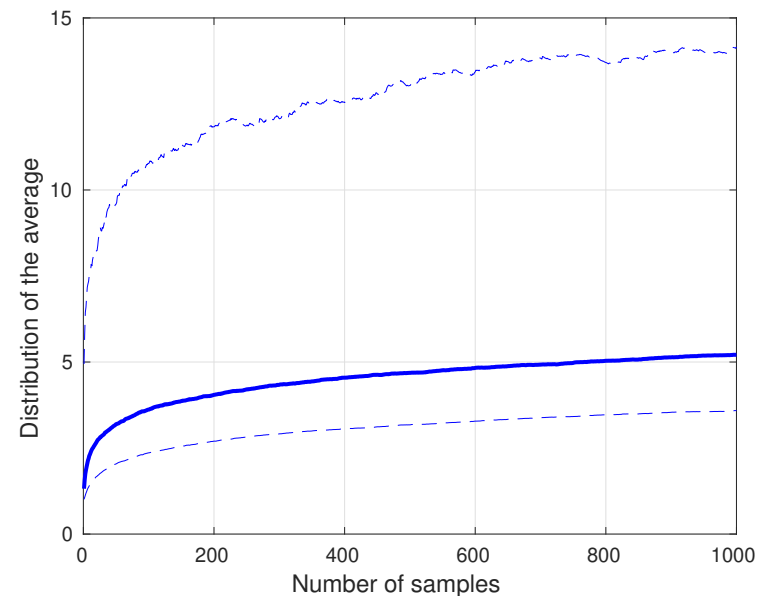
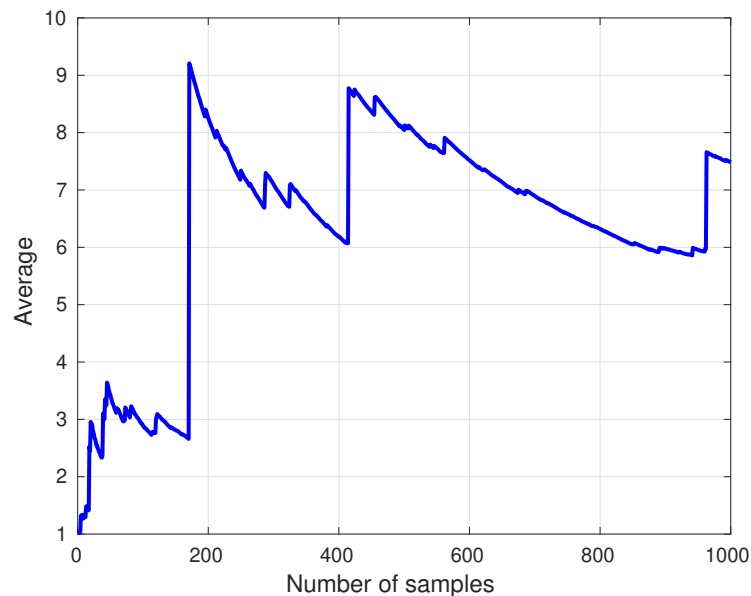
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for  $g(x) = \exp(0.6x^2)$

Left: sample average as a function of  $n$

Right: Variability (0.5 quantile: solid, 0.1 and 0.9 quantiles: dashed)



# Example

- ▶ Indicators that something is wrong:
  - ▶ Strong fluctuations in the sample average as  $n$  increases.
  - ▶ Large non-declining variability.
- ▶ Note: integral is not finite:

$$\begin{aligned}\int \exp(0.6x^2)\mathcal{N}(x; 0, 1)dx &= \frac{1}{\sqrt{2\pi}} \int \exp(0.6x^2) \exp(-0.5x^2)dx \\ &= \frac{1}{\sqrt{2\pi}} \int \exp(0.1x^2)dx \\ &= \infty\end{aligned}$$

but for any  $n$ , the sample average is finite and may be mistaken for a good approximation.

- ▶ Check variability when approximating the expected value by a sample average!



# Approximating general integrals

- ▶ If the integral does not correspond to an expectation, we can smuggle in a pdf  $q$  to rewrite it as an expected value with respect to  $q$

$$\begin{aligned} I &= \int g(\mathbf{x}) d\mathbf{x} = \int g(\mathbf{x}) \frac{q(\mathbf{x})}{q(\mathbf{x})} d\mathbf{x} \\ &= \int \frac{g(\mathbf{x})}{q(\mathbf{x})} q(\mathbf{x}) d\mathbf{x} \\ &= \mathbb{E}_{q(\mathbf{x})} \left[ \frac{g(\mathbf{x})}{q(\mathbf{x})} \right] \\ &\approx \frac{1}{n} \sum_{i=1}^n \frac{g(\mathbf{x}_i)}{q(\mathbf{x}_i)} \end{aligned}$$

with  $x_i \sim q(\mathbf{x})$  (iid)

- ▶ This is the basic idea of importance sampling.
- ▶  $q$  is called the importance (or proposal) distribution

# Choice of the importance distribution

- ▶ Call the approximation  $\hat{I}$ ,

$$\hat{I} = \frac{1}{n} \sum_{i=1}^n \frac{g(\mathbf{x}_i)}{q(\mathbf{x}_i)}$$

- ▶  $\hat{I}$  is unbiased by construction

$$\mathbb{E}[\hat{I}] = \mathbb{E} \left[ \frac{g(\mathbf{x})}{q(\mathbf{x})} \right] = \int \frac{g(\mathbf{x})}{q(\mathbf{x})} q(\mathbf{x}) d\mathbf{x} = \int g(\mathbf{x}) d\mathbf{x} = I$$

- ▶ Variance

$$\mathbb{V}[\hat{I}] = \frac{1}{n} \mathbb{V} \left[ \frac{g(\mathbf{x})}{q(\mathbf{x})} \right] = \frac{1}{n} \mathbb{E} \left[ \left( \frac{g(\mathbf{x})}{q(\mathbf{x})} \right)^2 \right] - \frac{1}{n} \underbrace{\left( \mathbb{E} \left[ \frac{g(\mathbf{x})}{q(\mathbf{x})} \right] \right)^2}_{I^2}$$

Depends on the second moment.

# Choice of the importance distribution

- ▶ The second moment is

$$\begin{aligned}\mathbb{E} \left[ \left( \frac{g(\mathbf{x})}{q(\mathbf{x})} \right)^2 \right] &= \int \left( \frac{g(\mathbf{x})}{q(\mathbf{x})} \right)^2 q(\mathbf{x}) d\mathbf{x} = \int \frac{g(\mathbf{x})^2}{q(\mathbf{x})} d\mathbf{x} \\ &= \int |g(\mathbf{x})| \frac{|g(\mathbf{x})|}{q(\mathbf{x})} d\mathbf{x}\end{aligned}$$

- ▶ Bad:  $q(\mathbf{x})$  is small when  $|g(\mathbf{x})|$  is large. Gives large variance.
- ▶ Good:  $q(\mathbf{x})$  is large when  $|g(\mathbf{x})|$  is large.
- ▶ Optimal  $q$  equals

$$q^*(\mathbf{x}) = \frac{|g(\mathbf{x})|}{\int |g(\mathbf{x})| d\mathbf{x}}$$

- ▶ Optimal  $q$  cannot be computed, but justifies the heuristic that  $q(\mathbf{x})$  should be large when  $|g(\mathbf{x})|$  is large, or that **the ratio  $|g(\mathbf{x})|/q(\mathbf{x})$  should be approximately constant** .

# Proof (not examinable)

Since the variance of a random variable  $|x|$  is non-negative and can be written as

$$\mathbb{V}[|x|] = \mathbb{E}[x^2] - (\mathbb{E}[|x|])^2,$$

we have

$$\mathbb{E}[x^2] \geq \mathbb{E}[|x|]^2$$

The smallest second moment achieves equality. We now verify that for  $q^*(\mathbf{x})$ , we have

$$\mathbb{E} \left[ \left( \frac{g(\mathbf{x})}{q^*(\mathbf{x})} \right)^2 \right] = \mathbb{E} \left[ \left| \frac{g(\mathbf{x})}{q^*(\mathbf{x})} \right| \right]^2$$

# Proof (not examinable)

Indeed, for the optimal  $q$ , we have

$$\begin{aligned}\mathbb{E} \left[ \left( \frac{g(\mathbf{x})}{q^*(\mathbf{x})} \right)^2 \right] &= \int |g(\mathbf{x})| \frac{|g(\mathbf{x})|}{q^*(\mathbf{x})} d\mathbf{x} \\ &= \int |g(\mathbf{x})| d\mathbf{x} \int |g(\mathbf{x})|^2 \frac{1}{|g(\mathbf{x})|} d\mathbf{x} \\ &= \left( \int |g(\mathbf{x})| d\mathbf{x} \right)^2\end{aligned}$$

and

$$\begin{aligned}\mathbb{E} \left[ \left| \frac{g(\mathbf{x})}{q^*(\mathbf{x})} \right| \right]^2 &= \left( \int \left| \frac{g(\mathbf{x})}{q^*(\mathbf{x})} \right| q^*(\mathbf{x}) d\mathbf{x} \right)^2 \\ &= \left( \int |g(\mathbf{x})| d\mathbf{x} \right)^2,\end{aligned}$$

which concludes the proof.

# Importance sampling to compute the partition function

We can use importance sampling to approximate the partition function for unnormalised models  $\tilde{p}(\mathbf{x}; \boldsymbol{\theta})$ .

$$\begin{aligned} Z(\boldsymbol{\theta}) &= \int \tilde{p}(\mathbf{x}; \boldsymbol{\theta}) d\mathbf{x} \\ &= \int \tilde{p}(\mathbf{x}; \boldsymbol{\theta}) \frac{q(\mathbf{x})}{q(\mathbf{x})} d\mathbf{x} \\ &= \int \frac{\tilde{p}(\mathbf{x}; \boldsymbol{\theta})}{q(\mathbf{x})} q(\mathbf{x}) d\mathbf{x} \\ &\approx \frac{1}{n} \sum_{i=1}^n \frac{\tilde{p}(\mathbf{x}_i; \boldsymbol{\theta})}{q(\mathbf{x}_i)} \quad (\mathbf{x}_i \sim q(\mathbf{x}) \text{ iid}) \end{aligned}$$

# Example

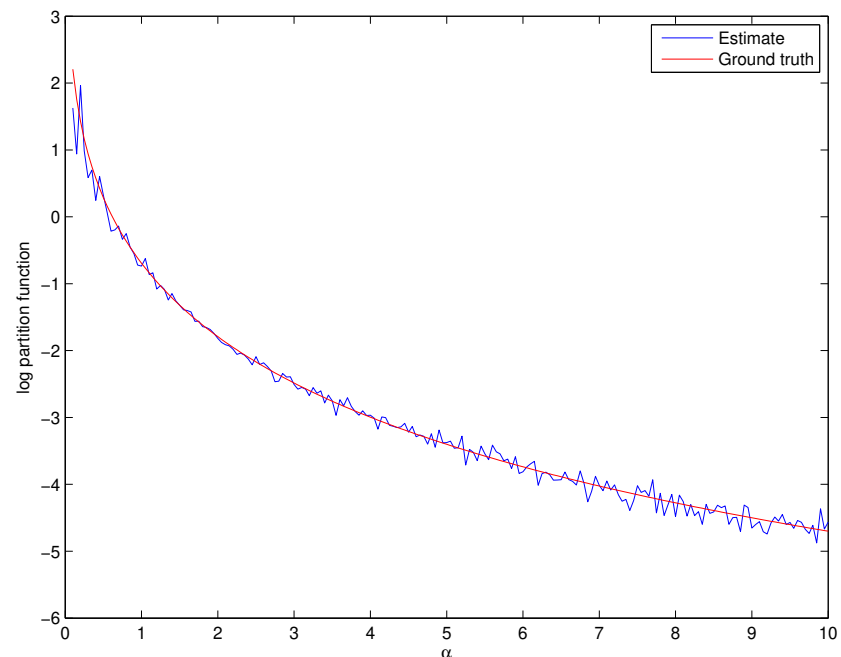
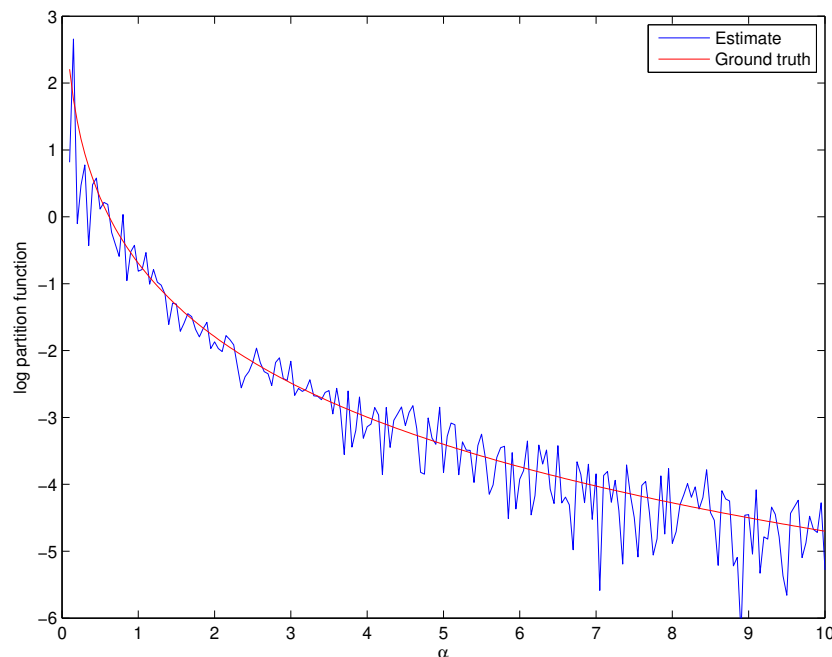
Approximating the log partition function of the unnormalised beta-distribution

$$\tilde{p}(x; \alpha, \beta) = x^{\alpha-1}(1-x)^{\beta-1}, \quad x \in [0, 1]$$

for  $\beta$  fixed to  $\beta = 2$ .

Importance distribution: uniform distribution on  $[0, 1]$

Left:  $n = 10$ , right:  $n = 100$ .



# Importance sampling to compute expectations

- ▶ Assume you would like to approximate  $\mathbb{E}_{p(\mathbf{x})}[g(\mathbf{x})]$  by a sample average but sampling from  $p(\mathbf{x})$  is difficult.
- ▶ We can write

$$\begin{aligned}\mathbb{E}_{p(\mathbf{x})}[g(\mathbf{x})] &= \int g(\mathbf{x})p(\mathbf{x})d\mathbf{x} \\ &= \int g(\mathbf{x})\frac{p(\mathbf{x})}{q(\mathbf{x})}q(\mathbf{x})d\mathbf{x} \\ &= \mathbb{E}_{q(\mathbf{x})}\left[g(\mathbf{x})\frac{p(\mathbf{x})}{q(\mathbf{x})}\right] \\ &\approx \frac{1}{n}\sum_{i=1}^n g(\mathbf{x}_i)\frac{p(\mathbf{x}_i)}{q(\mathbf{x}_i)}\end{aligned}$$

where  $\mathbf{x}_i \sim q(\mathbf{x})$  (iid)

- ▶ The  $w_i = p(\mathbf{x}_i)/q(\mathbf{x}_i)$  are called the importance weights.



# Normalised importance weights

- ▶ We can combine the above ideas to approximate

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

by importance sampling even if we only know  $\tilde{p}(\mathbf{x}) \propto p(\mathbf{x})$  and

$$p(\mathbf{x}) = \frac{\tilde{p}(\mathbf{x})}{\int \tilde{p}(\mathbf{x})d\mathbf{x}}$$

- ▶ Write

$$\begin{aligned} \int g(\mathbf{x})p(\mathbf{x})d\mathbf{x} &= \frac{\int g(\mathbf{x})\tilde{p}(\mathbf{x})d\mathbf{x}}{\int \tilde{p}(\mathbf{x})d\mathbf{x}} \\ &= \frac{\int g(\mathbf{x})\frac{\tilde{p}(\mathbf{x})}{q(\mathbf{x})}q(\mathbf{x})d\mathbf{x}}{\int \frac{\tilde{p}(\mathbf{x})}{q(\mathbf{x})}q(\mathbf{x})d\mathbf{x}} \\ &= \frac{\mathbb{E}_{q(\mathbf{x})} \left[ g(\mathbf{x})\frac{\tilde{p}(\mathbf{x})}{q(\mathbf{x})} \right]}{\mathbb{E}_{q(\mathbf{x})} \left[ \frac{\tilde{p}(\mathbf{x})}{q(\mathbf{x})} \right]} \end{aligned}$$

# Normalised importance weights

- ▶ Since

$$\begin{aligned}\int g(\mathbf{x})p(\mathbf{x})d\mathbf{x} &= \frac{\mathbb{E}_{q(\mathbf{x})} \left[ g(\mathbf{x}) \frac{\tilde{p}(\mathbf{x})}{q(\mathbf{x})} \right]}{\mathbb{E}_{q(\mathbf{x})} \left[ \frac{\tilde{p}(\mathbf{x})}{q(\mathbf{x})} \right]} \\ &= \frac{\mathbb{E}_{q(\mathbf{x})} \left[ g(\mathbf{x}) \frac{\tilde{p}(\mathbf{x})}{\tilde{q}(\mathbf{x})} \right]}{\mathbb{E}_{q(\mathbf{x})} \left[ \frac{\tilde{p}(\mathbf{x})}{\tilde{q}(\mathbf{x})} \right]}\end{aligned}$$

we only need to know the importance distribution  $q(\mathbf{x})$  up to normalisation constant.

- ▶ Approximate both expectations by a sample average

$$\int g(\mathbf{x})p(\mathbf{x})d\mathbf{x} \approx \frac{\frac{1}{n} \sum_{i=1}^n g(\mathbf{x}_i) \frac{\tilde{p}(\mathbf{x}_i)}{\tilde{q}(\mathbf{x}_i)}}{\frac{1}{n} \sum_{i=1}^n \frac{\tilde{p}(\mathbf{x}_i)}{\tilde{q}(\mathbf{x}_i)}}$$

where  $\mathbf{x}_i \sim q(\mathbf{x})$  (iid)

# Normalised importance weights

- ▶ With importance weights

$$w_i = \frac{\tilde{p}(\mathbf{x}_i)}{\tilde{q}(\mathbf{x}_i)},$$

where  $\mathbf{x}_i \stackrel{\text{iid}}{\sim} q(\mathbf{x})$ , we can write

$$\int g(\mathbf{x})p(\mathbf{x})d\mathbf{x} \approx \frac{\sum_{i=1}^n g(\mathbf{x}_i)w_i}{\sum_{i=1}^n w_i}$$

- ▶ Same weights in numerator and denominator.
- ▶ The quantities

$$\frac{w_i}{\sum_{i=1}^n w_i}$$

are called normalised importance weights.

# Program

## 1. Monte Carlo integration

- Approximating expectations by averages
- Importance sampling

## 2. Sampling

# Program

## 1. Monte Carlo integration

## 2. Sampling

- Simple univariate sampling
- Rejection sampling
- Ancestral sampling
- Gibbs sampling

# Assumption

- ▶ We assume that we are able to generate iid samples from the uniform distribution on  $[0, 1]$ .
- ▶ How to do that: see e.g.  
<https://statweb.stanford.edu/~owen/mc/Ch-unifrng.pdf>  
(not examinable)

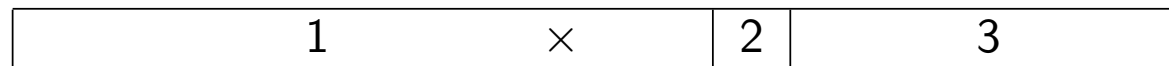
# Sampling for univariate discrete random variables

(Based on a slide from David Barber)

- ▶ Consider the one dimensional discrete distribution  $p(x)$  with  $x \in \{1, 2, 3\}$ , with

$$p(x) = \begin{cases} 0.6 & x = 1 \\ 0.1 & x = 2 \\ 0.3 & x = 3 \end{cases}$$

- ▶ Divide  $[0, 1]$  into chunks  $[0, 0.6)$ ,  $[0.6, 0.7)$ ,  $[0.7, 1]$



- ▶ We then draw a sample  $u$  uniformly from  $[0, 1]$
- ▶ We return the label of the partition in which  $u$  fell.
- ▶ Example: if  $u = 0.53$ , we return the sample “1”

# Sampling for univariate continuous random variables

- ▶ A similar method as the one above exists for continuous random variables.
- ▶ Called inverse transform sampling.
- ▶ Recall: the cumulative distribution function (cdf) of a random variable  $x$  with pdf  $p_x$  is

$$F_x(\alpha) = \mathbb{P}(x \leq \alpha) = \int_{-\infty}^{\alpha} p_x(u) du$$

- ▶ To generate  $n$  iid samples from  $x$  with cdf  $F_x$ :
  - ▶ calculate the inverse  $F_x^{-1}$
  - ▶ sample  $n$  iid random variables uniformly distributed on  $[0, 1]$ :  
 $y_i \sim \mathcal{U}(0, 1)$ ,  $i = 1, \dots, n$ .
  - ▶ transform each sample by  $F_x^{-1}$ :  $x_i = F_x^{-1}(y_i)$ ,  $i = 1, \dots, n$ .

(see Tutorial 8 for derivation)



# Basic principle of rejection sampling

- ▶ Assume you can draw iid samples  $\mathbf{x}_i \sim q(\mathbf{x})$ .
- ▶ For each sampled  $\mathbf{x}_i$ , you draw a Bernoulli random variable  $y_i \in \{0, 1\}$  whose success probability depends on  $\mathbf{x}_i$

$$\mathbb{P}(y_i = 1 | \mathbf{x}_i) = f(\mathbf{x}_i)$$

- ▶ You get samples  $(y_i, \mathbf{x}_i)$  with joint distribution

$$q(\mathbf{x})f(\mathbf{x})^y(1 - f(\mathbf{x}))^{(1-y)}$$

- ▶ Conditional pdf of  $\mathbf{x} | y = 1$  is proportional to  $q(\mathbf{x})f(\mathbf{x})$
- ▶ Keep or “accept” the  $\mathbf{x}_i$  with  $y_i = 1$ , “reject” those with  $y_i = 0$ .
- ▶ Accepted samples follow

$$\mathbf{x}_i \sim \frac{q(\mathbf{x})f(\mathbf{x})}{\int q(\mathbf{x})f(\mathbf{x})d\mathbf{x}}$$

# Sampling from the posterior by rejection sampling

- ▶ Conditional acceptance probability  $f(\mathbf{x}) \in [0, 1]$  can be used to shape the distribution of the samples from  $q(\mathbf{x})$
- ▶ Consider Bayesian inference: prior  $p(\boldsymbol{\theta})$ , likelihood  $L(\boldsymbol{\theta})$
- ▶ Using  $L(\boldsymbol{\theta})/(\max L(\boldsymbol{\theta}))$  as acceptance probability  $f$  transforms the samples  $\boldsymbol{\theta}_i$  from the prior into samples from the posterior.
- ▶ Accepted parameters follow

$$\boldsymbol{\theta}_i \sim \frac{p(\boldsymbol{\theta})L(\boldsymbol{\theta})}{\int p(\boldsymbol{\theta})L(\boldsymbol{\theta})d\boldsymbol{\theta}} = p(\boldsymbol{\theta}|\mathcal{D})$$

- ▶ More likely parameter configurations are more likely accepted.

# Sampling from the posterior by rejection sampling

- ▶ For discrete random variables  $L(\theta) = \mathbb{P}(\mathbf{x} = \mathcal{D}; \theta) \in [0, 1]$ .
- ▶ Accepting a  $\theta_i$  with probability  $L(\theta)$  can be implemented by checking whether data simulated from the model with parameter value  $\theta_i$  equals the observed data.
- ▶ Samples from the posterior = samples from the prior that produce data equal to the observed one.  
(see slides “Basic of Model-Based Learning”)

Side-note (not examinable): enables Bayesian inference when the likelihood is intractable (e.g. due to unobserved variables) but sampling from the model is possible. Forms the basis of a set of methods called approximate Bayesian computation, for an introductory review paper see <https://michaelgutmann.github.io/assets/papers/Lintusaari2017.pdf>.

# Standard formulation of rejection sampling

- ▶ Rejection sampling is typically presented (slightly) differently.
- ▶ Goal is to generate samples from a target distribution  $p(\mathbf{x})$  known up to normalisation constant when being able to sample from  $q(\mathbf{x})$ .
- ▶ Since accepted samples follow

$$\mathbf{x}_i \sim \frac{q(\mathbf{x})f(\mathbf{x})}{\int q(\mathbf{x})f(\mathbf{x})d\mathbf{x}}$$

choose conditional acceptance probability  $f(\mathbf{x}) \propto p(\mathbf{x})/q(\mathbf{x})$

- ▶ See Barber 27.1.2.

# Multivariate by univariate sampling

- ▶ Rejection sampling is limited to low-dimensional cases (see Barber 27.1.2)
- ▶ Sampling from high-dimensional multivariate distributions is generally difficult.
- ▶ One way to approach the problem of multivariate sampling is to translate it into the task of solving several lower dimensional sampling problems.
- ▶ Examples:
  - ▶ Ancestral sampling
  - ▶ Gibbs sampling

# Ancestral sampling

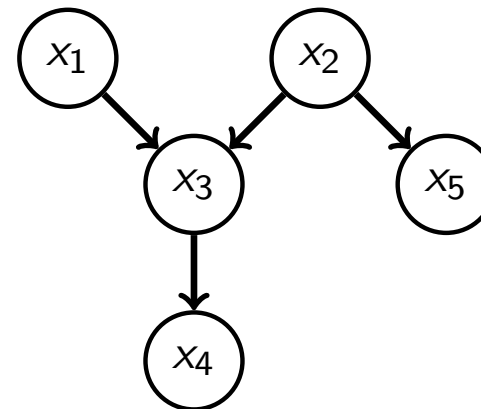
- ▶ Factorisation provides a recipe for data generation / sampling from  $p(\mathbf{x})$

- ▶ Example:

$$p(x_1, \dots, x_5) = p(x_1)p(x_2)p(x_3|x_1, x_2)p(x_4|x_3)p(x_5|x_2)$$

- ▶ We can generate samples from the joint distribution  $p(x_1, x_2, x_3, x_4, x_5)$  by sampling

1.  $x_1 \sim p(x_1)$
2.  $x_2 \sim p(x_2)$
3.  $x_3 \sim p(x_3|x_1, x_2)$
4.  $x_4 \sim p(x_4|x_3)$
5.  $x_5 \sim p(x_5|x_2)$



- ▶ Sets of univariate sampling problems.

# Gibbs sampling

(Based on a slide from David Barber)

- ▶ Gibbs sampling also reduces the problem of multivariate sampling to the problem of univariate sampling.
- ▶ Goal: generate samples  $\mathbf{x}^{(k)}$  from  $p(\mathbf{x}) = p(x_1, \dots, x_d)$ .
- ▶ By product rule

$$\begin{aligned} p(\mathbf{x}) &= p(x_i | x_1, \dots, x_{i-1}, x_{i+1}, \dots, x_d) p(x_1, \dots, x_{i-1}, x_{i+1}, \dots, x_d) \\ &= p(x_i | \mathbf{x}_{\setminus i}) p(\mathbf{x}_{\setminus i}) \end{aligned}$$

- ▶ Given a joint initial state  $\mathbf{x}^{(1)}$  from which we can read off the ‘parental’ state  $\mathbf{x}_{\setminus i}^{(1)}$

$$\mathbf{x}_{\setminus i}^{(1)} = (x_1^{(1)}, \dots, x_{i-1}^{(1)}, x_{i+1}^{(1)}, \dots, x_d^{(1)}),$$

we can draw a sample  $x_i^{(2)}$  from  $p(x_i | \mathbf{x}_{\setminus i}^{(2)})$ .

- ▶ We assume this distribution is easy to sample from since it is univariate.

# Gibbs sampling

(Based on a slide from David Barber)

- ▶ We call the new joint sample in which only  $x_i$  has been updated  $\mathbf{x}^{(2)}$ ,

$$\mathbf{x}^{(2)} = (x_1^{(1)}, \dots, x_{i-1}^{(1)}, x_i^{(2)}, x_{i+1}^{(1)}, \dots, x_d^{(1)}).$$

- ▶ One then selects another variable  $x_j$  to sample and, by continuing this procedure, generates a set  $\mathbf{x}^{(1)}, \dots, \mathbf{x}^{(n)}$  of samples in which each  $\mathbf{x}^{(k+1)}$  differs from  $\mathbf{x}^{(k)}$  in only a single component.
- ▶ Since  $p(x_i | \mathbf{x}_{\setminus i}) = p(x_i | \text{MB}(x_i))$ , we can sample from  $p(x_i | \text{MB}(x_i))$  which is easier.  
( $\text{MB}(x_i)$  denotes the Markov blanket of  $x_i$ , see slides on directed and undirected graphical models.)
- ▶ Samples are not independent.
- ▶ Gibbs sampling is an example of a Markov chain Monte Carlo method (see Barber 27.3 and 27.4).



# Program recap

## 1. Monte Carlo integration

- Approximating expectations by averages
- Importance sampling

## 2. Sampling

- Simple univariate sampling
- Rejection sampling
- Ancestral sampling
- Gibbs sampling