Sampling and Monte Carlo Integration

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Recap

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

Marginalisation

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

Expectations

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

for some function g.

For unobserved variables, likelihood and gradient of the log lik

$$egin{aligned} L(m{ heta}) &= p(\mathcal{D}; m{ heta}) = \int_{\mathbf{u}} p(\mathbf{u}, \mathcal{D}; m{ heta} \mathrm{d}\mathbf{u}), \
abla_{m{ heta}} \ell(m{ heta}) &= \mathbb{E}_{p(\mathbf{u}|\mathcal{D}; m{ heta})} \left[
abla_{m{ heta}} \log p(\mathbf{u}, \mathcal{D}; m{ heta})
ight] \end{aligned}$$

Recap

For unnormalised models with intractable partition functions

$$egin{aligned} \mathcal{L}(oldsymbol{ heta}) &= rac{ ilde{p}(\mathcal{D};oldsymbol{ heta})}{\int_{\mathbf{x}} ilde{p}(\mathbf{x};oldsymbol{ heta})\mathrm{d}\mathbf{x}} \
abla_{oldsymbol{ heta}}\ell(oldsymbol{ heta}) &\propto \mathbf{m}(\mathcal{D};oldsymbol{ heta}) - \mathbb{E}_{p(\mathbf{x};oldsymbol{ heta})}\left[\mathbf{m}(\mathbf{x};oldsymbol{ heta})
ight] \end{aligned}$$

- 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 and expectations using sampling.

Program

- 1. Monte Carlo integration
- 2. Sampling

Program

1. Monte Carlo integration

- Approximating expectations by averages
- Importance sampling
- Effective sample size

2. Sampling

Averages with iid samples

From exercises): 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 $x \sim p(x)$.

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

- ightharpoonup 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

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

(*: "identically distributed" assumption is used, not independence)

Variability

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

(*: independence assumption used)

 \triangleright Expected squared error decreases as 1/n

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

Averages with iid samples

Weak law of large numbers:

$$\mathbb{P}\left(|\bar{x}_n - \mathbb{E}[x]| \ge \epsilon\right) \le \frac{\mathbb{V}[x]}{n\epsilon^2}$$

- As $n \to \infty$, the probability for the sample average to deviate from the expected value goes to zero if the variance is finite.
- 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 $\mathbb{V}[s]$.

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

- ► This means that for *all* random variables with finite mean and variance:
 - 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.

Proofs (not examinable)

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

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

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

$$\mathbb{P}\left(|s - \mathbb{E}[s]| \ge t\right) = \mathbb{P}\left((s - \mathbb{E}[s])^2 \ge t^2\right) \ \le \frac{\mathbb{E}\left[(s - \mathbb{E}[s])^2\right]}{t^2}.$$

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,

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

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

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

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

$$\mathbb{E}[y] \ge \int_0^t u p(u) du + t \mathbb{P}(y \ge t)$$

 $\ge t \mathbb{P}(y \ge t),$

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

$$\mathbb{P}(y \ge t) \le \frac{\mathbb{E}(y)}{t} \quad (t > 0)$$

Averages with correlated samples

When computing the variance of the sample average

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

we assumed the samples are identically and independently distributed.

- ▶ The variance shrinks with increasing n and the average becomes more and more concentrated around $\mathbb{E}[x]$.
- ightharpoonup Corresponding results exist for the case of statistically dependent samples x_i . Known as "ergodic theorems".
- Out of scope for PMR but important for the theory of Markov chain Monte Carlo methods.

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 $\frac{1}{n}\mathbb{V}[g(\mathbf{x})]$

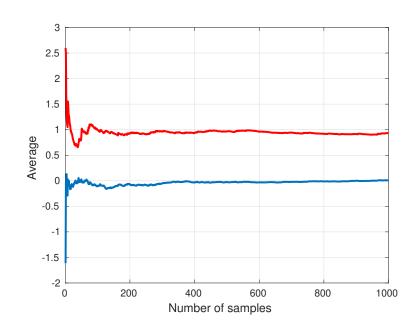
Example (Based on a slide from Amos Storkey)

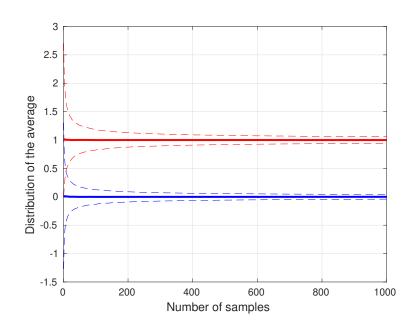
$$\mathbb{E}[g(x)] = \int g(x) \mathcal{N}(x; 0, 1) dx \approx \frac{1}{n} \sum_{i=1}^{n} g(x_i) \qquad (x_i \sim \mathcal{N}(x; 0, 1))$$

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)



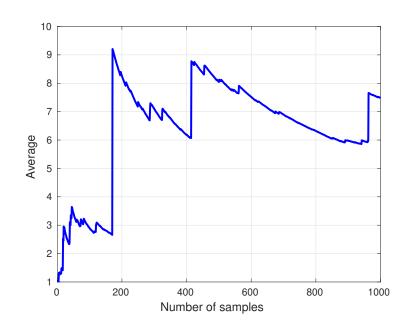


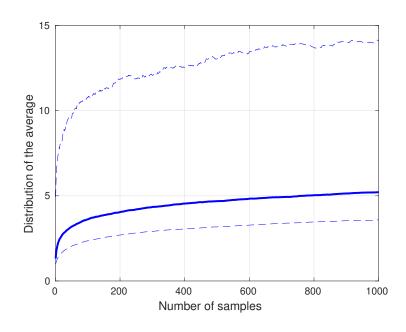
Example (Based on a slide from Amos Storkey)

$$\mathbb{E}[g(x)] = \int g(x)\mathcal{N}(x;0,1)\mathrm{d}x \approx \frac{1}{n}\sum_{i=1}^n g(x_i) \qquad (x_i \sim \mathcal{N}(x;0,1))$$
 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:
 - \triangleright Strong fluctuations in the sample average as n increases.
 - Large non-declining variability.
- ► Note: integral is not finite:

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

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!

Importance sampling to approximate 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*

$$I = \int g(\mathbf{x}) d\mathbf{x} = \int g(\mathbf{x}) rac{q(\mathbf{x})}{q(\mathbf{x})} d\mathbf{x}$$
 (assume $q(\mathbf{x}) > 0$ when $g(\mathbf{x}) > 0$)
$$= \int rac{g(\mathbf{x})}{q(\mathbf{x})} q(\mathbf{x}) d\mathbf{x}$$

$$= \mathbb{E}_{q(\mathbf{x})} \left[rac{g(\mathbf{x})}{q(\mathbf{x})}
ight]$$

$$pprox rac{1}{n} \sum_{i=1}^{n} rac{g(\mathbf{x}_{i})}{q(\mathbf{x}_{i})}$$

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

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

Choice of the importance distribution

ightharpoonup Call the approximation \hat{I} ,

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

 $\triangleright \hat{I}$ is unbiased by construction

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

Variance

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

Depends on the second moment.

Choice of the importance distribution

► The second moment is

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

- ▶ 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

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

Poptimal 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

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

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

$$\mathbb{E}\left[\left(rac{g(\mathsf{x})}{q^*(\mathsf{x})}
ight)^2
ight] = \mathbb{E}\left[\left|rac{g(\mathsf{x})}{q^*(\mathsf{x})}
ight|
ight]^2$$

Proof (not examinable)

Indeed, for the optimal q, we have

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

$$= \int |g(\mathbf{x})| d\mathbf{x} \int |g(\mathbf{x})|^2 \frac{1}{|g(\mathbf{x})|} d\mathbf{x}$$

$$= \left(\int |g(\mathbf{x})| d\mathbf{x}\right)^2$$

and

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

which concludes the proof.

Importance sampling to approximate the partition function

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

$$egin{aligned} Z(oldsymbol{ heta}) &= \int ilde{p}(\mathbf{x};oldsymbol{ heta}) \mathrm{d}\mathbf{x} \ &= \int ilde{p}(\mathbf{x};oldsymbol{ heta}) rac{q(\mathbf{x})}{q(\mathbf{x})} \mathrm{d}\mathbf{x} \qquad ext{(assume } q(\mathbf{x}) > 0 ext{ when } ilde{p}(\mathbf{x}) > 0) \ &= \int rac{ ilde{p}(\mathbf{x};oldsymbol{ heta})}{q(\mathbf{x})} q(\mathbf{x}) \mathrm{d}\mathbf{x} \ &pprox rac{1}{n} \sum_{i=1}^n rac{ ilde{p}(\mathbf{x}_i;oldsymbol{ heta})}{q(\mathbf{x}_i)} \qquad ext{(}\mathbf{x}_i \sim q(\mathbf{x}) ext{ iid)} \end{aligned}$$

Example

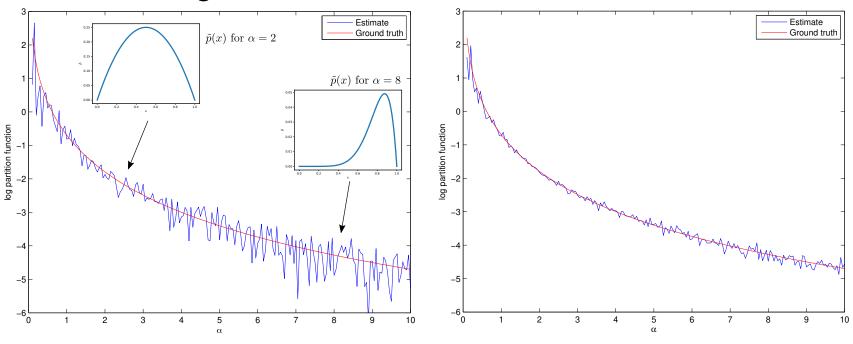
Approximating the log partition function of the unnormalised beta-distribution

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

for β fixed to $\beta = 2$.

Importance distribution: uniform distribution on [0,1]

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



Importance sampling to approximate 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

$$egin{aligned} \mathbb{E}_{p(\mathbf{x})}[g(\mathbf{x})] &= \int g(\mathbf{x}) p(\mathbf{x}) \mathrm{d}\mathbf{x} \ &= \int g(\mathbf{x}) rac{p(\mathbf{x})}{q(\mathbf{x})} q(\mathbf{x}) \mathrm{d}\mathbf{x} & ext{(assume } q(\mathbf{x}) > 0 ext{ when } g(\mathbf{x}) p(\mathbf{x}) > 0) \ &= \mathbb{E}_{q(\mathbf{x})} \left[g(\mathbf{x}) rac{p(\mathbf{x})}{q(\mathbf{x})}
ight] \ &pprox rac{1}{n} \sum_{i=1}^n g(\mathbf{x}_i) rac{p(\mathbf{x}_i)}{q(\mathbf{x}_i)} \end{aligned}$$

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.

Self/auto-normalised importance sampling

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

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

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

$$= \frac{\mathbb{E}_{q(\mathbf{x})}\left[g(\mathbf{x})\frac{\tilde{p}(\mathbf{x})}{q(\mathbf{x})}\right]}{\mathbb{E}_{q(\mathbf{x})}\left[\frac{\tilde{p}(\mathbf{x})}{q(\mathbf{x})}\right]}$$

Self/auto-normalised importance sampling

Since

$$\int g(\mathbf{x}) p(\mathbf{x}) d\mathbf{x} = rac{\mathbb{E}_{q(\mathbf{x})} \left[g(\mathbf{x}) rac{ ilde{p}(\mathbf{x})}{q(\mathbf{x})}
ight]}{\mathbb{E}_{q(\mathbf{x})} \left[rac{ ilde{p}(\mathbf{x})}{q(\mathbf{x})}
ight]} = rac{\mathbb{E}_{q(\mathbf{x})} \left[g(\mathbf{x}) rac{ ilde{p}(\mathbf{x})}{ ilde{q}(\mathbf{x})}
ight]}{\mathbb{E}_{q(\mathbf{x})} \left[rac{ ilde{p}(\mathbf{x})}{ ilde{q}(\mathbf{x})}
ight]}$$

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

Approximate both expectations by a sample average

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

where
$$w_i = rac{ ilde{p}(\mathbf{x}_i)}{ ilde{q}(\mathbf{x}_i)}$$
 and $\mathbf{x}_i \sim q(\mathbf{x})$ (iid)

Self/auto-normalised importance sampling

$$w_i = rac{ ilde{p}(\mathbf{x}_i)}{ ilde{q}(\mathbf{x}_i)}, \ \mathbf{x}_i \stackrel{\mathrm{iid}}{\sim} q(\mathbf{x})$$

Called self-normalised or auto-normalised importance sampling

$$\int g(\mathbf{x})p(\mathbf{x})d\mathbf{x} \approx \sum_{i=1}^{n} W_{i}g(\mathbf{x}_{i}) \qquad W_{i} = \frac{W_{i}}{\sum_{k=1}^{n} W_{k}}$$

Note: $\sum_{i=1}^{n} W_i = 1$

▶ Interpretation in terms of a Dirac-delta approximation of $p(\mathbf{x})$,

$$p(\mathbf{x}) \approx \sum_{i=1}^{n} W_i \delta_{\mathbf{x}_i}(\mathbf{x})$$
 $\mathbf{x}_i \stackrel{\text{iid}}{\sim} q(\mathbf{x})$

(\equiv mixture of Gaussians with mixture probabilities W_i , expected values \mathbf{x}_i , and infinitesimally small variances)

Effective sample size

$$w_i = rac{ ilde{p}(\mathsf{x}_i)}{ ilde{q}(\mathsf{x}_i)}, \; \mathsf{x}_i \stackrel{\mathsf{iid}}{\sim} q(\mathsf{x})$$

If the weights w_i are constants, the weighted average $\sum_{i=1}^{n} W_i g(\mathbf{x}_i)$ becomes a standard average

$$W_i = \frac{w_i}{\sum_{k=1}^{n} w_k} \stackrel{w_i = c}{=} \frac{c}{\sum_{k=1}^{n} c} = \frac{1}{N}$$

▶ But the w_i are typically not all equal, so that some x_i contribute more to the average than others, e.g.

$$w_1 = 10^6, w_k = 1, k > 1 \Longrightarrow W_1 \approx 1, W_k \approx 0, k > 1$$

We would effectively "average" over 1 data point!

► When working with a weighted average, always compute the "effective sample size" (ESS),

ESS =
$$\frac{\left(\sum_{i=1}^{n} w_i\right)^2}{\sum_{i=1}^{n} w_i^2} = \frac{1}{\sum_{i=1}^{n} W_i^2} \in [1, N]$$

Small ESS means the average is unreliable (high variance).

Program

1. Monte Carlo integration

- Approximating expectations by averages
- Importance sampling
- Effective sample size

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]



- \triangleright We then draw a sample u uniformly from [0,1]
- \triangleright We return the label of the partition in which u fell.
- ightharpoonup 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 :
 - ightharpoonup calculate the inverse F_{\times}^{-1}
 - sample n iid random variables uniformly distributed on [0,1]: $y_i \sim \mathcal{U}(0,1), i = 1, \ldots, n$.
 - ransform each sample by F_x^{-1} : $x_i = F_x^{-1}(y_i)$, i = 1, ..., n.

(see exercises 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

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

 \triangleright 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/"accept" the \mathbf{x}_i with $y_i = 1$, "reject" those with $y_i = 0$.
- Accepted samples follow

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

Denominator equals the marginal probability of acceptance

$$\mathbb{P}(y=1) = \mathbb{E}_{q(\mathbf{x})} \mathbb{P}(y=1|\mathbf{x}) = \int q(\mathbf{x}) f(\mathbf{x}) d\mathbf{x}$$

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 $q(\mathbf{x})$
- ightharpoonup Consider Bayesian inference: prior $p(\theta)$, likelihood $L(\theta)$
- ▶ Using $L(\theta)/(\max L(\theta))$ as acceptance probability f transforms the samples θ_i from the prior into samples from the posterior.
- Accepted parameters follow

$$egin{aligned} heta_i &\sim rac{p(heta)L(heta)}{\int p(heta)L(heta)\mathrm{d} heta} = p(heta|\mathcal{D}) \end{aligned}$$

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 θ_i with probability $L(\theta)$ can be implemented by checking whether data simulated from the model with parameter value θ_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 $p(\mathbf{x})$ 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 $f(\mathbf{x}) \propto p(\mathbf{x})/q(\mathbf{x})$

▶ To determine the proportionality factor, note that $f(\mathbf{x}) \leq 1$ since it is a conditional probability. Hence:

$$f(\mathbf{x}) = \frac{1}{M} \frac{p(\mathbf{x})}{q(\mathbf{x})}$$
 $M = \max_{\mathbf{x}} \frac{p(\mathbf{x})}{q(\mathbf{x})}$

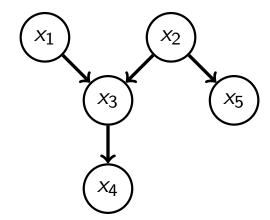
Acceptance probability: $\mathbb{P}(y=1) = \int q(\mathbf{x}) f(\mathbf{x}) d\mathbf{x} = \frac{1}{M}$.

Multivariate by univariate sampling

- ▶ Rejection sampling may scale poorly because *M* increases with dimensionality so that acceptance probability goes down.
- 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: $p(x_1,...,x_5) = p(x_1)p(x_2)p(x_3|x_1,x_2)p(x_4|x_3)p(x_5|x_2)$
- 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

$$p(\mathbf{x}) = p(x_i|x_1, \dots, x_{i-1}, x_{i+1}, \dots, x_d)p(x_1, \dots, x_{i-1}, x_{i+1}, \dots, x_d)$$

= $p(x_i|\mathbf{x}_{\setminus i})p(\mathbf{x}_{\setminus i})$

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

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

we can draw a sample $x_i^{(2)}$ from $p(x_i|\mathbf{x}_{\setminus i}^{(1)})$.

► 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 $p(x_i|\mathbf{x}_{\setminus i}) = p(x_i|\mathrm{MB}(x_i))$, we can sample from $p(x_i|\mathrm{MB}(x_i))$ which is easier. (MB(x_i) is the Markov blanket of x_i)
- Samples are not independent.
- ► Gibbs sampling is an example of a Markov chain Monte Carlo method (see Barber 27.4 and 27.3.1, and the exercises).

Program recap

1. Monte Carlo integration

- Approximating expectations by averages
- Importance sampling
- Effective sample size

2. Sampling

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