

# Projet de Monte Carlo P3

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## Question 23

Let's determine a rejection algorithm which produces a sample that follows the distribution of the random variable  $X$  conditioned on the event  $X \in A$ .

**Conditional density** Let  $X$  be a random variable with a density  $f_X(x)$  defined on a measurable space. The conditional density of  $X$  given  $X \in A$ , where  $A$  is a measurable subset of the space, is proportional to  $f_X(x)$  restricted to  $A$ . Specifically:

$$f_{X|X \in A}(x) \propto f_X(x)\mathbf{1}_A(x),$$

This ensures that the conditional density  $f_{X|X \in A}$  is obtained by restricting  $f_X$  to the set  $A$ , then normalizing it so that it integrates to 1.

**Rejection algorithm principle** The rejection algorithm proceeds as follows:

1. Generate a sequence of variables  $X_1, X_2, \dots$  i.i.d. according to the density  $f_X(x)$ .
2. Accept a sample  $X_i$  if and only if  $X_i \in A$ .
3. Collect the accepted samples as a new sequence  $\{Y_k\}$ , where  $Y_k$  is the  $k$ -th accepted sample.

**proof:** Each sample  $X_i$  is distributed according to  $f_X(x)$ . The rejection step ensures that only samples within  $A$  are retained. Since the density  $f_{X|X \in A}$  is proportional to  $f_X(x)$  restricted to  $A$ , the accepted samples  $\{Y_k\}$  are distributed according to:

$$f_{X|X \in A}(x) \propto f_X(x)\mathbf{1}_A(x).$$

The normalization is automatically achieved because the algorithm accepts samples with probability proportional to their density under  $f_X(x)$  within the set  $A$ .

**Independence of accepted samples:** The samples  $X_1, X_2, \dots$  are independent and identically distributed. Consequently, the accepted samples  $Y_k$  are also independent and identically distributed, following the conditional density  $f_{X|X \in A}(x)$ .

**Conclusion** The rejection algorithm is valid for simulating samples from the conditional distribution of  $X$  given  $X \in A$ , as it effectively restricts the samples to the set  $A$  and ensures that the resulting distribution is proportional to  $f_X(x)$  restricted to  $A$ , with proper normalization. This normalization actually gives us  $\mathbb{P}(X \in A)$ .

**Remark:** With respect to the question 24, we assume that  $\mathbb{P}(X \in A)$  is unknown. If it were known, we could have done a classical reject-algorithm to simulate a sample directly from the density  $f_{X|X \in A}(x)$ . This method is pretty naive and inefficient, especially for events of small probability. We will look at more effective methods in parts 2 and 3 of the project.

## Question 26

Let  $g(x)$  be the chosen instrumental density, a Cauchy distribution of parameters q and 1. Our choice is motivated by the fact the density g of a Cauchy is strictly positive for all real numbers, therefore:

$$\text{supp}(h \cdot f) \subseteq \text{supp}(g)$$

Furthermore, the Cauchy density of median=q is q-symmetric, so we can use an antithetic approach "Z.antithetic = -Z + 2\*q" and "1/(Z-q) + q" to reduce the variance of our estimator. The importance sampling estimator for  $h(X) = \mathbf{1}_{\{X>q\}}$  is derived as follows:

The expectation of  $h(X)$  under  $f$  can be written as:

$$\mathbb{E}_f[h(X)] = \int_{\mathbb{R}} h(x)f(x) dx = \int_q^{\infty} f(x) dx = \mathbb{P}(X > q).$$

Using importance sampling, we rewrite the integral in terms of the instrumental density  $g(x)$ :

$$\mathbb{E}_f[h(X)] = \int_q^{\infty} \frac{f(x)}{g(x)} g(x) dx.$$

The corresponding importance sampling estimator, based on  $n$  i.i.d. samples  $Z_1, Z_2, \dots, Z_n$  drawn from  $g(x)$ , is:

$$\hat{\mathbb{E}}_f[h(X)] = \frac{1}{n} \sum_{i=1}^n h(Z_i) \frac{f(Z_i)}{g(Z_i)} = \frac{1}{n} \sum_{i=1}^n \mathbf{1}(Z_i > q) \frac{f(Z_i)}{g(Z_i)}.$$

**Properties:** The estimator  $\hat{\mathbb{E}}_f[h(X)]$  is unbiased and converges to  $\mathbb{E}_f[h(X)]$  as  $n \rightarrow \infty$ .

## Question 27

The probability density function of a random variable following a Cauchy distribution with parameters  $x_0$  (the location parameter, representing the median and mode) and  $\gamma > 0$  (the scale parameter) is given by:

$$f(x; x_0, \gamma) = \frac{1}{\pi \gamma \left[ 1 + \left( \frac{x-x_0}{\gamma} \right)^2 \right]}$$

As mentioned in question 26, the choice of parameters q and 1 comes from the symmetric properties of the Cauchy distribution and the fact that it gives us bigger tails than the function f.

## Question 29

The score is the gradient of the log-likelihood function with respect to the parameter vector:  $\log \mathcal{L}(\Theta; x)$ . In other words the score is :

$$s(\Theta, x) = \nabla_{\Theta} \log \mathcal{L}(\Theta; x)$$

The likelihood function for the data  $x_1, \dots, x_n$  is given by:

$$\mathcal{L}(\mu_1, \mu_2, \sigma_1^2, \sigma_2^2 | x_1, \dots, x_n) = \prod_{i=1}^n \frac{1}{1-a} (f_1(x_i) - af_2(x_i))$$

where:

$$f_1(x_i) = \frac{1}{\sqrt{2\pi\sigma_1^2}} \exp\left(-\frac{(x_i - \mu_1)^2}{2\sigma_1^2}\right)$$

$$f_2(x_i) = \frac{1}{\sqrt{2\pi\sigma_2^2}} \exp\left(-\frac{(x_i - \mu_2)^2}{2\sigma_2^2}\right)$$

Taking the logarithm of the likelihood:

$$\log \mathcal{L}(\mu_1, \mu_2, \sigma_1^2, \sigma_2^2 | x_1, \dots, x_n) = \sum_{i=1}^n \log \left[ \frac{1}{1-a} (f_1(x_i) - af_2(x_i)) \right]$$

This simplifies to:

$$-n \log(1-a) - n \log(\sqrt{2\pi}) + \sum_{i=1}^n \log \left( \frac{1}{\sigma_1} \exp\left(-\frac{(x_i - \mu_1)^2}{2\sigma_1^2}\right) - \frac{a}{\sigma_2} \exp\left(-\frac{(x_i - \mu_2)^2}{2\sigma_2^2}\right) \right)$$

We compute the derivative of the log-likelihood with respect to  $\mu_1$ :

Expanding the derivative:

$$\frac{\partial}{\partial \mu_1} f_1(x_i) = \frac{x_i - \mu_1}{\sigma_1^3} \exp\left(-\frac{(x_i - \mu_1)^2}{2\sigma_1^2}\right)$$

$$\frac{\partial}{\partial \mu_1} f_2(x_i) = 0 \quad (\text{since } f_2 \text{ depends on } \mu_2, \text{ not } \mu_1)$$

$$\frac{\partial \log \mathcal{L}(x_1, \dots, x_n | \mu_1, \mu_2, \sigma_1^2, \sigma_2^2)}{\partial \mu_1} = \sum_{i=1}^n \frac{\frac{(x_i - \mu_1)}{\sigma_1^3} \exp\left(-\frac{(x_i - \mu_1)^2}{2\sigma_1^2}\right)}{\frac{1}{\sigma_1} \exp\left(-\frac{(x_i - \mu_1)^2}{2\sigma_1^2}\right) - a \frac{1}{\sigma_2} \exp\left(-\frac{(x_i - \mu_2)^2}{2\sigma_2^2}\right)}$$

The score function for  $\mu_1$  is defined as:

$$s_{\mu_1}(x | \theta_1, \theta_2) = \sum_{i=1}^n \frac{\frac{(x_i - \mu_1)}{\sigma_1^3} \exp\left(-\frac{(x_i - \mu_1)^2}{2\sigma_1^2}\right)}{\frac{1}{\sigma_1} \exp\left(-\frac{(x_i - \mu_1)^2}{2\sigma_1^2}\right) - a \frac{1}{\sigma_2} \exp\left(-\frac{(x_i - \mu_2)^2}{2\sigma_2^2}\right)}.$$

## Question 30

Let us take  $h_0 : x \mapsto \frac{\partial \log \mathcal{L}(x_1, \dots, x_n | \mu_1, \mu_2, \sigma_1^2, \sigma_2^2)}{\partial \mu}$ .

Then a control variate Monte Carlo Estimator is :

$$\hat{\delta}_n^{cv}(b) = \frac{1}{n} \sum_{i=1}^n 1_{X_i \geq q} - b[h_0(X_i) - m].$$

Since the expectation of the score function is 0, we have  $m = 0$ .

Thus :

$$\widehat{\delta}_n^{cv}(b) = \frac{1}{n} \sum_{i=1}^n \mathbb{1}_{X_i \geq q} - bh_0(X_i).$$

## Question 32

### 1 Naive method

The computational cost of this method can be evaluated based on the number of samples  $n$ . Specifically, the variance of the naive estimator is given by:

$$\hat{\theta}_{\text{naive}} = \frac{1}{n} \sum_{i=1}^n f(X_i),$$

where  $f(X_i)$  is an indicator function that equals 1 if  $X_i \geq q$  and 0 otherwise. The variance of this estimator can be calculated as:

$$\text{Var}(\hat{\theta}_{\text{naive}}) = \frac{1}{n} \cdot \text{Var}(X_i).$$

The computational cost of this method is also evaluated using `microbenchmark`, and by calculating the median execution time for a large number of samples  $n$ , we obtain a cost  $R_1 = 47758532$  (in time units per sample). This cost is based on the observed data from our experiment.

### 2 Importance Sampling Method

In our code, the function  $f(x)$  is defined as the difference between two densities, and the estimator based on importance sampling is computed using Cauchy weights. This method involves a weighting distribution  $g(x) = \text{Cauchy}(x, q, 1)$ , and the estimator is given by:

$$\hat{\theta}_{\text{IS}} = \frac{1}{n} \sum_{i=1}^n \frac{f(Z_i)}{g(Z_i)} \mathbb{1}(Z_i \geq q),$$

where  $Z_i$  are samples generated from a Cauchy distribution. The computational cost of this method is also evaluated using `microbenchmark`, and we obtain a cost  $R_2 = 532252.2$ .

### 3 Control Variate Method

The control variate method relies on the use of an auxiliary variable that is correlated with the variable of interest but has a known expectation. This method reduces the variance of the estimator by adjusting the samples based on the control variable.

The computational cost of this method is also evaluated using `microbenchmark`, and we obtain a cost  $R_3 = 401456$ .

### 4 Comparison of Methods

We now compare the three methods in terms of computational cost and estimation accuracy. The cost  $R_1$  of the naive method is compared to  $R_2$  for importance sampling and  $R_3$  for the control variate method. The following ratios are calculated to assess the relative efficiency of the methods:

$$\frac{R_1}{R_2} = \frac{47758532}{532252.2} \approx 89.73$$

This ratio shows that the importance sampling estimator is more efficient than the naive estimator in terms of computational cost, as  $\frac{R_1}{R_2} > 1$ .

$$\frac{R_2}{R_3} = \frac{532252.2}{401456} \approx 1.33$$

This ratio shows that the control variate method outperforms the importance sampling method in terms of computational cost, as  $\frac{R_2}{R_3} > 1$ .