## Efficient Computation of Upper Probabilities of Failure Using Monte Carlo Simulation and Reweighting Techniques

**Thomas Fetz** 

THOMAS.FETZ@UIBK.AC.AT

Unit for Engineering Mathematics Institute of Basic Sciences in Engineering Science Technikerstraße 13, 6020 Innsbruck, Austria

## **Poster Abstract**

Let  $(X_{\lambda})_{\lambda \in \Lambda}$  be a family of random variables. The upper probability  $\overline{p}_f$  of failure is the solution of the optimization problem

$$\overline{p}_f = \max_{\lambda \in \Lambda} \int_D \mathbb{1}_{g(x) \le 0} f_{X_\lambda}(x) \, \mathrm{d}x$$

where  $f_{X_{\lambda}}$  is the corresponding probability density of the random variable  $X_{\lambda}, g : D \to \mathbb{R}$  a limit state function and 1 the indicator function. A value  $g(x) \leq 0$  means failure of the underlying engineering structure and a value g(x) > 0 means that the engineering structure is safe. As an example,  $f_{X_{\lambda}}$  is the density function of a Gaussian random variable  $X_{\lambda} \sim \mathcal{N}(\mu(\lambda), \Sigma(\lambda))$  with expectation  $\mu$  and covariance matrix  $\Sigma$  parametrized by  $\lambda = (\lambda_1, \ldots, \lambda_n) \in \Lambda$ .

The objective function  $p(\lambda) = \int_D \mathbb{1}_{g(x) \le 0} f_{X_\lambda}(x) dx$  of the above optimization problem can be approximated using Monte Carlo simulation together with reweighting or importance sampling techniques using only one single sample  $x_1, \ldots, x_N$  distributed according to a "basic" random variable  $X^0$  with  $f_{X^0} > 0$ :

$$p(\lambda) = \int_{D} \mathbb{1}_{g(x) \le 0} f_{X_{\lambda}}(x) \, \mathrm{d}x = \int_{D} \mathbb{1}_{g(x) \le 0} \frac{f_{X_{\lambda}}(x)}{f_{X^{0}}(x)} f_{X^{0}}(x) \, \mathrm{d}x$$
$$\approx \frac{1}{N} \sum_{k=1}^{N} \mathbb{1}_{g(x_{k}) \le 0} \frac{f_{X_{\lambda}}(x_{k})}{f_{X^{0}}(x_{k})} =: p_{x_{1},\dots,x_{N}}^{X^{0}}(\lambda),$$

cf. Fetz and Oberguggenberger (2016). Then an approximation of the upper probability  $\overline{p}_f$  of failure is obtained by  $\overline{p}_f \approx \max_{\lambda \in \Lambda} p_{x_1,\dots,x_N}^{X^0}(\lambda)$ . This method needs only N function evaluations  $g(x_1), \dots, g(x_N)$  of the limit state function which is an advantage in cases where the evaluation of g is time consuming, e.g. finite element computations. Further,  $p_{x_1,\dots,x_N}^{X^0}$  depends continuously on  $\lambda$  (if  $f_{X_\lambda}$  is continuous, too) which makes maximization easier. We also note that we get different functions  $p_{x_1,\dots,x_N}^{X^0}$  for different samples  $x_1,\dots,x_N$  and basic random variables  $X^0$ .

The purpose of the poster presentation is to discuss and compare variants of the above approach such as iterating and strategies for choosing the random variable  $X^0$ , and to exemplify these methods by an engineering example.

## References

T. Fetz and M. Oberguggenberger. Imprecise random variables, random sets, and Monte Carlo simulation. *International Journal of Approximate Reasoning*, 78:252–264, 2016.