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STRUCTURAL RELIABILITY ANALYSIS FOR P-BOXES USING MULTI-LEVEL META-MODELS

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Structural reliability analysis for p-boxes using multi-level meta-models

Roland Schöbi¹ and Bruno Sudret¹

¹*Chair of Risk, Safety and Uncertainty Quantification, ETH Zurich, Stefano-Franscini-Platz 5, 8093 Zurich, Switzerland*

Abstract

In modern engineering, computer simulations are a popular tool to analyse, design, and optimize systems. Furthermore, concepts of uncertainty and the related reliability analysis and robust design are of increasing importance. Hence, an efficient quantification of uncertainty is an important aspect of the engineer's workflow. In this context, the characterization of uncertainty in the input variables is crucial. In this paper, input variables are modelled by probability-boxes, which account for both aleatory and epistemic uncertainty. Two types of probability-boxes are distinguished: free and parametric (also called distributional) p-boxes. The use of probability-boxes generally increases the complexity of structural reliability analyses compared to traditional probabilistic input models. In this paper, the complexity is handled by two-level approaches which use Kriging meta-models with adaptive experimental designs at different levels of the structural reliability analysis. For both types of probability-boxes, the extensive use of meta-models allows for an efficient estimation of the failure probability at a limited number of runs of the performance function. The capabilities of the proposed approaches are illustrated through a benchmark analytical function and two realistic engineering problems.

Keywords: Structural reliability analysis – Kriging – probability-boxes – design enrichment – uncertainty quantification

1 Introduction

2 Introduction

Nowadays, computer simulations are a popular engineering tool to design systems of ever increasing complexity and to determine their performance. These simulations aim at reproducing the physical process and hence provide a solution to the underlying governing equations. As an example, finite

element models have become a standard tool in modern civil and mechanical engineering. Such models exploit the available computer power, meaning that a single run of the model can take up to hours of computation or more.

At the same time, uncertainties are intrinsic to these physical processes and to simulation models. Typically, the model parameters are not perfectly known, but inferred from data and then modelled probabilistically. However, a common situation in practice is to have a limited budget for data acquisition and to end up with scarce datasets. This introduces epistemic uncertainty (lack of knowledge) alongside aleatory uncertainty (natural variability). Then, a more general framework is required to characterize the uncertainty in model parameters beyond probability theory.

Imprecise probabilities describes a set of methodologies which are more general than the well-established probability theory. They include Dempster-Shafer's theory of evidence (Dempster, 1967; Shafer, 1976), Bayesian hierarchical models (Gelman, 2006; Gelman et al., 2009), possibility theory (De Cooman et al., 1995; Dubois and Prade, 1988), probability-boxes (Ferson and Ginzburg, 1996; Ferson and Hajagos, 2004), info-gap models (Ben-Haim, 2006; Kanno and Takewaki, 2006), and fuzzy sets (Zadeh, 1978; Möller and Beer, 2004). Among the various methods, probability-boxes provide a simple structure to distinguish aleatory and epistemic uncertainty naturally.

In this context, engineers are concerned with the reliability of their designs, which is assessed by structural reliability analysis. The goal of such an analysis is the estimation of the failure probability, which describes the probability that the systems performs below a minimal acceptable threshold value. A typical value for the failure probability in engineering is in the range of 10^{-3} to 10^{-6} . Altogether, the complexity of structural reliability analysis comprises three main factors: (i) a high-fidelity computer simulation, (ii) uncertainty in the parameters, and (iii) the estimation of rare events (*i.e.* the estimation of small failure probabilities).

In the presence of probabilistic variables, a number of methods are known to estimate failure probabilities (Lemaire, 2009; Morio et al., 2014), including Monte Carlo simulation, subset simulation (Au and Beck, 2001, 2003), importance sampling, first/second order reliability method (FORM/SORM) (Breitung, 1989; Hohenbichler and Rackwitz, 1988), and line sampling (Koutsourelakis et al., 2004; De Angelis et al., 2015). In the presence of imprecise probabilities, however, fewer publications are available, some of which are mentioned here. (Eldred and Swiler, 2009; Hurtado, 2013) use nested Monte Carlo simulations to cope with imprecise probabilities. (Alvarez and Hurtado, 2014) combine subset simulation with random sets. (Zhang et al., 2015) generalize FORM/SORM for Dempster-Shafer's evidence theory. A combination of line sampling and imprecise probabilities is discussed in (De Angelis et al., 2015). These publications typically require a large number of model evaluations and hence rely on an inexpensive-to-evaluate performance function to make the analysis tractable.

A way to reduce the number of model evaluations and hence allowing expensive-to-evaluate limit-state functions are meta-models. Meta-modelling techniques approximate the expensive-to-evaluate limit-state function by a fast-to-evaluate surrogate based on a relatively small number of runs of the original one. The meta-model is then used in the structural reliability analysis. In particular, meta-

models with adaptive experimental designs have been proven to be efficient in the estimation of failure probabilities. As an example, (Bichon et al., 2008; Echard et al., 2011; Dubourg et al., 2013) combine Kriging (Gaussian process models) (Santner et al., 2003) and structural reliability analysis in the context of probability theory.

However, only a few authors combine meta-modelling techniques with imprecise probabilities for structural reliability analysis. (Balesdent et al., 2014; Morio and Balesdent, 2016) use an adaptive-Kriging importance sampling algorithm for distributions with interval-valued parameters. A combination of Karhunen-Loève expansions and evidence theory is presented in (Oberguggenberger, 2014). (Zhang et al., 2014) use radial basis expansions in the presence of evidence theory.

Based on these developments, a set of algorithms is proposed in this paper to further increase the efficiency of structural reliability analysis in the presence of epistemic uncertainties. In particular, free and parametric p-boxes are presented separately and compared later in this paper. The application of Kriging meta-models at several stages of the imprecise structural reliability analysis promises a reduction of the overall computational costs at the same level of accuracy.

The paper is organized as follows. Section 3 defines probability-boxes as a generalization of probability theory. In Section 4, the basic structural reliability analysis is set up and Monte Carlo simulation is presented as a solution algorithm. Then, the proposed algorithms to solve imprecise structural reliability analyses are discussed in Section 5 and 6 for free and parametric probability-boxes, respectively. Finally, the performance of the proposed algorithms is assessed on a benchmark analytical function and realistic engineering problem settings, presented in Section 7.

3 Probability-boxes

3.1 Definition

A probability space is defined as the triplet $(\Omega, \mathcal{F}, \mathbb{P})$, where Ω is the universal event space, \mathcal{F} is the associated σ -algebra, and \mathbb{P} is the probability measure. A random variable X defines the mapping $X(\omega) : \omega \in \Omega \mapsto \mathcal{D}_X \subset \mathbb{R}$, where $\omega \in \Omega$ is an elementary event, \mathcal{D}_X is the support of X , and \mathbb{R} is the set of real numbers. In the context of probability theory, X is typically characterized by its cumulative distribution function (CDF) F_X and, in the case of continuous variables, by its probability density function (PDF) f_X . The CDF describes the probability of the event $\{X \leq x\}$, *i.e.* $F_X(x) = \mathbb{P}(X \leq x)$. Its PDF is defined as $f_X(x) = dF_X(x)/dx$ and describes the likelihood of X being in the neighbourhood of a given $x \in \mathcal{D}_X$.

Probability theory is based on the assumption that the probability measure \mathbb{P} is known precisely, resulting in a well-defined CDF. In cases of incomplete and/or sparse knowledge on X , however, aleatory uncertainty (natural variability) and epistemic uncertainty (lack of knowledge) can occur simultaneously. Epistemic uncertainty may not be considered adequately in the context of probability theory alongside aleatory uncertainty Beer et al. (2016). Hence, a more general formulation of the probability

measure, such as in Dempster-Shafer's theory of evidence and probability-boxes, is required to capture the uncertainty fully.

Dempster-Shafer's theory of evidence accounts for the epistemic uncertainty by replacing the probability measure \mathbb{P} with two values: belief and plausibility (Dempster, 1967; Shafer, 1976). *Belief* measures the minimum amount of probability that must be associated to an event, whereas *plausibility* measures the maximum amount of probability that could be associated to the same event. A *probability-box* (p-box) is a special case of the theory of evidence, considering only the nested set of events $\{X \leq x\}$, which in turns is related to the definition of a CDF of a probabilistic random variable.

P-boxes define the CDF of a variable X by lower and upper bounds denoted by \underline{F}_X and \overline{F}_X , respectively (Ferson and Ginzburg, 1996; Ferson and Hajagos, 2004). For any value $x \in \mathcal{D}_X$, the true-but-unknown CDF value lies within these bounds such that $\underline{F}_X(x) \leq F_X(x) \leq \overline{F}_X(x)$. The two boundary curves form an intermediate area, hence the name *probability-box*. In the literature, two types of p-boxes are distinguished, namely free and parametric p-boxes, which are discussed in the following.

3.2 Free p-box

Free p-boxes are defined as introduced in the previous section, *i.e.* only by lower and upper bounds of the CDF. This implies that the true CDF can have any arbitrary shape as long as it fulfils the characteristics of a generic CDF and lies within the bounds of the p-box. Figure 1(A) shows the boundary curves of a free p-box and a number of possible realizations of the true CDF. Because the shape of the true CDF is not specified, different types of curves are possible, including non-smooth ones (see realization #3 in Figure 1(A)).

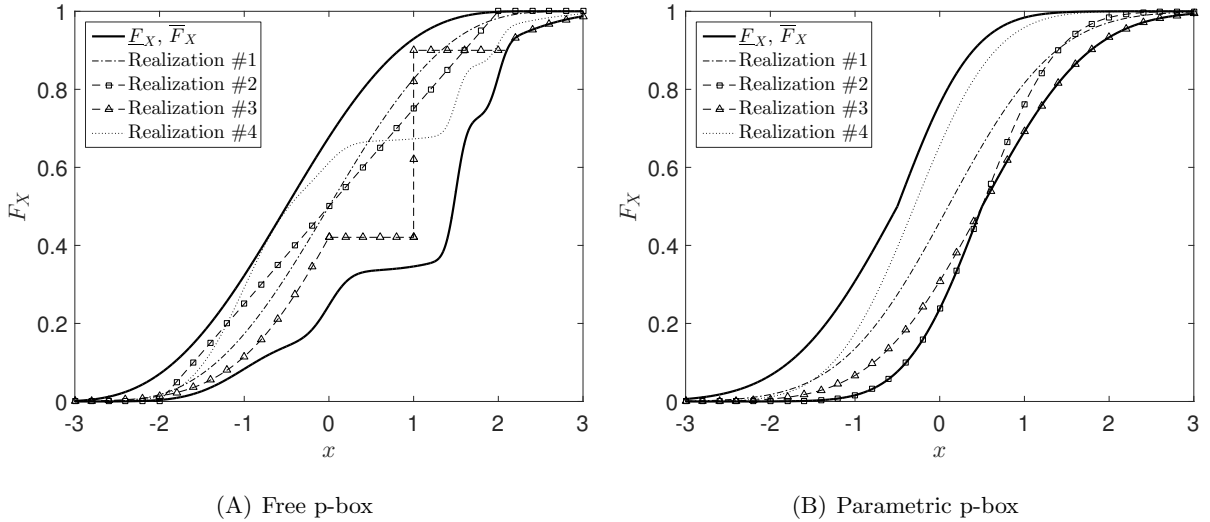


Figure 1: Definition of p-box types and realizations of possible CDFs

3.3 Parametric p-box

Parametric p-boxes (a.k.a. distributional p-boxes) are defined as cumulative distribution function families the parameters of which are known in intervals:

$$F_X(x) = F_X(x|\boldsymbol{\theta}), \quad \text{s.t.} \quad \boldsymbol{\theta} \in \mathcal{D}_{\boldsymbol{\theta}} \subset \mathbb{R}^{n_{\boldsymbol{\theta}}}, \quad (1)$$

where $\mathcal{D}_{\boldsymbol{\theta}}$ denotes the interval domain of the distribution parameters of dimension $n_{\boldsymbol{\theta}} = |\boldsymbol{\theta}|$. When the intervals are independent to each other, $\mathcal{D}_{\boldsymbol{\theta}} = [\underline{\theta}_1, \bar{\theta}_1] \times \dots \times [\underline{\theta}_{n_{\boldsymbol{\theta}}}, \bar{\theta}_{n_{\boldsymbol{\theta}}}]$ denotes a hyper-rectangular domain.

Parametric p-boxes allow for a clear separation of aleatory and epistemic uncertainty: aleatory uncertainty is represented by the distribution function family, whereas epistemic uncertainty is represented by the intervals in the distribution parameters. However, parametric p-boxes are more restrictive than free p-boxes because of the required knowledge on the distribution family. In other words, free p-boxes can be interpreted as a generalization of parametric p-boxes where the distribution family has $n_{\boldsymbol{\theta}} \rightarrow \infty$ parameters.

Parametric p-boxes are related to different concepts in the field of imprecise probabilities including Bayesian hierarchical models and fuzzy distributions. The parametric p-box construction resembles a Bayesian hierarchical model (Gelman, 2006) in which the distribution of the hyper-parameters $\boldsymbol{\theta}$ is replaced by an interval. From a different point of view, the parametric p-box is a fuzzy distribution function where the membership function of the distribution parameters is equal to one for $\boldsymbol{\theta} \in \mathcal{D}_{\boldsymbol{\theta}}$ and zero otherwise (Möller and Beer, 2004).

Figure 1(B) illustrates a parametric p-box, consisting of a Gaussian distribution family with mean value $\mu \in [-0.5, 0.5]$ and standard deviation $\sigma \in [0.7, 1.0]$. The lower and upper boundary curves of the parametric p-box drawn in Figure 1(B) are obtained by:

$$\underline{F}_X(x) = \min_{\boldsymbol{\theta} \in \mathcal{D}_{\boldsymbol{\theta}}} F_X(x|\boldsymbol{\theta}), \quad \overline{F}_X(x) = \max_{\boldsymbol{\theta} \in \mathcal{D}_{\boldsymbol{\theta}}} F_X(x|\boldsymbol{\theta}), \quad (2)$$

where $\boldsymbol{\theta} = (\mu, \sigma)^T$. Hence, the boundary curves shown in Figure 1(B) have the characteristics of a CDF but are not necessarily a realization with a specific parameter combination $\boldsymbol{\theta}_0$. They can consist of sections of different realizations. As an example, the lower boundary CDF is a combination of realization #2 ($\mu = 0.5, \sigma = 0.7$) and realization #3 ($\mu = 0.5, \sigma = 1.0$).

4 Structural reliability analysis

4.1 Limit-state function

A limit-state function describes the performance of a process or system as a function of a set of input parameters. The deterministic mapping is defined as:

$$\mathcal{G} : \mathbf{x} \in \mathcal{D}_{\mathbf{x}} \subset \mathbb{R}^M \mapsto y = \mathcal{G}(\mathbf{x}) \in \mathbb{R}, \quad (3)$$

where \mathbf{x} is a M -dimensional vector defined in the input domain $\mathcal{D}_{\mathbf{X}}$ and y is the output scalar indicating the performance. The sign of y , and hence $\mathcal{G}(\mathbf{x})$, determines whether an input vector \mathbf{x} corresponds to a safe system ($\mathcal{G}(\mathbf{x}) > 0$) or a failed system ($\mathcal{G}(\mathbf{x}) \leq 0$). The limit-state function \mathcal{G} is interpreted as a *black box* of which only the input vector \mathbf{x} and the corresponding response $y = \mathcal{G}(\mathbf{x})$ are accessible.

4.2 Failure probability

Due to uncertainties in the definition of the input vector, \mathbf{X} is modelled traditionally by probability distributions and in this paper \mathbf{X} is modelled by p-boxes. Considering a probabilistic input vector \mathbf{X} , the *failure probability* is defined as the probability that the limit-state function takes negative values:

$$P_f = \mathbb{P}(Y \leq 0) = \mathbb{P}(\mathcal{G}(\mathbf{X}) \leq 0), \quad (4)$$

which can be recast as an integral:

$$P_f = \int_{\mathcal{D}_f} f_{\mathbf{X}}(\mathbf{x}) d\mathbf{x}, \quad (5)$$

where $\mathcal{D}_f = \{\mathbf{x} \in \mathcal{D}_{\mathbf{X}} : \mathcal{G}(\mathbf{x}) \leq 0\}$ is the failure domain and $f_{\mathbf{X}}$ is the joint probability density function of the input vector \mathbf{X} . The integration in Eq. (5) cannot be performed analytically in the general case where the failure domain \mathcal{D}_f has a complex shape. Hence, numerical estimates for the failure probability were developed such as Monte Carlo simulation. Considering a large sample of \mathbf{X} denoted by $\mathcal{S} = \{\mathbf{x}_1, \dots, \mathbf{x}_n\}$, the failure probability can be estimated by:

$$\hat{P}_f = \frac{n_f}{n} = \frac{1}{n} \sum_{i=1}^n \mathbb{I}_{\mathcal{G}(\mathbf{x}) \leq 0}(\mathbf{x}_i), \quad (6)$$

where n_f is the number of failure samples $\mathbf{x}_i \in \mathcal{D}_f$, $n = |\mathcal{S}|$ is the total number of samples and \mathbb{I} is the indicator function with $\mathbb{I} = 1$ for a true subscript statement and $\mathbb{I} = 0$ otherwise. Note that the use of \mathbb{I} transforms the structural reliability problem into a classification problem where only the sign of $\mathcal{G}(\mathbf{x})$ is relevant.

In the context of p-boxes, the joint probability density function $f_{\mathbf{X}}$ is not defined deterministically. Hence, Eq. (5) leads to a range of failure probabilities $P_f \in [\underline{P}_f, \overline{P}_f]$, the bounds of which are defined as:

$$\underline{P}_f = \underset{f_{\mathbf{X}}}{\text{"min"}} \int_{\mathcal{D}_f} f_{\mathbf{X}}(\mathbf{x}) d\mathbf{x}, \quad \overline{P}_f = \underset{f_{\mathbf{X}}}{\text{"max"}} \int_{\mathcal{D}_f} f_{\mathbf{X}}(\mathbf{x}) d\mathbf{x}, \quad (7)$$

where "min" (resp. "max") means that the optimization would be carried out over all PDF $f_{\mathbf{X}}$ that satisfy some constraints related to the definition of the underlying p-box. This *imprecise structural reliability analysis* (ISRA) is not straightforward as it involves an optimization with a multi-dimensional distribution $f_{\mathbf{X}}$ as the argument of the objective function. In the following sections, solution algorithms for the two types of p-boxes, *i.e.* free and parametric p-boxes, are discussed.

5 Imprecise structural reliability analysis (ISRA) for free p-boxes

5.1 Problem conversion

For the case of free p-boxes, the imprecise structural reliability problem can be recast as two structural reliability problems in the probabilistic sense (Zhang et al., 2010; Zhang, 2012; Schöbi and Sudret, 2015a). These can be solved by conventional structural reliability methods, such as Monte Carlo simulations.

Consider a random vector \mathbf{C} of length M which follows a uniform distribution in the unit-hypercube domain $[0, 1]^M$. The component C_i shall describe the CDF value of X_i . For the sake of simplicity, the input variables X_i are assumed to be statistically independent throughout this paper. When X_i is modelled by a free p-box, each $c_i \in \mathcal{D}_{C_i}$ can be transformed into an interval in the domain \mathcal{D}_{X_i} by applying the inverse CDF of the p-box bounds:

$$\underline{x}_i = \overline{F}_{X_i}^{-1}(c_i), \quad \overline{x}_i = \underline{F}_{X_i}^{-1}(c_i). \quad (8)$$

For a given realization of $\mathbf{c} \in \mathcal{D}_{\mathbf{C}}$, let us denote by $\mathcal{D}_{\mathbf{c}} = [\underline{x}_1(c_1), \overline{x}_1(c_1)] \times \dots \times [\underline{x}_M(c_M), \overline{x}_M(c_M)]$. The boundary curves of the p-box of the response variable Y are then obtained by optimization:

$$\overline{Y} = \underline{\mathcal{G}}(\mathbf{C}) = \min_{\mathbf{x} \in \mathcal{D}_{\mathbf{c}}} \mathcal{G}(\mathbf{x}), \quad \underline{Y} = \overline{\mathcal{G}}(\mathbf{C}) = \max_{\mathbf{x} \in \mathcal{D}_{\mathbf{c}}} \mathcal{G}(\mathbf{x}). \quad (9)$$

Note that due to the definition of $\overline{\mathcal{G}}$, taking the maximum response value naturally leads to a realization of the lower bound CDF of the response \underline{Y} and vice versa. Based on those boundary curves and the definition of the failure domain \mathcal{D}_f , the range of failure probabilities is obtained by:

$$\underline{P}_f = \mathbb{P}(\underline{Y} \leq 0) = \mathbb{P}(\overline{\mathcal{G}}(\mathbf{C}) \leq 0), \quad \overline{P}_f = \mathbb{P}(\overline{Y} \leq 0) = \mathbb{P}(\underline{\mathcal{G}}(\mathbf{C}) \leq 0). \quad (10)$$

Then, the bounds of the failure probability can be estimated by sampling the probabilistic input vector \mathbf{C} . Note that the two boundary values can be computed independently.

A drawback of this problem conversion and the subsequent Monte Carlo simulation, however, is the high computational costs due to three distinct causes. Firstly, the computational model can be an expensive-to-evaluate model in many applications, such as finite element models. Secondly, the number of model evaluations may be large due to the optimization operations in Eq. (9), for each $\mathbf{c} \in \mathcal{D}_{\mathbf{C}}$. And thirdly, failure probabilities are generally low, typically $P_f \approx 10^{-3} \dots 10^{-6}$ in engineering applications, which requires a large number of samples in Monte Carlo simulation to accurately estimate the failure probability, typically $100/P_f$ for a 10 % accuracy in the estimate. In order to cope with the three main aspects contributing to the total computational effort, meta-models are introduced at different levels of the ISRA for free p-boxes.

5.2 Adaptive Kriging Monte Carlo simulation

5.2.1 Kriging

Consider again the limit-state function in Eq. (3) and a probabilistic input vector \mathbf{X} . Kriging is a meta-modelling approach that considers the limit-state function to be a realization of a Gaussian process (Santner et al., 2003):

$$Y^{(K)}(\mathbf{x}) = \boldsymbol{\beta}^\top \mathbf{f}(\mathbf{x}) + \sigma^2 Z(\mathbf{x}, \omega), \quad (11)$$

where $\boldsymbol{\beta}^\top \mathbf{f}(\mathbf{x})$ is the mean value of the Gaussian process consisting of regression functions $\mathbf{f}(\mathbf{x}) = \{f_j(\mathbf{x}), j = 1, \dots, p\}$ and a vector of coefficients $\boldsymbol{\beta} \in \mathbb{R}^p$, σ^2 is the variance of the Gaussian process, and $Z(\mathbf{x}, \omega)$ is a zero-mean, unit-variance stationary Gaussian process, characterized by an autocorrelation function $R(|\mathbf{x} - \mathbf{x}'|; \boldsymbol{\rho})$ and its hyper-parameter(s) $\boldsymbol{\rho}$.

Training a Kriging meta-model is based on a set of input realizations $\mathcal{X} = \{\boldsymbol{\chi}^{(j)}, i = 1, \dots, N\}$ and the corresponding values of the limit-state function $\mathcal{Y} = \{\mathcal{Y}^{(j)} = \mathcal{G}(\boldsymbol{\chi}^{(j)}), j = 1, \dots, N\}$. Then, the Kriging parameters are obtained by the generalized least-squares solution (Santner et al., 2003):

$$\boldsymbol{\beta}(\boldsymbol{\rho}) = \left(\mathbf{F}^\top \mathbf{R}^{-1} \mathbf{F} \right)^{-1} \mathbf{F}^\top \mathbf{R}^{-1} \mathcal{Y}, \quad (12)$$

$$\sigma_y^2(\boldsymbol{\rho}) = \frac{1}{n} (\mathcal{Y} - \mathbf{F}\boldsymbol{\beta})^\top \mathbf{R}^{-1} (\mathcal{Y} - \mathbf{F}\boldsymbol{\beta}), \quad (13)$$

where $\mathbf{R}_{ij} = R(|\boldsymbol{\chi}^{(i)} - \boldsymbol{\chi}^{(j)}|; \boldsymbol{\rho})$ is the correlation matrix and $\mathbf{F}_{ij} = f_j(\boldsymbol{\chi}^{(i)})$ is the information matrix. When the correlation parameters $\boldsymbol{\rho}$ are unknown, their values can be estimated by maximum likelihood or cross validation (Bachoc, 2013).

Given an arbitrary point \mathbf{x} of the input domain $\mathcal{D}_{\mathbf{X}}$, the prediction value of the limit-state function is a Gaussian variable with mean value and variance:

$$\mu_{\hat{\mathcal{Y}}}(\mathbf{x}) = \mathbf{f}^\top(\mathbf{x})\boldsymbol{\beta} + \mathbf{r}^\top(\mathbf{x})\mathbf{R}^{-1}(\mathcal{Y} - \mathbf{F}\boldsymbol{\beta}), \quad (14)$$

$$\sigma_{\hat{\mathcal{Y}}}^2(\mathbf{x}) = \sigma_y^2 \left(1 - \mathbf{r}^\top(\mathbf{x})\mathbf{R}\mathbf{r}(\mathbf{x}) + \mathbf{u}^\top(\mathbf{x}) \left(\mathbf{F}^\top \mathbf{R}^{-1} \mathbf{F} \right)^{-1} \mathbf{u}(\mathbf{x}) \right), \quad (15)$$

where $r_i(\mathbf{x}) = R(|\mathbf{x} - \boldsymbol{\chi}^{(i)}|; \boldsymbol{\rho})$ and $\mathbf{u}(\mathbf{x}) = \mathbf{F}^\top \mathbf{R}^{-1} \mathbf{r}(\mathbf{x}) - \mathbf{f}(\mathbf{x})$. The Kriging predictor interpolates the experimental design, meaning that $\mu_{\hat{\mathcal{Y}}}(\boldsymbol{\chi}^{(i)}) = \mathcal{G}(\boldsymbol{\chi}^{(i)})$ and $\sigma_{\hat{\mathcal{Y}}}(\boldsymbol{\chi}^{(i)}) = 0$ for $\boldsymbol{\chi}^{(i)} \in \mathcal{X}$.

5.2.2 Adaptive experimental designs

Kriging is an interpolation algorithm which approximates the limit-state function most accurately close to the points of the experimental design \mathcal{X} . However, these points are not necessarily suitable to approximate the boundary between the safety and failure domain, *i.e.* the *limit-state surface* $\mathcal{G}(\mathbf{x}) = 0$, and hence to estimate the failure probability. Then, an active learning algorithm can improve the estimation of P_f by enriching the experimental design in a guided way.

Adaptive Kriging Monte Carlo simulation (AK-MCS) combines Kriging with an enrichment scheme and Monte Carlo simulation, as proposed in Echard et al. (2011) and further discussed in Schöbi et al. (2016); Marelli et al. (2015). The main steps of AK-MCS are described here:

1. Generate a small initial experimental design \mathcal{X} and compute the corresponding response of the limit-state function: $\mathcal{Y}^{(i)} = \mathcal{G}(\mathcal{X}^{(i)})$.
2. Train a Kriging meta-model $\mathcal{G}^{(K)}$ based on $\{\mathcal{X}, \mathcal{Y}\}$. In this paper, ordinary Kriging meta-models are used based on a constant (yet unknown) trend β .
3. Generate a large set of candidate samples $\mathcal{S} = \{\mathbf{x}_1, \dots, \mathbf{x}_n\}$ from \mathbf{X} and evaluate the response of the meta-model: $\mu_{\hat{Y}}(\mathbf{x})$ and $\sigma_{\hat{Y}}^2(\mathbf{x})$.
4. Compute the failure probability and its confidence values:

$$\hat{P}_f = \mathbb{P}(\mu_{\hat{Y}}(\mathbf{x}) \leq 0), \quad (16)$$

$$\hat{P}_f^+ = \mathbb{P}(\mu_{\hat{Y}}(\mathbf{x}) - k\sigma_{\hat{Y}}(\mathbf{x}) \leq 0), \quad \hat{P}_f^- = \mathbb{P}(\mu_{\hat{Y}}(\mathbf{x}) + k\sigma_{\hat{Y}}(\mathbf{x}) \leq 0). \quad (17)$$

In practice, $k = 2$ is selected which approximately corresponds to a point-wise 95 % confidence interval on $\hat{Y}(\mathbf{x})$.

5. Check the convergence criterion for the estimate of the failure probability:

$$\frac{\hat{P}_f^+ - \hat{P}_f^-}{\hat{P}_f} \leq \epsilon_{P_f}. \quad (18)$$

It has been shown that $\epsilon_{P_f} = 5\%$ gives accurate results at reasonable costs (Schöbi et al., 2016). If the criterion is fulfilled, terminate the adaptive algorithm and return the last meta-model $\mathcal{G}^{(K)}$. Otherwise, continue with the next step.

6. Compute the probability of misclassification for every $\mathbf{x}_i \in \mathcal{S}$ Bect et al. (2012):

$$P_m(\mathbf{x}_i) = \Phi\left(-\frac{|\mu_{\hat{Y}}(\mathbf{x}_i)|}{\sigma_{\hat{Y}}(\mathbf{x})}\right), \quad (19)$$

where $\Phi(\cdot)$ is the CDF of the standard normal variable. Select the best next sample to be added to the experimental design, which maximizes the probability of misclassification:

$$\mathbf{x}^* = \arg \max_{i=1, \dots, n} P_m(\mathbf{x}_i). \quad (20)$$

7. Add \mathbf{x}^* to the experimental design and compute $\mathcal{Y}^* = \mathcal{G}(\mathbf{x}^*)$. Then, go back to step 2 to update the meta-model with the enriched experimental design.

After the termination of the adaptive Kriging iterations, the failure probability is estimated based on the last meta-model $\mathcal{G}^{(K)}$ and the Monte Carlo sample \mathcal{S} .

5.3 Meta-modelling the limit-state surface $\mathcal{G}(\mathbf{x}) = 0$

A first meta-model is applied to the limit-state function \mathcal{G} and in particular to model the limit-state surface $\{\mathbf{x} \in \mathcal{D}_{\mathbf{X}} : \mathcal{G}(\mathbf{x}) = 0\}$. However, in order to conduct an AK-MCS analysis, a probabilistic input vector is required. When the input is modelled by free p-boxes in ISRA, an auxiliary input vector $\widetilde{\mathbf{X}}$

is created by *condensation* (Schöbi and Sudret, 2016). In this paper, the auxiliary input variables \tilde{X}_i are characterized by the average value of the boundary curves of the p-box:

$$F_{\tilde{X}_i}(x_i) = \frac{1}{2} (F_{X_i}(x_i) + \bar{F}_{X_i}(x_i)). \quad (21)$$

The auxiliary distribution covers the shape of the p-box. Hence, the resulting meta-model $\mathcal{G}^{(K)}$ is accurate in the very neighbourhood of the limit-state surface that contributes to the value of the failure probability in the p-box setting.

5.4 Meta-modelling the limit-state surfaces $\underline{\mathcal{G}}(\mathbf{x}) = 0$ and $\bar{\mathcal{G}}(\mathbf{x}) = 0$

The second meta-model is applied to the limit-state functions $\underline{\mathcal{G}}$ and $\bar{\mathcal{G}}$ and the estimation of the bounds of the failure probability (see Eqs. (9) and (10)). By using the approximation of \mathcal{G} , Eq. (9) may be replaced by:

$$\underline{\mathcal{G}}(\mathbf{c}) \approx \min_{\mathbf{x} \in \mathcal{D}_c} \mathcal{G}^{(K)}(\mathbf{x}), \quad \bar{\mathcal{G}}(\mathbf{c}) \approx \max_{\mathbf{x} \in \mathcal{D}_c} \mathcal{G}^{(K)}(\mathbf{x}), \quad (22)$$

where $\mathcal{G}^{(K)}$ is the meta-model resulting from the first-level approximation. Therefore, the bounds of the failure probability \underline{P}_f and \bar{P}_f can be estimated by two independent AK-MCS analyses of $\bar{\mathcal{G}}$ and $\underline{\mathcal{G}}$, respectively, *i.e.* $\underline{P}_f = \mathbb{P}(\bar{\mathcal{G}}(\mathbf{C}) \leq 0)$ and $\bar{P}_f = \mathbb{P}(\underline{\mathcal{G}}(\mathbf{C}) \leq 0)$ (see also Eq. (10)). Note that here the input vector consists of the probabilistic input vector \mathbf{C} , which makes the definition of an auxiliary distribution as in Section 5.3 unnecessary.

To improve the convergence of the AK-MCS algorithms, however, an auxiliary distribution may be beneficial. In fact, the auxiliary distribution $\tilde{\mathbf{X}}$ defined in Eq. (21) is suitable when p-boxes are unbounded, as shown in Schöbi and Sudret (2015a). Then, an isoprobabilistic transform T maps $\tilde{\mathbf{X}}$ to \mathbf{C} and vice versa. The failure probability is estimated by $\underline{P}_f = \mathbb{P}(\bar{\mathcal{G}}(T(\tilde{\mathbf{X}})) \leq 0)$ and $\bar{P}_f = \mathbb{P}(\underline{\mathcal{G}}(T(\tilde{\mathbf{X}})) \leq 0)$, respectively.

5.5 Two-level meta-modelling approach

Figure 2 summarizes the procedure, which consists of two sequential meta-modelling levels connected by a model conversion step. The first-level meta-model surrogates the original limit-state function \mathcal{G} , whereas the two second-level ones surrogate the limit-state functions for estimating the lower and upper bounds of the failure probability. Both levels use auxiliary random variables, *i.e.* $\tilde{\mathbf{X}}$ on the first level and \mathbf{C} on the second level.

6 Imprecise structural reliability analysis (ISRA) for parametric p-boxes

6.1 Nested algorithm

The definition of parametric p-boxes indicates a hierarchical model where the distribution of \mathbf{X} is defined conditionally on its distribution parameters. Hence, nested simulation algorithms can be applied

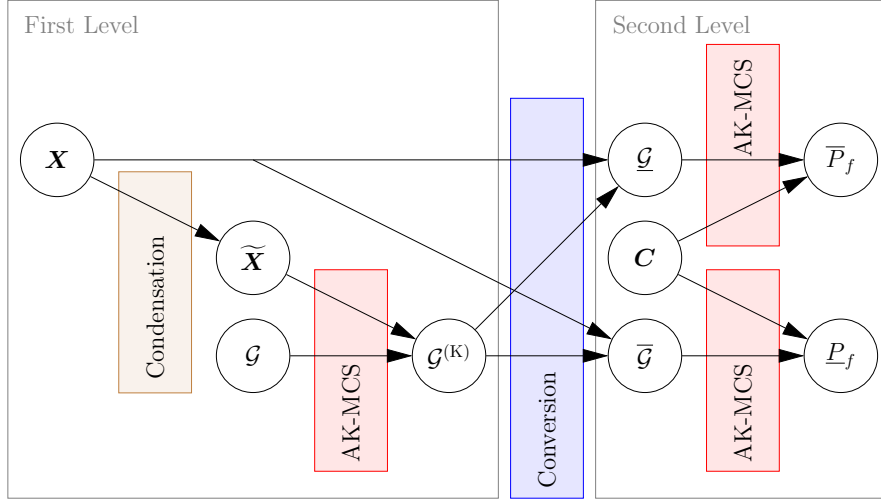


Figure 2: Imprecise structural reliability analysis for free p-boxes – flowchart of the two-level meta-modelling algorithm

in the context of ISRA, as discussed in Eldred and Swiler (2009); Schöbi and Sudret (2015b). In other words, the bounds of the failure probability can be found by estimating the conditional failure probabilities $P_{f|\theta} = \mathbb{P}(\mathcal{G}(\mathbf{X}_\theta) \leq 0)$, where \mathbf{X}_θ is a conditional distribution with $F_{\mathbf{X}_\theta}(\mathbf{x}) = F_{\mathbf{X}}(\mathbf{x}|\theta)$, and minimizing/maximizing it with respect to θ to get the bounds $\underline{P}_f = \min_{\theta \in \mathcal{D}_\Theta} P_{f|\theta}$ and $\bar{P}_f = \max_{\theta \in \mathcal{D}_\Theta} P_{f|\theta}$.

Making use of the same tools as in Section 5 for free p-boxes, the failure probability can be found with the help of Kriging meta-models. The conditional failure probabilities $P_{f|\theta}$ are estimated by AK-MCS, whereas the efficient global optimization (EGO) Jones et al. (1998) algorithm is used to optimize on θ . Figure 3 illustrates the main components of the proposed algorithm. In the following two sections, AK-MCS and EGO are discussed more in depth.

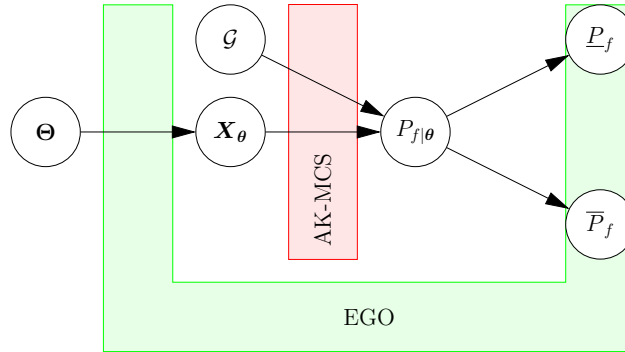


Figure 3: Imprecise structural reliability analysis for parametric p-boxes – overview of the nested algorithm

6.2 AK-MCS for conditional failure probabilities

In the context of parametric p-boxes, the core task consists of estimating efficiently the conditional failure probability $P_{f|\theta}$ by AK-MCS. Given a vector θ and consequently the conditional variable \mathbf{X}_θ , AK-MCS can be applied to estimate the conditional failure probability as described in Section 5.2.2.

6.3 Adaptive-Kriging for the lower bound \underline{P}_f

EGO is a global optimization algorithm which is based on Kriging meta-models with a design enrichment strategy similar to AK-MCS (Echard et al., 2011; Jones et al., 1998). In the context of ISRA for parametric p-boxes, EGO is used to find the extreme values of the failure probability by optimizing on the distribution parameters θ . The main steps for finding the minimum failure probability are discussed here:

1. Given the distribution parameter domain $\theta \in \mathcal{D}_\theta$, generate a small initial experimental design $\mathcal{T} = \{\tau^{(1)}, \dots, \tau^{(N)}\}$.
2. Compute the corresponding failure probabilities $\mathcal{P}^{(i)} = \mathbb{P}(\mathcal{G}(\mathbf{X}_{\tau^{(i)}}) \leq 0)$, $i = 1, \dots, N$ by AK-MCS. Note that for the first evaluation (*i.e.* $\theta = \tau^{(1)}$) of such a conditional failure probability, the procedure of Section 5.2.2 is applicable one-to-one with the distribution $\mathbf{X}_{\tau^{(1)}}$. However, for $i = 2, \dots, n$, the limit-state function evaluations of the previous AK-MCS analyses (*i.e.* the samples of the final experimental design \mathcal{X}) are kept as initial experimental design. The difference between two samples $\tau^{(i)}$ and $\tau^{(j)}$ lies then solely in a modified $\mathbf{X}_{\tau^{(i)}}$ and hence a modified $\mathcal{S}_{\tau^{(i)}}$. Due to the similarity of the MC populations $\mathcal{S}_{\tau^{(i)}}$, the number of limit-state function evaluations can be kept small (see also the application in Section 7.2).
3. Train a Kriging meta-model $\mathcal{M}^{(K)}$ based on $\{\mathcal{T}, \mathcal{P}\}$. This meta-model approximates the estimation of the conditional failure probability $P_{f|\theta}(\theta)$.
4. Search for the optimal samples to be added to the experimental design for minimizing the failure probability:

$$\tau_{min}^* = \arg \max_{\theta \in \mathcal{D}_\theta} [\text{EI}_{min}(\theta)], \quad (23)$$

where the expected improvement is defined as (Jones et al., 1998; Mockus et al., 1978):

$$\text{EI}_{min}(\theta) = \left(\mathcal{P}_{min} - \mu_{\hat{P}_f}(\theta) \right) \Phi \left(\frac{\mathcal{P}_{min} - \mu_{\hat{P}_f}(\theta)}{\sigma_{\hat{P}_f}(\theta)} \right) + \sigma_{\hat{P}_f}(\theta) \varphi \left(\frac{\mathcal{P}_{min} - \mu_{\hat{P}_f}(\theta)}{\sigma_{\hat{P}_f}(\theta)} \right), \quad (24)$$

where EI_{min} is the expected improvement for minimizing P_f , $\mathcal{P}_{min} = \min_i [\mathcal{P}^{(i)}]$ is the current minimum failure probability, $\Phi(\cdot)$ and $\varphi(\cdot)$ are the CDF and PDF of a standard normal variable, respectively. As explained in Jones et al. (1998), the expected improvement allows for a balance between local and global search.

5. Check the convergence criterion proposed by Jones et al. (1998), applied to the minimum estimate of the failure probability:

$$\text{EI}_{min}(\tau_{min}^*) \leq \epsilon_{EI}. \quad (25)$$

If the criterion is fulfilled, terminate the optimization algorithm and return the last estimate of the minimum failure probability: $\underline{P}_f = \min_i [\mathcal{P}^{(i)}]$. Otherwise, continue with Step 6. Note that a value of $\epsilon_{EI} = 10^{-5}$ was identified as a reliable choice for the applications in Section 7.

6. Add the best next sample $\boldsymbol{\tau}_{min}^*$ to the experimental design \mathcal{T} according to Eq. (23).
7. Evaluate the failure probability corresponding to $\boldsymbol{\tau}_{min}^*$ and add it to \mathcal{P} . Note that the limit-state function evaluations computed in previous iterations in AK-MCS can be recycled. They can be used as the initial experimental design in AK-MCS for estimating \mathcal{P}^* corresponding to $\boldsymbol{\tau}^*$ (see also Step 2 for further details). Finally, continue with Step 3.

6.4 Adaptive Kriging for the upper bound \overline{P}_f

The upper boundary value of the failure probability \overline{P}_f can be estimated by an EGO algorithm analogously to the lower boundary value as follows. The best next sample is determined by replacing Eq. (23) by:

$$\boldsymbol{\tau}_{max}^* = \arg \max_{\boldsymbol{\theta} \in \mathcal{D}_{\boldsymbol{\Theta}}} [\text{EI}_{max}(\boldsymbol{\theta})], \quad (26)$$

where the expected improvement is defined by:

$$\text{EI}_{max}(\boldsymbol{\theta}) = \left(\mu_{\hat{P}_f}(\boldsymbol{\theta}) - \mathcal{P}_{max} \right) \Phi \left(\frac{\mu_{\hat{P}_f}(\boldsymbol{\theta}) - \mathcal{P}_{max}}{\sigma_{\hat{P}_f}(\boldsymbol{\theta})} \right) + \sigma_{\hat{P}_f}(\boldsymbol{\theta}) \varphi \left(\frac{\mu_{\hat{P}_f}(\boldsymbol{\theta}) - \mathcal{P}_{max}}{\sigma_{\hat{P}_f}(\boldsymbol{\theta})} \right), \quad (27)$$

where $\mathcal{P}_{max} = \max_i [\mathcal{P}^{(i)}]$ is the current maximum failure probability. Furthermore, the algorithm is terminated with a stopping criterion similar to Eq. (25):

$$\text{EI}_{max}(\boldsymbol{\tau}_{max}^*) \leq \epsilon_{EI}. \quad (28)$$

When this is not fulfilled, the best next sample $\boldsymbol{\tau}_{max}^*$ is added to the experimental design \mathcal{T} . After termination of the iterative algorithm, the maximum failure probability is estimated by $\overline{P}_f = \max_i [\mathcal{P}^{(i)}]$.

6.5 Joint optimization

The optimization for the minimum and maximum failure probabilities are discussed separately in Sections 6.3 and 6.4, respectively. However, the Steps 4 and 5 of the EGO algorithm allow for a *simultaneous* optimization for the bounds of the failure probability. For simultaneous optimization, both boundary values of the failure probability are estimated with the same meta-model, $\{\boldsymbol{\tau}_{min}^*, \boldsymbol{\tau}_{max}^*\}$ are determined by Eq. (23) and (26), respectively, and are added in each iteration of the EGO algorithm. In fact, depending on the two stopping criteria, $\{\boldsymbol{\tau}_{min}^*, \boldsymbol{\tau}_{max}^*\}$ is added if Eqs. (25) and (28) are both *not* fulfilled, $\boldsymbol{\tau}_{min}^*$ is added if only Eq. (28) is fulfilled, or $\boldsymbol{\tau}_{max}^*$ is added if only Eq. (25) is fulfilled. If both equations are fulfilled, the iterative algorithm is terminated.

For *separate* optimization, the lower and upper bounds of the failure probability are estimated one after the other as presented in the Section 6.3 and 6.4. The effect of the different optimization modes is discussed in Section 7.1.

6.6 Comparison to free p-boxes

ISRA for parametric p-boxes consists of two interconnected Kriging models, the experimental designs of which are enriched quasi simultaneously, as shown in Figure 3. In the context of free p-boxes, two levels of meta-models are computed subsequently (see Figure 2). Additionally, an optimization operation is required for free p-boxes, which increases the complexity and potentially the computational effort in this case. The increased complexity originates from the basic definition of the free p-box, which can be interpreted as a generalization of the parametric p-box, as discussed in Section 3.3.

Assuming, however, that the computational costs are dominated by the evaluation of the limit-state function, the total computational costs for estimating the boundary values of the failure probabilities may be similar for free and parametric p-boxes. The total number of limit-state function evaluations may be similar due to the same method (*i.e.* AK-MCS) used to surrogate the limit-state surface $\mathcal{G}(\mathbf{x}) = 0$.

7 Applications

7.1 Two-dimensional toy function

7.1.1 Problem statement

The first example is a simple two-dimensional problem used to illustrate the machinery of the proposed approaches. The limit-state function is defined as:

$$g_1(\mathbf{x}) = x_1 - x_2^2. \quad (29)$$

Failure is defined as $g_1(\mathbf{x}) \leq 0$ and hence the failure probability is $P_f = \mathbb{P}(g_1(\mathbf{x}) \leq 0)$. The two variables X_i are modelled by p-boxes. In the following, two cases are distinguished: (i) X_i is modelled by free p-boxes and (ii) X_i is modelled by parametric p-boxes. In the former case, the bounds of the free p-box are defined as:

$$\underline{F}_{X_i}(x_i) = F_{\mathcal{N}}(x_i | \mu = 2.5, \sigma = 1), \quad \overline{F}_{X_i}(x_i) = F_{\mathcal{N}}(x_i | \mu = 1.5, \sigma = 1), \quad i = 1, 2, \quad (30)$$

where $F_{\mathcal{N}}(x | \mu, \sigma)$ denotes the CDF of a Gaussian distribution with mean value μ and standard deviation σ . For the parametric p-box case, the distribution parameters θ are chosen so that the boundary curves are the same as in the case of free p-boxes:

$$F_{X_i}(x_i) = F_{\mathcal{N}}(x_i | \mu, \sigma = 1), \quad \mu \in [1.5, 2.5]. \quad (31)$$

7.1.2 Analysis

The settings for the ISRA are the following. The meta-models and Monte Carlo simulations are computed using the Matlab-based uncertainty quantification framework UQLAB (Marelli and Sudret, 2014), including its implementation of AK-MCS (Marelli et al., 2015). All Kriging models are defined

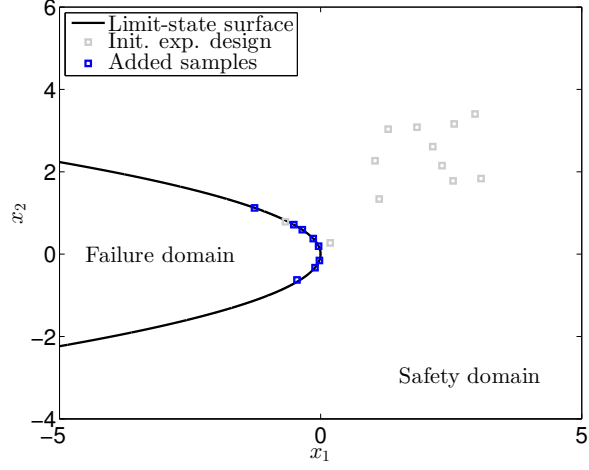
by a Gaussian autocorrelation function and a constant trend ($f(\mathbf{x}) = 1$, a.k.a. ordinary Kriging). The number of samples in the initial experimental design are $N_1^0 = 12$ and $N_2^0 = 4$ for AK-MCS and EGO, respectively. The initial experimental designs are generated by the Latin-hypercube sampling method (McKay et al., 1979) in the unit hypercube $[0, 1]^M$ and subsequent isoprobabilistic transform to the corresponding input domains. For free p-boxes, the auxiliary input distribution is defined as $\tilde{X}_i \sim \mathcal{N}(x_i|2, 1)$. For parametric p-boxes and EGO, the threshold value of the stopping criterion is set to $\epsilon_{EI} = 10^{-5}$. The failure probabilities are estimated with a set of $n = 10^6$ Monte Carlo samples. The analysis is repeated 50 times with different initial experimental designs in order to obtain statistically significant results.

7.1.3 Results

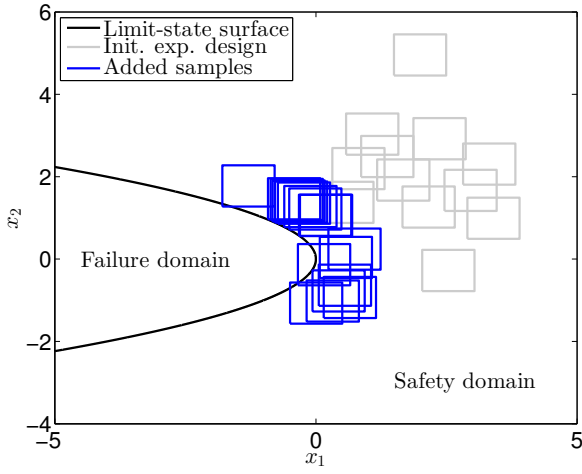
Free p-boxes Figure 4 visualizes a typical realization of the ISRA for free p-boxes. Note that the stopping criterion in AK-MCS was deactivated in this simulation run to visualize the general behaviour of the method. Figure 4(A) illustrates the experimental design of the final first-level meta-model. The initial experimental design is marked by the grey squares. The additional samples (blue squares) group around the limit-state surface (black solid line) in the physical domain, which implies an efficient selection of additional samples. Further, Figures 4(B) and 4(C) show the constraint optimization domains \mathcal{D}_c in Eq. (22) in the physical space $\mathcal{D}_{\mathbf{X}}$ for the experimental design \mathfrak{C} on the second-level meta-model. Again, the optimization domains are selected in an efficient manner, as seen from the positioning of the blue rectangles. For $\underline{\mathcal{G}}$ in Figure 4(B), the minimum value of the limit-state function within each blue rectangular is achieved at a point which lies close to the limit-state surface. Thus, the rectangles are mostly contained in the safety domain. Analogously in Figure 4(C) for $\overline{\mathcal{G}}$, the maximum value of the limit-state function within each blue rectangle is close to the limit-state surface. Hence, the rectangles are mostly contained in the failure domain.

Accounting for the 50 replications of the ISRA, the estimates of the failure probability are summarized in Table 1. The reference value of the failure probability (denoted by $P_{f,ref}$) is obtained by an Importance Sampling analysis of $\underline{\mathcal{G}}$ and $\overline{\mathcal{G}}$ with a sample size of $n = 10^6$. The table shows that all three meta-models accurately estimate the corresponding failure probabilities in terms of mean estimate. However, in terms of coefficient of variation, the second-level meta-models are less accurate than the expected Monte Carlo sampling-related values: for \underline{P}_f the expected value is $\text{CoV}[\underline{P}_f] = \sqrt{(1 - \underline{P}_f)/(n_{MC} \cdot \underline{P}_f)} = 11.3\% < 13.8\%$ and for \overline{P}_f the expected value is $\text{CoV}[\overline{P}_f] = 0.9\% < 7.4\%$. The large variation in \overline{P}_f is caused by the larger number of samples in the failure domain and hence the larger length of the limit-state surface relevant for the failure probability estimation, when comparing to the auxiliary space $\mathcal{D}_{\tilde{\mathbf{X}}}$. Selecting a different auxiliary distribution $\tilde{\mathbf{X}}$ may possibly reduce this phenomenon.

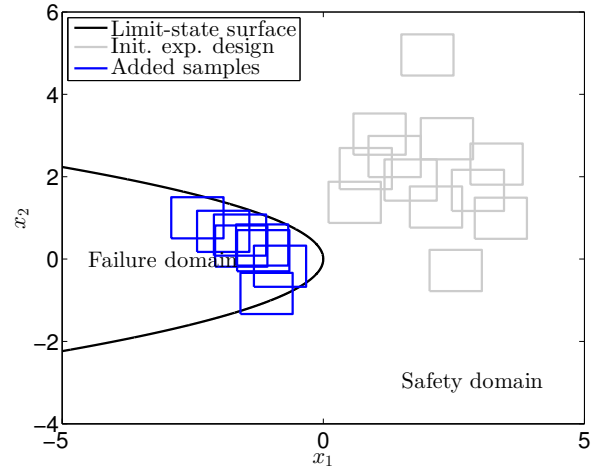
The lower part of Table 1 shows the number of samples in the experimental design of the final meta-models. N_1 and N_2 denote the number of model evaluations on the first-level and second-level meta-model, respectively (average over 50 replications of the full algorithm). N_1 is low due to the



(A) Experimental design for \mathcal{G} (first-level meta-model)



(B) Optimization domains in $\mathcal{D}_{\mathbf{X}}$ for $\underline{\mathcal{G}}$ (second-level meta-model)



(C) Optimization domains in $\mathcal{D}_{\mathbf{X}}$ for $\overline{\mathcal{G}}$ (second-level meta-model)

Figure 4: Two-dimensional toy function – free p-boxes – adaptive experimental designs of a single run of the analysis

Table 1: Two-dimensional toy function – free p-boxes – failure probability estimation (results of 50 replications of the same analysis with different sample sets and initial experimental designs)

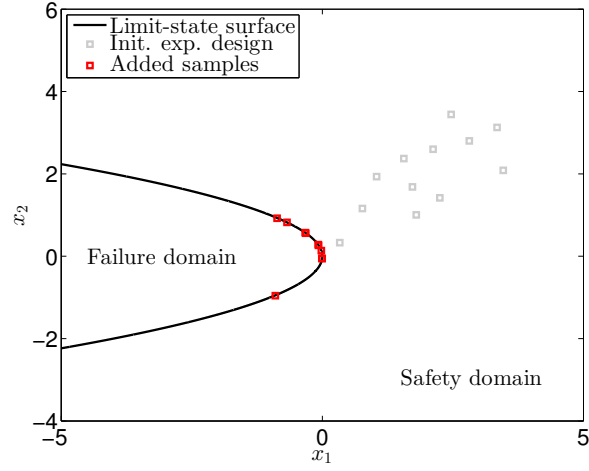
	\tilde{P}_f	\underline{P}_f	\overline{P}_f
$P_{f,ref}$	$1.68 \cdot 10^{-3}$	$7.78 \cdot 10^{-5}$	$1.27 \cdot 10^{-2}$
$\mathbb{E} [\hat{P}_f]$	$1.67 \cdot 10^{-3}$	$7.62 \cdot 10^{-5}$	$1.25 \cdot 10^{-2}$
$\text{CoV} [\hat{P}_f]$	2.5%	13.8%	7.4%
$\mathbb{E} [N_1]$	$12 + 2.7 = 14.7$	-	-
$\text{Std} [N_1]$	0.6	-	-
$\mathbb{E} [N_2]$	-	$4 + 36.9 = 40.9$	$4 + 19.4 = 23.4$
$\text{Std} [N_2]$	-	13.3	17.7

smooth shape of the limit-state surface. In fact, an addition of $\Delta N_1 = 2.7$ points on average is sufficient to model it with sufficient accuracy. On the second level, the number of samples N_2 is larger due to the shape of $\underline{\mathcal{G}}$ and $\overline{\mathcal{G}}$ and the corresponding limit-state surfaces.

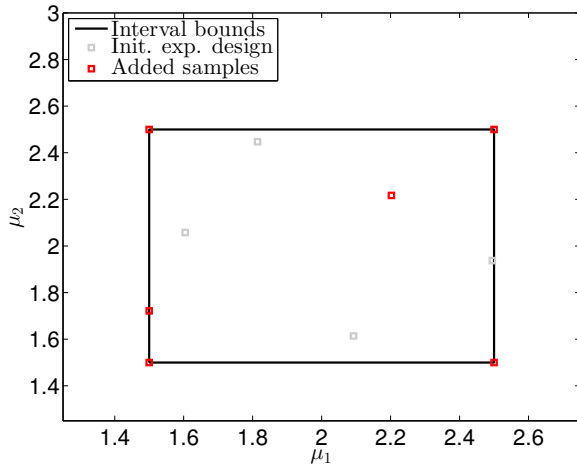
Parametric p-boxes A typical realization of the ISRA for parametric p-boxes is visualized in Figure 5. Figure 5(A) shows the experimental design of the final AK-MCS meta-model. Similar to Figure 4(A), the additional samples (red squares) group efficiently around the limit-state surface which is represented by the solid black line. The experimental design of the EGO algorithm is shown in Figure 5(B), which illustrates the exploratory behaviour of the EGO algorithm. Despite the small number of added samples, the boundary values of the failure probability are estimated accurately, as seen in Figure 5(C). The left side of Figure 5(C) shows the evolution of the bounds $[\underline{P}_f, \overline{P}_f]$ of the failure probability using the initial experimental design (*i.e.* $N_1^0 = 4$ samples in \mathcal{X}^0 are generated simultaneously and the corresponding $\mathcal{Y}^{(i)}$ are computed sequentially), whereas the right side shows them during each iteration of EGO. After iteration 3, the optimization algorithm converges to a stable value of the failure probability bounds.

The results of the 50 replications of the analysis are summarized in Table 2 in terms of failure probability estimates and experimental design sizes. Further, it is distinguished whether the EGO algorithm is used to simultaneously optimize for \underline{P}_f and \overline{P}_f (left block of values) or whether EGO is performed separately for \underline{P}_f and \overline{P}_f (right block of values). Note that for the separate optimization case, samples from one optimization are *not* re-used in the other one. In terms of the failure probability, both settings result in accurate estimates. Moreover, the coefficient of variation is close to the expected value of $\text{CoV}[P_f] = 8.0\%$ and 0.9% for \underline{P}_f and \overline{P}_f , respectively, for a Monte Carlo simulation with $n = 10^6$ samples.

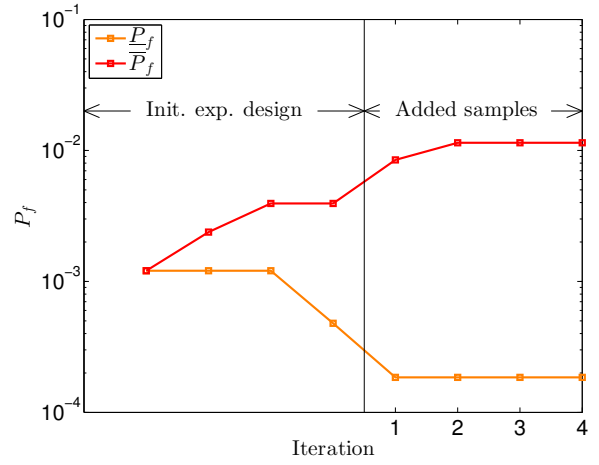
Further, the number of evaluations of the true limit-state function N_1 is similar for the simultaneous and separate optimization cases. The number of samples added during AK-MCS is generally



(A) Experimental design \mathcal{X} for \mathcal{G} (first-level meta-model)



(B) Experimental design \mathcal{T} in \mathcal{D}_{Θ} (EGO)



(C) Evolution of the estimate of the failure probability bounds (EGO)

Figure 5: Two-dimensional toy function – parametric p-boxes – single realization of the ISRA

Table 2: Two-dimensional toy function – parametric p-boxes – failure probability estimation (results of 50 replications of the same analysis with different sample sets and initial experimental designs)

	Simultaneous optimization		Separate optimization	
	\underline{P}_f	\overline{P}_f	\underline{P}_f	\overline{P}_f
$P_{f,ref}$	$1.57 \cdot 10^{-4}$	$1.14 \cdot 10^{-2}$	$1.57 \cdot 10^{-4}$	$1.14 \cdot 10^{-2}$
$\mathbb{E}[\hat{P}_f]$	$1.58 \cdot 10^{-4}$	$1.14 \cdot 10^{-2}$	$1.58 \cdot 10^{-4}$	$1.14 \cdot 10^{-2}$
$\text{CoV}[\hat{P}_f]$	8.0%	1.2%	8.1%	1.2%
$\mathbb{E}[N_1]$	12 + 3.4 = 15.4		12 + 3.4 = 15.4	12 + 3.2 = 15.2
$\text{Std}[N_1]$	0.9		1.0	0.9
$\mathbb{E}[N_2]$	4 + 10.4 = 14.3		4 + 4.7 = 8.7	4 + 2.0 = 6.0
$\text{Std}[N_2]$	1.4		1.2	1.1

low. At the level of EGO, however, the simultaneous optimization requires more samples (N_2) as the separate optimization. Even when adding up the samples for the separate optimization case, $N_2^{tot} = 4 + 4.7 + 2.0 = 10.7$ is less than $N_2 = 14.3$. Hence, the separate optimization is more efficient than the simultaneous optimization in this case. This behaviour has been also observed on other toy functions. Based on this observation, the remaining application examples are analysed using separate optimizations.

Comparison The algorithms for free and parametric p-boxes originate from two different paradigms: interval analysis and nested Monte Carlo simulations, respectively. However, the cost statistics shows that the two proposed two-level meta-modelling algorithms perform similarly in terms of model evaluations. Free p-boxes required $N_1 = 14.7$ evaluations of the limit-state function whereas parametric p-boxes required $N_1 = 15.4$ on average. This can be intuitively understood due to the identical limit-state surface in both settings. Contrarily to N_1 , the number of evaluations on the second-level meta-models, *i.e.* N_2 , are not comparable for the case of free and parametric p-boxes, due to the different algorithms used.

7.2 Single-degree-of-freedom oscillator

7.2.1 Problem setup

In order to visualize the effects of non-monotone functions on the failure probability, the following example is analysed. Consider the non-linear undamped single-degree-of-freedom (SDOF) oscillator sketched in Figure 6 (Echard et al., 2011, 2013; Schueremans and Van Gemert, 2005). The corresponding limit-state function reads:

$$g_{SDOF}(r, F_1, t_1, k_1, k_2, m) = 3r - \left| \frac{2F_1}{m\omega_0^2} \sin\left(\frac{\omega_0 T_1}{2}\right) \right|, \quad (32)$$

where m is the mass, k_1, k_2 are the spring constants of the primary and secondary springs, r is the displacement at which the secondary spring yields, t_1 is the duration of the loading, F_1 is the amplitude of the force and $\omega_0 = \sqrt{\frac{k_1+k_2}{m}}$ is the natural frequency of the oscillator. Failure is defined as $g_{SDOF} \leq 0$ and the associated failure probability is $P_f = \mathbb{P}(g_{SDOF} \leq 0)$.

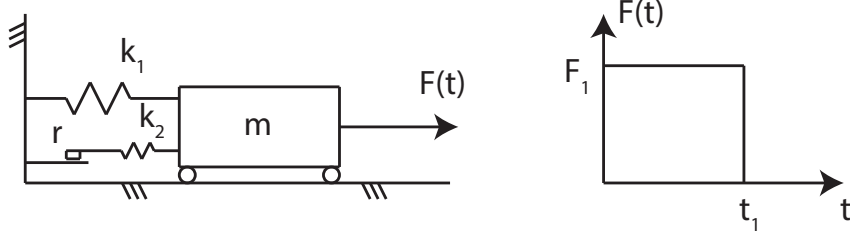


Figure 6: SDOF oscillator – geometry sketch and definition of the variables

The input vector is modelled by a mix of probabilistic variables and p-boxes accounting for the different levels of knowledge. The description of the input variables is provided in Table 3. It is assumed that the spring stiffnesses and the mass are known well. Hence $\{k_1, k_2, m\}$ are modelled by precise CDFs. On the other side, knowledge on $\{r, F_1, t_1\}$ is scarce. Hence, these variables are modelled by p-boxes. The two cases of free and parametric p-boxes are distinguished and compared in the following. As seen in Table 3, the parametric p-box is characterized by a distribution function with interval-valued mean value. For the case of free p-boxes, the same p-box boundary curves are used as for parametric p-boxes. Note that $F_{\mathcal{N}}(x|\mu, \sigma)$ refers to a Gaussian distribution with mean value μ and standard deviation σ .

Table 3: Oscillator – definition of the input vector for the cases of parametric and free p-boxes

Variable	Distribution	Parametric p-box		Free p-box	
		Mean	Std. dev.	\underline{F}_{X_i}	\overline{F}_{X_i}
r	Gaussian	[0.49, 0.51]	0.05	$F_{\mathcal{N}}(r 0.51, 0.05)$	$F_{\mathcal{N}}(r 0.49, 0.05)$
F_1	Gaussian	[-0.2, 0.2]	0.5	$F_{\mathcal{N}}(F_1 0.2, 0.5)$	$F_{\mathcal{N}}(F_1 -0.2, 0.5)$
t_1	Gaussian	[0.95, 1.05]	0.2	$F_{\mathcal{N}}(t_1 1.05, 0.2)$	$F_{\mathcal{N}}(t_1 0.95, 0.2)$
k_1	Gaussian	1	0.1	$F_{\mathcal{N}}(k_1 1, 0.1)$	
k_2	Gaussian	0.1	0.01	$F_{\mathcal{N}}(k_2 0.1, 0.01)$	
m	Gaussian	1	0.05	$F_{\mathcal{N}}(m 1, 0.05)$	

7.2.2 Analysis

The settings for the imprecise structural reliability analyses are kept the same as in Section 7.1, except for the size of the initial experimental designs. The size of the initial experimental designs for all meta-models (including EGO) is set to $N^0 = 12$. Further, the threshold value in EI is set to $\epsilon_{EI} = 10^{-5}$.

For free p-boxes, the auxiliary distributions for $\{r, F_1, t_1\}$ are defined by $F_{\mathcal{N}}(r|0.5, 0.05)$, $F_{\mathcal{N}}(F_1|0, 0.5)$, and $F_{\mathcal{N}}(t_1|1.00, 0.2)$, respectively.

7.2.3 Results

Efficiency of the proposed approaches Table 4 summarizes the results of a single run of the ISRA. The estimates of the failure probability \hat{P}_f are compared to a reference Monte Carlo simulation obtained with $n_{MC} = 10^7$ samples and the exact limit-state function. The corresponding optimizations are using MATLAB's genetic algorithm. All four failure probabilities are estimated accurately. Further, the number of evaluations of the limit-state function N_1 is low. The case of parametric p-boxes results in a larger $N_1 = 197$ compared to the free p-box ($N_1 = 132$). The reason for this is the added samples during the iterations of the EGO. In comparison, the two levels of the free p-box analysis are independent and hence the total N_1 is lower.

Table 4: Oscillator – results of the ISRA (reference solution obtained by Monte Carlo simulation with $n_{MC} = 10^7$ samples)

	Parametric p-box		Free p-box	
	\underline{P}_f	\overline{P}_f	\underline{P}_f	\overline{P}_f
$P_{f,ref}$	$2.42 \cdot 10^{-3}$	$9.04 \cdot 10^{-3}$	$7.08 \cdot 10^{-4}$	$1.63 \cdot 10^{-2}$
\hat{P}_f	$2.41 \cdot 10^{-3}$	$9.04 \cdot 10^{-3}$	$6.69 \cdot 10^{-4}$	$1.61 \cdot 10^{-2}$
N_1	$12 + 185 = 197$	$12 + 185 = 197$	$12 + 120 = 132$	
N_2	$6 + 3 = 9$	$6 + 6 = 12$	$12 + 60 = 82$	$12 + 153 = 165$
μ_r^*	1.02	0.98	-	-
$\mu_{F_1}^*$	-0.01	0.20	-	-
$\mu_{t_1}^*$	0.95	1.05	-	-

The number of samples in the second-level meta-models are varying more than the first-level meta-models. In the case of parametric p-boxes, a few samples $N_2 = 3 \dots 6$ are sufficient to find the optimal distribution parameters. In the case of free p-boxes, N_2 has the same order of magnitude as N_1 because the analysis is of the same type, *i.e.* AK-MCS. However, the difference in N_2 of estimating \underline{P}_f and \overline{P}_f is large, which represents the different complexity of the limit-state surfaces $\overline{\mathcal{G}} = 0$ and $\underline{\mathcal{G}} = 0$, respectively.

For the case of parametric p-boxes, Table 4 shows also the optimal mean values for $\{r, F_1, t_1\}$, denoted by $\{\mu_r^*, \mu_{F_1}^*, \mu_{t_1}^*\}$, which result from the EGO optimization. The results confirm the intuition: the largest maximal failure probability is obtained by the smallest yield displacement in combination with a long activation time and an eccentric force ($\mu_{F_1} \approx 0.2$ or $\mu_{F_1} \approx -0.2$). On the other hand, the minimum failure probability is obtained by the largest yield displacement in combination with a short activation time and a concentric force ($\mu_{F_1} \approx 0$).

Evolution of experimental designs As mentioned above, the evolution of the experimental design is different for the two types of p-boxes. For parametric p-boxes, the details are shown in Figure 7. The size of N_1 is plotted as a function of the iteration in EGO in Figure 7(B). Interestingly, the major part of evaluations are required to estimate the first sample of EGO’s initial experimental design. Only a few additional samples are evaluated in order to refine the conditional failure probability estimates. Hence, the recycling of limit-state function evaluations is an efficient tool to keep the computational costs at a low level.

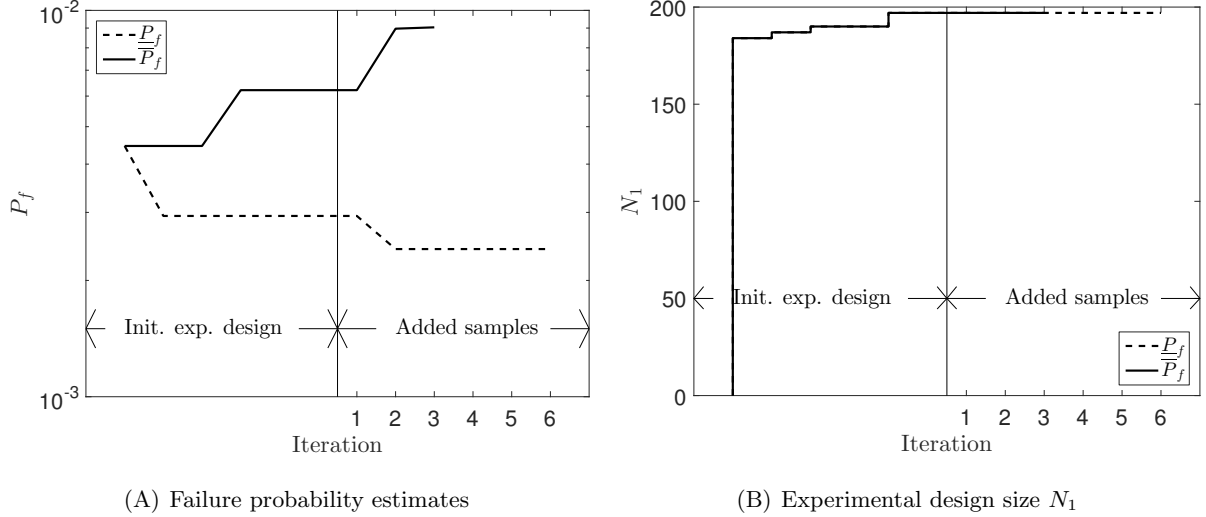


Figure 7: Oscillator – parametric p-boxes – convergence of the ISRA

In the same figure, the evolution of the boundary values of the failure probability are also given (see Fig. 7(A)). Interestingly, the failure probabilities evolve with each iteration in EGO where at the same time N_1 remains constant. In other words, the limit-state surface $\mathcal{G} = 0$ is modelled accurately enough to estimate the extreme failure probabilities, *i.e.* for different realizations $\theta \in \mathcal{D}_\theta$.

Effect of non-monotonicity on failure probability In this example, the two types of p-boxes have the same boundary curves, as seen in Table 3. In case the limit-state function was a monotone function, the failure probabilities would be identical. In other words, the same realizations of the input CDF would lead to one of the failure probability boundary values.

In the case of the oscillator though, parametric and free p-box models result in significantly different failure probability bounds, as seen from Table 4. In fact, the imprecise failure probabilities of the free p-box case encapsulate the ones of the parametric p-box case. As free p-boxes are more general than parametric p-boxes by definition, more extreme failure probabilities are possible. In this example, free p-boxes result in a 3.4-times lower \underline{P}_f and in a 1.8-times larger \bar{P}_f compared to the parametric p-boxes.

7.3 Two-dimensional truss structure

7.3.1 Problem statement

Hurtado Hurtado (2013) introduced a two-dimensional linear-elastic truss structure, whose geometry is shown in Figure 8. The truss is subjected to seven loads P_i which are modelled with p-boxes. The parametric p-boxes are defined by lognormal distributions with mean value $\mu_{P_i} \in [95, 105]$ kN and standard deviation $\sigma_{P_i} \in [13, 17]$ kN. The geometry of the structure and the material properties are assumed constant. The modulus of elasticity is $E = 200 \cdot 10^9$ Pa, whereas the cross section area varies along the different bars: $A = 0.00535$ m² for bars marked by \bullet , $A = 0.0068$ m² for bars marked by \circ , and $A = 0.004$ m² for the remaining bars.

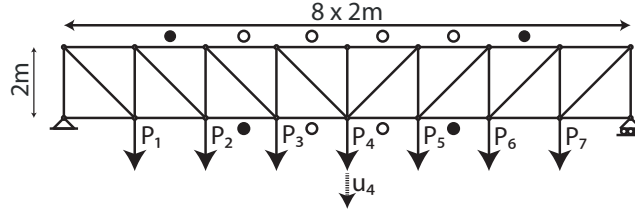


Figure 8: Frame structure – sketch of the geometry and definition of the input variables

In the following analysis, a second scenario is considered, where the former parametric p-boxes are modelled by free p-boxes. As in the other examples, the boundary curves of the p-boxes should coincide, hence:

$$\underline{F}_{P_i}(p_i) = \min_{\substack{\mu_{P_i} \in [95, 105] \\ \sigma_{P_i} \in [13, 17]}} F_{P_i}(p_i | \mu_{P_i}, \sigma_{P_i}), \quad (33)$$

$$\overline{F}_{P_i}(p_i) = \max_{\substack{\mu_{P_i} \in [95, 105] \\ \sigma_{P_i} \in [13, 17]}} F_{P_i}(p_i | \mu_{P_i}, \sigma_{P_i}). \quad (34)$$

In the context of structural reliability analysis, the limit-state function is defined as

$$g_{truss}(\mathbf{p}) = 0.029 \text{ m} - u_4(\mathbf{p}), \quad (35)$$

where u_4 is the deflection of the truss at midspan, as indicated in Figure 8. Then, the failure probability describes the probability that the deflection of the truss exceeds 0.029 m, *i.e.* $P_f = \mathbb{P}(u_4(\mathbf{P}) \geq 0.029 \text{ m})$.

7.3.2 Analysis

The deflection of the truss is computed by a MATLAB-based finite element model (FEM), which is connected to the software framework UQLAB (Marelli and Sudret, 2014). The FEM interprets each bar as a bar element, whereas the loads are modelled as point loads at the intersections of the corresponding bars as indicated in Figure 8.

The ISRA settings are kept the same as in the previous examples. However, the initial experimental design is set to $N_1^0 = N_2^0 = 12$ Latin-hypercube samples and $\epsilon_{EI} = 10^{-5}$. For free p-boxes, the auxiliary

distributions \tilde{X}_i are chosen as lognormal distributions with mean value $\mu_{\tilde{P}_i} = 100$ kN and standard deviation $\mu_{\tilde{P}_i} = 15$ kN.

7.3.3 Results

Table 5 summarizes the results in terms of failure probability estimates and number of model evaluations. The reference values are obtained by Monte Carlo simulation with $n = 10^7$ samples, by evaluating the exact limit-state function, and by making use of the monotonicity property of the limit-state function. Using the proposed two-level algorithms, the failure probabilities are estimated accurately for both cases of parametric and free p-boxes. The numbers of FEM evaluations N_1 are low for both cases, too. Hence, the proposed multi-level meta-modelling techniques provide efficient algorithms for ISRA in this application example.

Table 5: Truss – results of the ISRA (reference solution obtained by Monte Carlo simulation with $n_{MC} = 10^7$ samples)

	Parametric p-box		Free p-box	
	\underline{P}_f	\overline{P}_f	\underline{P}_f	\overline{P}_f
$P_{f,ref}$	$2.49 \cdot 10^{-4}$	$7.62 \cdot 10^{-2}$	$2.21 \cdot 10^{-4}$	$9.22 \cdot 10^{-2}$
\hat{P}_f	$2.64 \cdot 10^{-4}$	$7.62 \cdot 10^{-2}$	$2.36 \cdot 10^{-4}$	$9.16 \cdot 10^{-2}$
N_1	$12 + 130 = 142$	$12 + 129 = 141$	$12 + 106 = 118$	$12 + 106 = 118$
N_2	$12 + 174 = 186$	$12 + 40 = 52$	$12 + 116 = 128$	$12 + 192 = 204$
N_2^*	$12 + 26 = 38$	$12 + 4 = 16$	-	-

In the case of parametric p-boxes, the minimum failure probability is obtained by $\mu_{P_i} = 95$ kN and $\sigma_{P_i} = 13$ kN, whereas the maximum failure probability is obtained by $\mu_{P_i} = 105$ kN and $\sigma_{P_i} = 17$ kN, $i = 1, \dots, 7$, as one could expect. EGO is capable of identifying these two extreme cases and terminating the adaptive algorithm in this application in 174 and 40 iterations, which correspond to $N_2 = 186$ and $N_2 = 52$ samples in the EGO-related Kriging model, respectively. In this run, the number of iterations required to find those extremes is indicated in Table 5 by N_2^* . The values are lower than the final number of samples N_2 , which are used to terminate the EGO algorithm. The additional samples beyond N_2^* show the exploratory behaviour of the EGO algorithm in this application example and might be reduced by a more appropriate stopping criterion.

Hurtado Hurtado (2013) analysed the same truss structure for the parametric p-box case. However, due to its use of a Monte Carlo simulation on the second level of the algorithm, the published failure probability range is narrower, *i.e.* $P_f \in [3.2 \cdot 10^{-3}, 2.37 \cdot 10^{-2}]$, than reported in Table 5, which is arguably not conservative. This indicates the importance of using a proper optimization algorithm in order to find the extreme failure probabilities.

8 Conclusions

In this paper, structural reliability analysis is discussed in the presence of deterministic, black-box performance functions. Traditionally, uncertainty in the input variables of such analyses is described by probability theory. In engineering practice however, a typical scenario consists of expensive data acquisition and thus limited data. Due to the sparsity of calibration data, a more general framework is required to characterize the input uncertainty.

In this paper, the input variables are modelled as imprecise probabilities accounting for aleatory (natural variability) and epistemic uncertainty (lack of knowledge). Probability-boxes (p-boxes) are one methodology of the class of imprecise probabilities that captures both types of uncertainty in a clear setting. In particular, two types of p-boxes are distinguished: free and parametric p-boxes. Free p-boxes model a variable by providing lower and upper bounds for the cumulative distribution function. Parametric p-boxes consist of a distribution function family whose parameters are modelled within intervals.

In the context of p-boxes, structural reliability analyses are more complex than in the conventional probabilistic setting. Moreover, free and parametric p-boxes require different approaches. In order to cope with the complexity of imprecise structural reliability analyses, meta-models are used at different stages. In this paper, two two-level meta-modelling approaches are proposed for the cases of free and parametric p-boxes, respectively, making use of Kriging meta-models in combination with active learning algorithms. These approaches allow for an accurate estimation of the imprecise failure probabilities using only a limited number of runs of the performance function.

The capabilities of the two-level approaches are illustrated on a benchmark analytical function and two realistic engineering problems. In all examples, the proposed approaches are capable of estimating the imprecise failure probabilities in an efficient manner. In fact, the imprecise and conventional structural reliability analyses require a comparable number of runs of the performance function when using Kriging meta-models with adaptive experimental designs. This is of importance when the evaluations of the performance function are the dominating computational costs.

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