Failure Probability Estimation of a Class of Series Systems by Multidomain Line Sampling

- Marcos A. Valdebenito^{a,*}, Pengfei Wei^{b,d}, Jingwen Song^{c,d}, Michael Beer^{d,e,f}, Matteo Broggi^d
- ^aFaculty of Engineering and Sciences, Universidad Adolfo Ibáñez, Av. Padre Hurtado 750, 2562340 Viña del Mar, Chile
- ^bSchool of Mechanics, Civil Engineering and Architecture, Northwestern Polytechnical University, Xi'an 710072, China
 - ^cAdvanced Research Laboratories, Tokyo City University, 1-28-1 Tamazutsumi Setagaya-ku, Tokyo 158-8557, Japan
 - ^dInstitute for Risk and Reliability, Leibniz Universität Hannover, Callinstr. 34, 30167 Hannover, Germany ^eInstitute for Risk and Uncertainty and School of Engineering, University of Liverpool, Peach Street, Liverpool L69 7ZF, UK
 - ^fInternational Joint Research Center for Engineering Reliability and Stochastic Mechanics, Tongji University, 1239 Siping Road, Shanghai 200092, P.R. China

Abstract

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This contribution proposes an approach for the assessment of the failure probability associated with a particular class of series systems. The type of systems considered involves components whose response is linear with respect to a number of Gaussian random variables. Component failure occurs whenever this response exceeds prescribed deterministic thresholds. We propose multidomain Line Sampling as an extension of the classical Line Sampling to work with a large number of components at once. By taking advantage of the linearity of the performance functions involved, multidomain Line Sampling explores the interactions that occur between failure domains associated with individual components in order to produce an estimate of the failure probability. The performance and effectiveness of multidomain Line Sampling is illustrated by means of two test problems and an application example, indicating that this technique is amenable for treating problems comprising both a large number of random variables and a large number of components. Keywords: Line sampling, Multidomain, Linear performance function, Failure probability, Series system

^{*}E-mail: marcos.valdebenito@uai.cl

29 Highlights:

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- Failure probability of series system is calculated by multidomain Line Sampling.
- Knowledge on individual component failure domains is exploited.
 - Several important directions are considered simultaneously.
 - Lines explore interaction between failure events associated with components.

1. Introduction

An engineering system can be seldom described precisely, as different sources of uncertainty 35 may affect its performance. Whenever the nature of uncertainty is of the aleatory type, it is possible to resort to probability theory for analyzing such system [1]. In this way, some input parameters of the system are modeled as random variables (or random processes or random fields, in case time or spatial correlations are present). In turn, such description of the uncertain input parameters causes that the performance of the system becomes random as well. Due to design, operation or maintenance purposes, it is of interest assessing the level of safety associated with 41 the performance of a system, for example, in terms of a failure probability, that measures the chances of an undesirable behavior. At this point, it should be noted that an engineering system usually comprises a number of components, each of which may possess a different failure probability and whose performance may be correlated with that of other components. Depending on the configuration of those components within the system, it may be of interest calculating the failure probability associated with different types of system events: simultaneous failure of all components of the system (parallel event), failure of one or more components (series event), etc. Often, quantifying such failure probability is far from trivial and hence, a number of specialized approaches have been developed for calculating it, for example: bounds based on failure probabili-50 ties of individual components and interactions between two [2] or three components [3]; application of surrogate models [4, 5, 6, 7, 8]; linear programming [9, 10, 11] or binary programming [12] techniques; approximation concepts by compounding individual component failure events [13, 14, 15], sampling approaches [16, 17, 18, 19, 20], etc. Although the previous list of contributions is far from being extensive, it demonstrates that calculation of failure probabilities involving different system events is a field of active research.

This contribution focuses on the calculation of the failure probability associated with a series event

of a system. In other words, the objective is calculating the probability that the performance of one or more components of a system exceeds a prescribed threshold level. The class of problems considered herein pertain components whose response is characterized as a linear combination of Gaussian random variables. Failure of the component occurs whenever the response falls below prescribed lower threshold or exceeds a prescribed upper threshold. This class of problems has attracted considerable attention in the literature due to its applications in, e.g. reliability of time-variant systems [21, 22, 23], stochastic linear dynamics [24, 25], seismic fragility analysis [10], geotechnical applications [26, 27], network analysis [28], etc. The focus is on problems that involve a large number of random variables and a large number of components, possibly in the order of thousands. It should be noted that the type of problems considered in this contribution possesses a distinctive geometry in the standard Gaussian space, where the boundary of the failure domain involves a number of hyperplanes [24]. Such distinctive geometry has allowed to design simulation schemes that allow calculating the sought failure probability with high accuracy and efficiency using either 71 concepts of Importance Sampling [29], Domain Decomposition with averages [30] and Directional Importance Sampling [31, 32]. This work builds on that knowledge of the failure domain and 73 proposes multidomain Line Sampling (mLS), which is a novel extension of Line Sampling [33, 34]. The salient feature of mLS is that it is capable of dealing with failure domains associated with 75 multiple components. This is achieved by introducing multiple search directions instead of a sin-76 gle search direction as usually considered in classical Line Sampling. In addition, lines simulated 77 during the calculation of probability are conditioned to lie in the failure domain, which allows exploring the interaction between failure events associated with individual components along that line. The rest of this work is organized as follows. Section 2 formulates the failure probability problem associated with a series event of a system. Section 3 presents the description and formulation of multidomain Line Sampling. Section 4 illustrates the application of multidomain Line Sampling to two test problems and an application example, the latter involving a large number of random variables and components. The paper closes with a conclusions and outlook for future developments in Section 5.

2. Formulation of the Problem

2.1. General Aspects

Consider a series system involving a total of n_c components. The behavior of each of these components is described in terms of a response $r_i(\boldsymbol{x}), i = 1, \ldots, n_c$, that depends on a parameter vector \boldsymbol{x} of dimension n. The component exhibits an acceptable behavior whenever its response lies within prescribed thresholds, that is $b_i^L < r_i(\boldsymbol{x}) < b_i^U$. In other words, failure of the component occurs whenever the response either falls below b_i^L or exceeds b_i^U . Note that no particular restrictions must be imposed regarding the thresholds other than $b_i^L < b_i^U$, $i = 1, \ldots, n_c$.

It is assumed that the parameter vector \boldsymbol{x} is uncertain and is characterized by means of a random variable vector \boldsymbol{X} that follows a Gaussian multivariate distribution with mean $\boldsymbol{\mu}$ and (positive definite) covariance matrix \boldsymbol{C} . The parameter vector \boldsymbol{x} can be represented in the standard Gaussian space as:

$$x = \mu + Bz \tag{1}$$

where z is a realization of Z, which follows an n-dimensional standard Gaussian distribution; and where B is a matrix that can be calculated, for example, using Cholesky decomposition or spectral 100 representation. In case the latter is applied, it is noted that ${m B}={m \Phi}{m \Lambda}^{1/2},$ where the columns of 101 matrix Φ contain the eigenvectors of C while the diagonal of matrix Λ contains the corresponding 102 eigenvalues of C. It is further assumed that the response associated with each component depends linearly on \boldsymbol{x} , that is $r_i(\boldsymbol{x}) = \boldsymbol{a}_i^T \boldsymbol{x}$, where \boldsymbol{a}_i is an *n*-dimensional vector with real entries and $(\cdot)^T$ 104 denotes transpose. 105 Taking into account the previous assumptions, it is possible to formulate two performance functions 106 associated with the i-th component: one for monitoring whenever the response falls below the 107 threshold b_i^L and the other one for monitoring whenever the response exceeds the threshold b_i^U . These functions are equal to:

$$g_{2i-1}(\boldsymbol{z}) = \beta_i^L + \boldsymbol{\alpha}_i^T \boldsymbol{z}, \ i = 1, \dots, n_c$$
 (2)

$$g_{2i}(\mathbf{z}) = \beta_i^U - \boldsymbol{\alpha}_i^T \mathbf{z}, \ i = 1, \dots, n_c$$
(3)

where $\beta_i^L = \left(\boldsymbol{a}_i^T \boldsymbol{\mu} - b_i^L\right) / ||\boldsymbol{a}_i^T \boldsymbol{B}||$, $\beta_i^U = \left(b_i^U - \boldsymbol{a}_i^T \boldsymbol{\mu}\right) / ||\boldsymbol{a}_i^T \boldsymbol{B}||$, $\boldsymbol{\alpha}_i = \boldsymbol{a}_i^T \boldsymbol{B} / ||\boldsymbol{a}_i^T \boldsymbol{B}||$ and $||\cdot||$ denotes Euclidean norm. Note that the formulation of the performance function in eq. (2) is

actually equal to the subtraction between the response r_i (which has been expressed in terms of vector \mathbf{z} applying eq. (1)) and the threshold b_i^L , divided by the Euclidean norm of vector $\mathbf{a}_i^T \mathbf{B}$. This ensures that $g_{2i-1}(\mathbf{z})$ assumes a value equal or smaller than zero whenever the response equals or is below the threshold b_i^L . In a similar way, eq. (3) is constructed as the subtraction between the threshold level b_i^U and the response r_i , divided by the Euclidean norm of vector $\mathbf{a}_i^T \mathbf{B}$. Thus, $g_{2i}(\mathbf{z})$ assumes a value equal or smaller than zero whenever the response equals or exceeds the threshold b_i^U . Note that in the previous definitions of the performance functions, normalization by the Euclidean norm of vector $\mathbf{a}_i^T \mathbf{B}$ is enforced as this ensures that $||\mathbf{a}_i|| = 1, i = 1, \ldots, n_c$.

o 2.2. Failure Probability Associated with Individual Component

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Different realizations z of the random vector Z may cause failure of the *i*-th component. The 121 set of all of these realizations is denoted as the failure domain F_i . In turn, this failure domain 122 is the union of a negative (F_i^-) and a positive failure domain (F_i^+) , that is $F_i = F_i^- \cup F_i^+$. The 123 negative failure domain F_i^- is a set that groups all realizations \boldsymbol{z} such that the response of the i-th component is equal to or below the threshold b_i^L , that is $F_i^- = \{ z \in \mathbb{R}^n : g_{2i-1}(z) \leq 0 \}$, i = 1 $1, \ldots, n_c$. In a similar way, the positive elementary failure domain F_i^+ groups all realizations zsuch that the response of the i-th component is equal to or exceeds the threshold b_i^U , that is $F_i^+ = \{ \boldsymbol{z} \in \mathbb{R}^n : g_{2i}(\boldsymbol{z}) \leq 0 \}, \ i = 1, \dots, n_c.$ In view of the linearity of each of the performance functions g_i with respect to z as noted from eqs. (2) and (3), negative and positive failure domains are bounded by hyperplanes. Furthermore, the negative and positive failure domains are fully described by their corresponding design points. Recall that the design point z^* is the realization of Z with smallest Euclidean norm with respect to the origin that causes failure. It is straightforward to demonstrate that the design points associated with each performance function are equal to $\mathbf{z}_{2i-1}^* = -\beta_i^L \boldsymbol{\alpha}_i$ and $\mathbf{z}_{2i}^* = \beta_i^U \boldsymbol{\alpha}_i$, $i = 1, \dots, n_c$, respectively [24, 29]. Figure 1 provides a schematic representation of the negative and positive failure domains as well as their corresponding design points for the specific case where n=2 and $n_c=1$. 136

The probability of failure of the *i*-th component is denoted as $p_{F,i}$ and is defined as:

$$p_{F,i} = \int_{\boldsymbol{z} \in \mathbb{R}^n} I_{F_i}(\boldsymbol{z}) f_{\boldsymbol{Z}}(\boldsymbol{z}) d\boldsymbol{z}$$
(4)

where $f_{\mathbf{Z}}(\mathbf{z})$ is the standard Gaussian probability density function in n dimensions; and where $I_{F_i}(\mathbf{z})$ is the indicator function associated with the i-the failure event, which is equal to $I_{F_i}(\mathbf{z}) = 1$

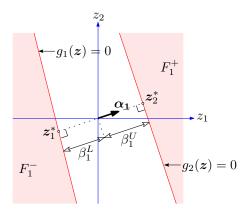


Figure 1: Schematic representation of negative and positive failure domains (F_i^- and F_i^+ , respectively) and their design points ($n = 2, n_c = 1$).

in case $z \in F_i$ and zero, otherwise. In view of the linearity of the performance functions g_{2i-1} and g_{2i} with respect to z, the probability integral in eq. (4) possesses an analytic solution [2], which is equal to:

$$p_{F,i} = \Phi\left(-\beta_i^L\right) + \Phi\left(-\beta_i^U\right) \tag{5}$$

where $\Phi(\cdot)$ is the standard Gaussian cumulative density function. Note that β_i^L and β_i^U are actually the reliability indexes associated with F_i^- and F_i^+ , respectively [2, 24]. In other words, they are the Euclidean norm of the corresponding design points, that is $\beta_i^L = ||\boldsymbol{z}_{2i-1}^*||$ and $\beta_i^U = ||\boldsymbol{z}_{2i}^*||$, $i = 1, \ldots, n_c$ [24].

2.3. Failure Probability Associated with Series Event

The failure event associated with a series system implies that one or more of its components fail. The failure domain F groups all realizations z of the random variable vector Z that cause the failure event, that is, $F = F_1 \cup F_2 \cup ... \cup F_{n_c}$. The probability of failure associated with the systems event is denoted as p_F and is defined as:

$$p_F = \int_{\boldsymbol{z} \in \mathbb{R}^n} I_F(\boldsymbol{z}) f_{\boldsymbol{Z}}(\boldsymbol{z}) d\boldsymbol{z}$$
 (6)

where $I_F(z)$ is the indicator function, which is equal to $I_F(z) = 1$ in case $z \in F$ and zero, otherwise. It is important to note that for most cases of practical interest, eq. (6) cannot be solved in closed form [13]. This is due to the fact that interactions between the failure domains associated with individual components cannot be analyzed analytically. In addition to the issue of interactions, the number of random variables n and of performance functions n_c associated with the probability integral may be considerable (in the order of hundreds or thousands). These two issues favor the application of simulation methods for calculating the failure probability [35].

Line Sampling is a simulation technique which was developed for calculating failure probabil-

3. Multidomain Line Sampling

160 3.1. Line Sampling

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ities in problems involving a large number of random variables [33, 36]. It is closely related to 162 another simulation technique known as Axis Orthogonal Sampling [37, 38]. Most of the appli-163 cations of Line Sampling which are available in the literature focus on the assessment of failure 164 probabilities associated with weakly or moderately nonlinear performance functions of an individ-165 ual component, see e.g. [39]. 166 The practical implementation of Line Sampling requires that the reliability problem is formulated 167 in the standard Gaussian space by means of a suitable projection [40]. After that, it is necessary 168 to identify the so-called important direction γ , which is a vector of unit Euclidean norm located at the origin of the standard normal space that points towards the failure domain. Several criteria 170 have been proposed for determining such direction [33, 41]. Then, taking advantage of the rota-171 tional invariance of the standard Gaussian distribution, a rotated coordinate system is introduced, 172 such that:

$$z = Rz^{\perp} + \gamma z^{\parallel} \tag{7}$$

where z^{\perp} is a vector of dimension (n-1) that represents coordinates in the hyperplane orthogonal to γ ; z^{\parallel} is a scalar denoting the coordinate parallel to γ ; and R is a matrix of dimension $n \times (n-1)$. The square matrix $[\mathbf{R}, \boldsymbol{\gamma}]$ forms an orthonormal basis and thus, it is straightforward 176 to demonstrate that $z^{\parallel} = \boldsymbol{\gamma}^T \boldsymbol{z}$ and $\boldsymbol{z}^{\perp} = \boldsymbol{R}^T \boldsymbol{z}$. Note that for practical implementation, there is no 177 need to determine matrix R in explicit form. In addition, note that the probability distributions 178 associated with z^{\perp} and z^{\parallel} are standard Gaussian distributions in (n-1) dimensions and one dimension, respectively. 180 Line Sampling takes advantage of the rotated coordinate system associated with eq. (7) by com-181 bining simulation with numerical integration. That is, random samples are generated in the hyper-182 plane orthogonal to the important direction γ . These samples are denoted as $\mathbf{z}^{\perp,(j)}$, $j=1,\ldots,N$, where N denotes the number of samples. Then, one-dimensional numerical integration is performed along the line $l^{(j)}$, $j=1,\ldots,N$, that is parallel to the important direction and that contains the sample $\mathbf{z}^{\perp,(j)}$. The aim of this one-dimensional integration is determining which portions of the line contribute to the failure probability integral. The whole procedure is depicted schematically in Figure 2, where the dimension of the problem is n=2 and the number of lines is set equal to N=2. Note that in this figure, the performance function is denoted as $g(\mathbf{z})$.

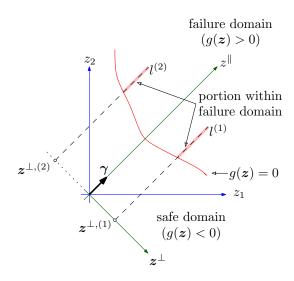


Figure 2: Schematic representation of Line Sampling considering two lines (n = N = 2).

3.2. Formulation of multidomain Line Sampling

Recall that the objective of this work is formulating a simulation scheme for calculating fail-191 ure probabilities of series systems involving a large number of components whose performance 192 functions are linear with respect to a set of parameters following a Gaussian distribution. In 193 order to develop such simulation scheme, note that the summation of the failure probabilities of 194 individual components provides an upper bound for the failure probability of the series event, 195 that is $p_F \leq p_{F,1} + p_{F,2} + \ldots + p_{F,n_c}$ [2, 29]. Such inequality can be understood with the help of the schematic representation in Figure 3, where it is assumed for simplicity that $n = n_c = 2$. 197 When examining this figure, it is noted that failure domains associated with components F_1 and F_2 exhibit overlap; in fact, this overlap occurs at each of the four corners of the figure, that is $F_1^+ \cap F_2^+$ (upper-right corner), $F_1^+ \cap F_2^-$ (lower-right corner), $F_1^- \cap F_2^-$ (lower-left corner) and $F_1^- \cap F_2^+$ (upper-left corner). This implies that the quantity $p_{F,1} + p_{F,2}$ must be necessarily larger than p_F , as the probability content associated with those overlapping regions is being counted twice. In other words, direct summation of the probabilities of failure of individual components does not take into account the possible interactions between the failure domains associated with each component, which in Figure 3 correspond to the realizations of z that belong to any of the sets $F_1^+ \cap F_2^+$, $F_1^+ \cap F_2^-$, $F_1^- \cap F_2^-$ and $F_1^- \cap F_2^+$.

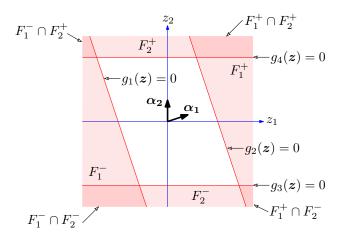


Figure 3: Schematic representation of failure domains associated with components $(n = n_c = 2)$.

The overlap between failure domains associated with individual components can be taken into account by explicitly modeling their interaction [29, 30]. For that purpose, let p_i be equal to:

$$p_i = \int_{\boldsymbol{z} \in F_i} \frac{1}{\sum_{k=1}^{n_c} I_{F_k}(\boldsymbol{z})} f_{\boldsymbol{z}}(\boldsymbol{z}) d\boldsymbol{z}, i = 1, \dots, n_c$$
 (8)

Note that p_i is the integral over the set F_i of the standard Gaussian probability density function divided by the sum of individual failure events that are associated with a particular realization z.

Clearly, p_i as defined in eq. (8) is different from the probability of failure of the i-th component $p_{F,i}$ (see eq. (4)). In fact, the quantity p_i can be loosely interpreted as the effective contribution of the i-th component to the failure probability associated with the series event, where interaction with other components is discounted by means of the factor $1/\sum_{k=1}^{n_c} I_{F_k}(z)$. It is readily seen that:

$$\sum_{i=1}^{n_c} p_i = \sum_{i=1}^{n_c} \left(\int_{\mathbf{z} \in F_i} \frac{1}{\sum_{k=1}^{n_c} I_{F_k}(\mathbf{z})} f_{\mathbf{Z}}(\mathbf{z}) d\mathbf{z} \right)$$

$$= \sum_{i=1}^{n_c} \left(\int_{\mathbf{z} \in F} \frac{I_{F_i}(\mathbf{z})}{\sum_{k=1}^{n_c} I_{F_k}(\mathbf{z})} f_{\mathbf{Z}}(\mathbf{z}) d\mathbf{z} \right)$$

$$= \int_{\mathbf{z} \in F} \frac{\sum_{i=1}^{n_c} I_{F_i}(\mathbf{z})}{\sum_{k=1}^{n_c} I_{F_k}(\mathbf{z})} f_{\mathbf{Z}}(\mathbf{z}) d\mathbf{z}$$

$$= \int_{\mathbf{z} \in F} f_{\mathbf{Z}}(\mathbf{z}) d\mathbf{z} = \int_{\mathbf{z} \in \mathbb{R}^n} I_{F}(\mathbf{z}) f_{\mathbf{Z}}(\mathbf{z}) d\mathbf{z} = p_F$$

$$(9)$$

which implies that the summation of all p_i , $i = 1, ..., n_c$ is equal to the failure probability 216 p_F . For a better understanding of the above equation, consider again Figure 3. The quantity p_1 217 would be equal to the integral of the standard Gaussian probability density function over the set $F_1 \setminus F_2$ (where $(\cdot) \setminus (\cdot)$ denotes set subtraction) plus one half of the standard Gaussian probability distribution over set $F_1 \cap F_2$. In a similar way, p_2 would be equal to the integral of the standard 220 Gaussian probability density function over the set $F_2 \setminus F_1$ plus one half of the standard Gaussian 221 probability distribution over set $F_1 \cap F_2$. Clearly, the summation of p_1 and p_2 would be equal to 222 p_F , as in this case, the interaction between components has been accounted for by means of the factor $1/\sum_{k=1}^{n_c} I_{F_k}(\boldsymbol{z})$. 224 The calculation of the probability of failure of a series event as proposed in eq. (9) demands calculating the quantities p_i , $i = 1, ..., n_c$. These quantities can be evaluated by means of Line 226 Sampling. For that purpose, consider the failure domain associated with the i-th component: an obvious choice for the important direction would be $\gamma = \alpha_i$. Then, a rotated coordinate system is introduced such that:

$$\boldsymbol{z} = \boldsymbol{R}_i \boldsymbol{z}_i^{\perp} + \boldsymbol{\alpha}_i \boldsymbol{z}_i^{\parallel} \tag{10}$$

where \mathbf{z}_i^{\perp} and z_i^{\parallel} denote the set of coordinates which are orthogonal and parallel to $\boldsymbol{\alpha}_i$, respectively; and where \boldsymbol{R}_i denotes the corresponding matrix for coordinate rotation. Figure 4 provides a schematic illustration of the different rotated coordinate systems for a case where $n = n_c = 2$.

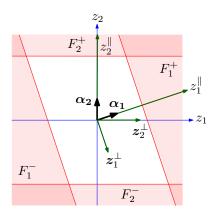


Figure 4: Schematic representation of multidomain Line Sampling considering two rotated coordinate systems $(n = n_c = 2)$.

Assuming that the square matrix $[\mathbf{R}_i, \boldsymbol{\alpha}_i]$ forms an orthonormal basis, the probability distributions associated with \boldsymbol{z}_i^{\perp} and $\boldsymbol{z}_i^{\parallel}$ are standard Gaussian in (n-1) dimensions and one dimension, respectively. Thus, taking into account eqs. (8) and (10), the integral associated with the quantity

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 p_i is recast as:

$$p_{i} = \int_{\boldsymbol{z}_{i}^{\perp} \in \mathbb{R}^{n-1}} \int_{\left(\boldsymbol{z}_{i}^{\parallel}, \boldsymbol{z}_{i}^{\perp}\right) \in F_{i}} \frac{1}{\sum_{k=1}^{n_{c}} I_{F_{k}} \left(\boldsymbol{R}_{i} \boldsymbol{z}_{i}^{\perp} + \boldsymbol{\alpha}_{i} \boldsymbol{z}_{i}^{\parallel}\right)} f_{Z_{i}^{\parallel}} \left(\boldsymbol{z}_{i}^{\parallel}\right) f_{\boldsymbol{Z}_{i}^{\perp}} \left(\boldsymbol{z}_{i}^{\perp}\right) d\boldsymbol{z}_{i}^{\parallel} d\boldsymbol{z}_{i}^{\perp}, \ i = 1, \dots, n_{c} \ (11)$$

The last equation can be further simplified by taking into account that the performance functions associated with the *i*-th failure domain in the rotated coordinates are equal to $g_{2i-1}\left(z_i^{\parallel}, \boldsymbol{z}_i^{\perp}\right) = \beta_i^L + z_i^{\parallel}$ and $g_{2i}\left(z_i^{\parallel}, \boldsymbol{z}_i^{\perp}\right) = \beta_i^U - z_i^{\parallel}$, yielding:

$$p_{i} = \int_{\boldsymbol{z}_{i}^{\perp} \in \mathbb{R}^{n-1}} \int_{-\infty}^{-\beta_{i}^{L}} \frac{1}{\sum_{k=1}^{n_{c}} I_{F_{k}} \left(\boldsymbol{R}_{i} \boldsymbol{z}_{i}^{\perp} + \boldsymbol{\alpha}_{i} \boldsymbol{z}_{i}^{\parallel}\right)} f_{\boldsymbol{Z}_{i}^{\parallel}} \left(\boldsymbol{z}_{i}^{\parallel}\right) f_{\boldsymbol{Z}_{i}^{\perp}} \left(\boldsymbol{z}_{i}^{\perp}\right) d\boldsymbol{z}_{i}^{\parallel} d\boldsymbol{z}_{i}^{\perp} + \int_{\boldsymbol{z}_{i}^{\perp} \in \mathbb{R}^{n-1}} \int_{\beta_{i}^{U}}^{\infty} \frac{1}{\sum_{k=1}^{n_{c}} I_{F_{k}} \left(\boldsymbol{R}_{i} \boldsymbol{z}_{i}^{\perp} + \boldsymbol{\alpha}_{i} \boldsymbol{z}_{i}^{\parallel}\right)} f_{\boldsymbol{Z}_{i}^{\parallel}} \left(\boldsymbol{z}_{i}^{\parallel}\right) f_{\boldsymbol{Z}_{i}^{\perp}} \left(\boldsymbol{z}_{i}^{\perp}\right) d\boldsymbol{z}_{i}^{\parallel} d\boldsymbol{z}_{i}^{\perp}, \ i = 1, \dots, n_{c}$$
 (12)

The above equation provides an expression for calculating p_i within the framework of Line Sampling. As integration along the parallel direction is carried out taking into account interactions of 241 the failure events associated with the different components, eq. (12) is denoted as a multidomain 242 Line Sampling (mLS) expression for calculating p_i . 243 Direct calculation of the expression for the failure probability as proposed in eq. (9) demands calculating each individual quantity p_i , $i = 1, ..., n_c$ by means of mLS. As it is expected that 245 n_c can be in the order of hundreds or even thousands, calculating each term in the summation 246 can become extremely demanding. As an alternative, this summation can be estimated by means 247 of simulation, following the approach proposed in [30]. For that purpose, consider the following 248 variant of eq. (9):

$$p_F = \sum_{i=1}^{n_c} \left(\frac{1}{\omega_i} p_i\right) \omega_i \tag{13}$$

where ω_i , $i = 1, ..., n_c$ is a weight factor such that $\omega_i > 0$ and $\sum_{i=1}^{n_c} \omega_i = 1$. This set of weights can be interpreted as the probability mass function of a discrete random variable. A possible criterion for selecting the *i*-th weight is to set it proportional to the failure probability associated with the *i*-th component, as considered in Importance Sampling using design points [42]. This leads to the following expression for calculating the weights.

$$\omega_i = \frac{p_{F,i}}{\sum_{k=1}^{n_c} p_{F,k}}, \ i = 1, \dots, n_c$$
 (14)

Thus, eq. (13) becomes an expression that involves summation over a discrete random variable as well as integration over a number of continuous random variables. Within the context of simulation, p_F is estimated by generating samples of the discrete and continuous random variables, that is:

$$p_F \approx \tilde{p}_F = \frac{1}{N} \sum_{j=1}^{N} \left(\frac{1}{\omega_{i^{(j)}}} p_{i^{(j)}} \left(\boldsymbol{z}_{i^{(j)}}^{\perp,(j)} \right) \right)$$
 (15)

where \tilde{p}_F is an estimate of p_F ; N denotes the total number of samples; $i^{(j)}$, $j=1,\ldots,N$ are independent and identically distributed samples drawn with replacement from the set $I=\{1,2,\ldots,n_c\}$ with probability mass function ω_i , $i=1,\ldots,n_c$; $\boldsymbol{z}_{i^{(j)}}^{\perp,(j)}$, $j=1,\ldots,N$ are independent and identically distributed samples that follow $f_{\boldsymbol{Z}_{i^{(j)}}^{\perp}}\left(\boldsymbol{z}_{i^{(j)}}^{\perp}\right)$; and where $p_{i^{(j)}}\left(\boldsymbol{z}_{i^{(j)}}^{\perp,(j)}\right)$, $j=1,\ldots,N$ represents an estimate of quantity $p_{i^{(j)}}$ evaluated at the sample $\boldsymbol{z}_{i^{(j)}}^{\perp,(j)}$, that is:

$$p_{i(j)}\left(\boldsymbol{z}_{i(j)}^{\perp,(j)}\right) = \int_{-\infty}^{-\beta_{i(j)}^{L}} \frac{1}{\sum_{k=1}^{n_{c}} I_{F_{k}}\left(\boldsymbol{R}_{i(j)}\boldsymbol{z}_{i(j)}^{\perp,(j)} + \boldsymbol{\alpha}_{i(j)}\boldsymbol{z}_{i(j)}^{\parallel}\right)} f_{Z_{i(j)}^{\parallel}}\left(\boldsymbol{z}_{i(j)}^{\parallel}\right) d\boldsymbol{z}_{i(j)}^{\parallel} + \int_{\beta_{i(j)}^{U}}^{\infty} \frac{1}{\sum_{k=1}^{n_{c}} I_{F_{k}}\left(\boldsymbol{R}_{i(j)}\boldsymbol{z}_{i(j)}^{\perp,(j)} + \boldsymbol{\alpha}_{i(j)}\boldsymbol{z}_{i(j)}^{\parallel}\right)} f_{Z_{i(j)}^{\parallel}}\left(\boldsymbol{z}_{i(j)}^{\parallel}\right) d\boldsymbol{z}_{i(j)}^{\parallel}, \ j = 1, \dots, N$$
 (16)

It is seen that the last equation corresponds to an estimate of the quantity p_i calculated by means of mLS. It represents the integral over the line that passes through the sample $\mathbf{z}_{i^{(j)}}^{\perp,(j)}$ and which is parallel to $\alpha_{i^{(j)}}$ and whose argument is the standard Gaussian univariate probability distribution divided over the number of components that fail at a given point of that line. Thus, eq. (16) can be interpreted as a means of exploring the interactions that occur between the behavior of different components along the line. Details about the numerical evaluation of eq. (16) are discussed in Section 3.3.2.

271 It is straightforward to demonstrate that the coefficient of variation of the probability estimate of

eq. (15) (which is denoted as δ_{p_F}) is equal to:

$$\delta_{p_F} = \frac{1}{\tilde{p}_F} \sqrt{\frac{1}{N(N-1)} \sum_{j=1}^{N} \left(\left(\frac{1}{\omega_{i^{(j)}}} p_{i^{(j)}} \left(\boldsymbol{z}_{i^{(j)}}^{\perp,(j)} \right) \right) - \tilde{p}_F \right)^2}$$
 (17)

273 3.3. Practical Implementation

Practical implementation of eq. (16) demands solving two issues: the generation of samples $z_{i^{(j)}}^{\perp,(j)}$, $j=1,\ldots,N$ and the calculation of the line integral associated with mLS. These two issues are discussed in the following.

277 3.3.1. Generation of samples $oldsymbol{z}_{i^{(j)}}^{\perp,(j)}$

Regarding the first implementation issue, recall that $\mathbf{z}_{i^{(j)}}^{\perp,(j)}$, $j=1,\ldots,N$ are independent and identically distributed samples that follow $f_{\mathbf{Z}_{i^{(j)}}^{\perp}}\left(\mathbf{z}_{i^{(j)}}^{\perp}\right)$. These samples can be conveniently generated by means of the following algorithm.

- 1. Set j = 1.
- 282 2. Draw an element from the set $I = \{1, 2, ..., n_c\}$ with probability ω_i , $i = 1, ..., n_c$. The drawn element is denoted as $i^{(j)}$.
- 3. Generate a random sample $z^{(j)}$ following a *n*-dimensional standard Gaussian distribution.
- 285 4. Calculate [33, 34]:

$$\boldsymbol{R}_{i(j)}\boldsymbol{z}_{i(j)}^{\perp,(j)} = \boldsymbol{z}^{(j)} - \left(\boldsymbol{\alpha}_{i(j)}^T \boldsymbol{z}^{(j)}\right) \boldsymbol{\alpha}_{i(j)}$$
(18)

5. In case j = N, stop the algorithm. Otherwise, return to step 2 with j = j + 1.

The core of the algorithm described above lies in step 4, which is represented schematically in Figure 5, where it has been assumed for simplicity that $n = n_c = 2$ and that $i^{(j)} = 1$. As noted from Figure 5, eq. (18) consists of subtracting the projection of the random sample $\mathbf{z}^{(j)}$ over the important direction $\boldsymbol{\alpha}_{i^{(j)}}$ (that is, $(\boldsymbol{\alpha}_{i^{(j)}}^T \mathbf{z}^{(j)}) \boldsymbol{\alpha}_{i^{(j)}}$) from the random sample $\mathbf{z}^{(j)}$ itself [33, 34]. It should be noted that such step does not produce $\mathbf{z}_{i^{(j)}}^{\perp,(j)}$ but instead, it leads to $\mathbf{R}_{i^{(j)}} \mathbf{z}_{i^{(j)}}^{\perp,(j)}$. This is quite convenient from a numerical viewpoint, as all calculations associated with mLS demand knowledge of $\mathbf{R}_{i^{(j)}} \mathbf{z}_{i^{(j)}}^{\perp,(j)}$ (and not of $\mathbf{z}_{i^{(j)}}$). Hence, explicit calculation of the rotation matrices \mathbf{R}_{i} , $i = 1, \ldots, n_c$ is avoided.

295 3.3.2. Evaluation of integral along line

Eq. (16) corresponds to a one-dimensional integral along the line $l^{(j)}$ that passes through

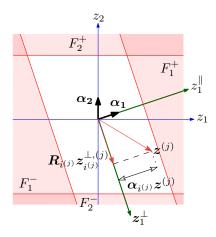


Figure 5: Schematic representation of generation of samples for multidomain Line Sampling $(n = n_c = 2, i^{(j)} = 1)$.

the sample $m{z}_{i^{(j)}}^{\perp,(j)}$ and which is parallel to $m{lpha}_{i^{(j)}}$. The argument of this integral is the standard Gaussian probability density function as a function of the coordinate $z_{i^{(j)}}^{\parallel}$ divided by the number of components that fail at realization $R_{i^{(j)}} z_{i^{(j)}}^{\perp,(j)} + \alpha_{i^{(j)}} z_{i^{(j)}}^{\parallel}$; the latter number is given by the formula $\sum_{k=1}^{n_c} I_{F_k} \left(\boldsymbol{R}_{i^{(j)}} \boldsymbol{z}_{i^{(j)}}^{\perp,(j)} + \boldsymbol{\alpha}_{i^{(j)}} \boldsymbol{z}_{i^{(j)}}^{\parallel} \right). \text{ As } \boldsymbol{R}_{i^{(j)}} \boldsymbol{z}_{i^{(j)}}^{\perp,(j)} \text{ is a fixed vector for a given line (see eq. (18)), the}$ challenge for calculating the line integral in eq. (16) lies precisely in calculating the number of failed 301 components as a function of $z_{i^{(j)}}^{\parallel}$. For a better understanding of this issue, consider the schematic 302 representation in Figure 6, that depicts a particular case where $n = n_c = 2$, N = 1 and $i^{(1)} = 2$. In 303 particular, Figure 6(a) illustrates the problem in the space of standard Gaussian random variables, where $l^{(1)}$ denotes the line that passes through the sample $\boldsymbol{z}_2^{\perp,(1)}$ and which is parallel to $\boldsymbol{\alpha}_2$. It is seen that line $l^{(1)}$ intersects the different failure domains associated with each of the two 306 components considered. Such issue must be considered when solving the integral associated with 307 that line, as this affects the indicator functions of each component $I_{F_k}\left(\mathbf{R}_2\mathbf{z}_2^{\perp,(1)}+\boldsymbol{\alpha}_2z_2^{\parallel}\right),\ k=1,2,$ 308 as depicted schematically in Figures 6(b) and 6(c), respectively, as well as the compound indicator function $\sum_{k=1}^{2} I_{F_k} \left(\mathbf{R}_2 \mathbf{z}_2^{\perp,(1)} + \boldsymbol{\alpha}_2 z_2^{\parallel} \right)$, as seen in Figure 6(d). A close examination of the indicator 310 functions associated with the components as shown in Figures 6(b) and 6(c) reveals that they can 311 be represented as the summation of two unit step functions. In a general case, the indicator 312 function associated with component k is cast as:

$$I_{F_k}\left(\boldsymbol{R}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\perp,(j)} + \boldsymbol{\alpha}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\parallel}\right) = u\left(-g_{2k-1}\left(\boldsymbol{R}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\perp,(j)} + \boldsymbol{\alpha}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\parallel}\right)\right) + u\left(-g_{2k}\left(\boldsymbol{R}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\perp,(j)} + \boldsymbol{\alpha}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\parallel}\right)\right), \ k = 1, \dots, n_c, \ j = 1, \dots, N$$
 (19)

where $u(\cdot)$ denotes the unit step function. Replacing the expressions for the performance functions

 $g_{2k-1}(z)$ and $g_{2k}(z)$ (see eqs. (2) and (3), respectively) into the above equation leads to the following expression.

$$I_{F_{k}}\left(\boldsymbol{R}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\perp,(j)} + \boldsymbol{\alpha}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\parallel}\right) = u\left(-\beta_{k}^{L} - \boldsymbol{\alpha}_{k}^{T}\boldsymbol{R}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\perp,(j)} - \boldsymbol{\alpha}_{k}^{T}\boldsymbol{\alpha}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\parallel}\right) + u\left(-\beta_{k}^{U} + \boldsymbol{\alpha}_{k}^{T}\boldsymbol{R}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\perp,(j)} + \boldsymbol{\alpha}_{k}^{T}\boldsymbol{\alpha}_{i^{(j)}}\boldsymbol{z}_{i^{(j)}}^{\parallel}\right), \ k = 1, \dots, n_{c}, \ j = 1, \dots, N$$

$$(20)$$

Under the assumption that $\alpha_k^T \alpha_{i^{(j)}} \neq 0$, it is possible to express the last equation in a more compact format, that is:

$$I_{F_k}\left(\mathbf{R}_{i^{(j)}}\mathbf{z}_{i^{(j)}}^{\perp,(j)} + \boldsymbol{\alpha}_{i^{(j)}}\mathbf{z}_{i^{(j)}}^{\parallel}\right) = u\left(\xi_{2k-1}^{(j)}\left(\mathbf{z}_{i^{(j)}}^{\parallel} - c_{2k-1}^{(j)}\right)\right) + u\left(\xi_{2k}^{(j)}\left(\mathbf{z}_{i^{(j)}}^{\parallel} - c_{2k}^{(j)}\right)\right),$$

$$k = 1, \dots, n_c, \ j = 1, \dots, N$$
(21)

where:

$$c_{2k-1}^{(j)} = -\frac{\beta_k^L + \boldsymbol{\alpha}_k^T \boldsymbol{R}_{i(j)} \boldsymbol{z}_{i(j)}^{\perp,(j)}}{\boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i(j)}}$$
(22)

$$c_{2k}^{(j)} = \frac{\beta_k^U - \boldsymbol{\alpha}_k^T \boldsymbol{R}_{i^{(j)}} \boldsymbol{z}_{i^{(j)}}^{\perp,(j)}}{\boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}}}$$
(23)

$$\xi_{2k-1}^{(j)} = -\operatorname{sgn}\left(\boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}}\right) \tag{24}$$

$$\xi_{2k}^{(j)} = \operatorname{sgn}\left(\boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}}\right) \tag{25}$$

and where $\operatorname{sgn}(\cdot)$ represents the sign function. As noted from eq. (21), $c_{2k-1}^{(j)}$ and $c_{2k}^{(j)}$ denote the coordinate $z_{i(j)}^{\parallel}$ for which the corresponding unit step function changes its value. Such concept is represented schematically in Figures 6(b) and 6(c). In fact, $c_{2k-1}^{(j)}$ and $c_{2k}^{(j)}$ denote the Euclidean distances from the sample $\boldsymbol{z}_{i(j)}^{\perp,(j)}$ to the limit state functions $g_{2k-1}(\boldsymbol{z}) = 0$ and $g_{2k}(\boldsymbol{z}) = 0$, respectively, measured along the line $l^{(j)}$. Additionally, $\xi_{2k-1}^{(j)}$ and $\xi_{2k}^{(j)}$ are variables whose value is either -1 or 1 depending on the sign of the dot product $\boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}}$. It should be recalled that eqs. (22) to (25) were deduced under the assumption that $\boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}} \neq 0$. However, these expressions can be generalized for the case where $\boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}} = 0$, as shown in detail in Appendix A.

The characterization of the indicator function associated with the k-th component as shown in eq. (21) allows a straightforward estimation of the sought line integral. For that purpose,

let $\left\{q_1^{(j)}, q_2^{(j)}, \dots, q_{2n_c}^{(j)}\right\}$ denote the sequence of integers such that $c_{q_1^{(j)}}^{(j)} \leq c_{q_2^{(j)}}^{(j)} \leq \dots \leq c_{q_{2n_c}^{(j)}}^{(j)}$; in

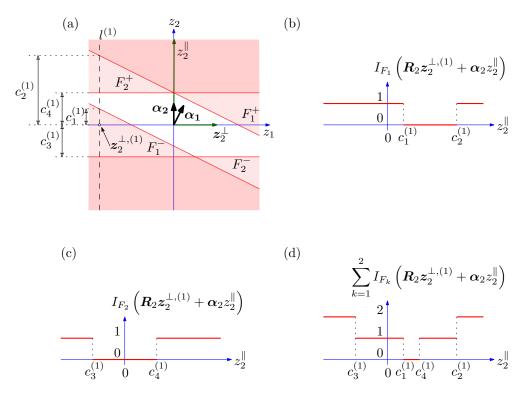


Figure 6: Schematic representation of line associated with the application of multidomain Line Sampling $(n = n_c = 2)$. (a) Representation in standard Gaussian space. (b) Behavior of indicator function associated with the first component $I_{F_1}(\cdot)$ with respect to z_2^{\parallel} along line $l^{(1)}$. (c) Behavior of indicator function associated with the second component $I_{F_2}(\cdot)$ with respect to z_2^{\parallel} along line $l^{(1)}$.(d) Behavior of the compound indicator function with respect to z_2^{\parallel} along line $l^{(1)}$.

addition, let $c_{q_0^{(j)}}^{(j)} \to -\infty$ and $c_{q_{2n_c+1}}^{(j)} \to \infty$. Furthermore, let $m_s^{(j)}$ be a natural number (including 0), which is defined as:

$$m_s^{(j)} = m_0^{(j)} + \sum_{l=1}^s \xi_{q_s^{(j)}}^{(j)}, \ s = 0, \dots, 2n_c, \ j = 1, \dots, N$$
 (26)

where $m_0^{(j)}$ counts the number of times that $\xi_q^{(j)} = -1, q = 1, \ldots, 2n_c$. The role of $m_s^{(j)}$, $s = 0, \ldots, 2n_c$ is expressing the number of components that fail at a point that belongs to the line $l^{(j)}$ and whose coordinate $z_{i^{(j)}}^{\parallel}$ lies within the interval $\left(c_{i^{(j)}}^{(j)}, c_{i^{(j)}}^{(j)}\right)$. In other words, $m_s^{(j)}$, $s = 0, \ldots, 2n_c$ contains all the values that the function $\sum_{k=1}^{n_c} I_{F_k} \left(\mathbf{R}_{i^{(j)}} \mathbf{z}_{i^{(j)}}^{\perp,(j)} + \boldsymbol{\alpha}_{i^{(j)}} \mathbf{z}_{i^{(j)}}^{\parallel}\right)$ assumes along the line $l^{(j)}$. For a better understanding of this point, consider Figure 6(d) that illustrates the function $\sum_{k=1}^{n_c} I_{F_k} \left(\mathbf{R}_{i^{(j)}} \mathbf{z}_{i^{(j)}}^{\perp,(j)} + \boldsymbol{\alpha}_{i^{(j)}} \mathbf{z}_{i^{(j)}}^{\parallel}\right)$ associated with line $l^{(1)}$ of Figure 6(a). It is seen that this function presents a staircase pattern, as depending on the value of z_2^{\parallel} , the number of failed components varies between 0, 1 and 2. This staircase pattern is reproduced by the quantity $m_s^{(j)}$ in eq. (26), as shown in Table 1. Note that for preparing this Table and according to the qualitative

information in Figures 6(b) and 6(c), it is considered that $c_3^{(1)} < c_1^{(1)} < c_4^{(1)} < c_2^{(1)}, \, \xi_1^{(1)} = \xi_3^{(1)} = -1$ and $\xi_2^{(1)} = \xi_4^{(1)} = 1$.

s	$q_s^{(1)}$	$\xi_{q_s^{(1)}}^{(1)}$	$\left(c_{q_s^{(1)}}^{(1)}, c_{q_{s+1}^{(1)}}^{(1)}\right)$	$m_s^{(1)}$
0	-	-	$\left(-\infty, c_3^{(1)}\right)$	-2
1	3	-1	$(c_3^{(1)}, c_1^{(1)})$	-1
2	1	-1	$\left(c_1^{(1)}, c_4^{(1)}\right)$	0
3	4	1	$\left(c_4^{(1)}, c_2^{(1)}\right)$	1
4	2	1	$\left(c_2^{(1)},\infty\right)$	2

Table 1: Values that variable $m_s^{(1)}$ assumes along line $l^{(1)}$ associated with the schematic illustration in Figure 6.

Taking into account the above definitions, eq. (16) is calculated in closed form as shown below:

$$p_{i^{(j)}}\left(\boldsymbol{z}_{i^{(j)}}^{\perp,(j)}\right) = \sum_{s=0}^{s^{L}-1} \frac{\Phi\left(c_{q_{s+1}}^{(j)}\right) - \Phi\left(c_{q_{s}}^{(j)}\right)}{m_{s}^{(j)}} + \sum_{s=s^{U}}^{2n_{c}} \frac{\Phi\left(c_{q_{s+1}}^{(j)}\right) - \Phi\left(c_{q_{s}}^{(j)}\right)}{m_{s}^{(j)}}$$
(27)

where s^L is an integer such that $c_{q_{sL}^{(j)}}^{(j)} = -\beta_{i^{(j)}}^L$ and s^U is another integer such that $c_{q_{sU}^{(j)}}^{(j)} = \beta_{i^{(j)}}^U$.

It is noted that from a numerical viewpoint, the computation of eq. (27) demands evaluating the response of each of the components twice, as it is necessary to solve two dot products involving the unit vector $\boldsymbol{\alpha}_l$ (see eqs. (22) to (25)). Therefore, assessing the estimator \tilde{p}_F in eq. (15) demands 2N evaluations of each component response.

350 3.4. Summary

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- The application of multidomain Line Sampling for calculating the failure probability of a series system as considered in this contribution involves the following steps.
- 1. Define the basic information of the problem. This implies setting up the probabilistic characterization of the parameter vector \boldsymbol{x} (of dimension n) in terms of its mean $\boldsymbol{\mu}$ and (positive definite) covariance matrix \boldsymbol{C} . Additionally, define vector $\boldsymbol{a}_i, i = 1, \dots, n_c$ that characterizes the response of the i-th component as well as the allowable lower and upper bounds for the response (b_i^L and b_i^U , respectively).
- 2. Set up the performance functions in standard normal space by applying eqs. (1), (2) and (3). Calculate the reliability indexes β_i^L and β_i^U , $i = 1, ..., n_c$ as well as the unit vectors $\alpha_i, i = 1, ..., n_c$.

- 3. Calculate the weights ω_i , $i = 1, ..., n_c$ by means of eq. (14).
- 4. Sample (with replacement) a total of N integers $i^{(j)}$, j = 1, ..., N from the set $I = \{1, 2, ..., n_c\}$ with probability ω_i , $i = 1, ..., n_c$. Generate samples $\boldsymbol{z}_{i^{(j)}}^{\perp, (j)}$, j = 1, ..., N applying the procedure described in Section 3.3.1 (see eq. (18)).
- 5. Estimate $p_{i^{(j)}}\left(\boldsymbol{z}_{i^{(j)}}^{\perp,(j)}\right), \ j=1,\ldots,N$ by means of eq. (27).
- 6. Calculate the estimator of the failure probability as well as its coefficient of variation applying eqs. (15) and (17).

368 4. Examples

369 4.1. Test Example 1

This first test example is borrowed from [15, 43]. It comprises the calculation of the failure probability of a series system, where the response of its i-th component is defined as:

$$r_i(\boldsymbol{x}) = x_i, \ i = 1, \dots, n_c \tag{28}$$

where x_i is a realization of a Gaussian random variable with zero mean, unit standard deviation and pairwise correlation coefficient 0.5 with all random variables (other than itself). That is, the correlation matrix \mathbf{R} of dimension $n_c \times n_c$ is defined as:

$$\mathbf{R} = \begin{bmatrix} 1 & 0.5 & \dots & 0.5 \\ 0.5 & 1 & \dots & 0.5 \\ \vdots & \vdots & \ddots & \vdots \\ 0.5 & 0.5 & \dots & 1 \end{bmatrix}$$
 (29)

Note that there are $n = n_c$ random variables. The threshold levels associated with the performance of each component are set such that $b_i^L \to -\infty$, $i = 1, ..., n_c$ and $b_i^U = \beta$, $i = 1, ..., n_c$ (where β is a real number), respectively. The failure probability associated with this series system can be expressed in terms of the following one-dimensional integral [15, 43], which can be accurately calculated by means of an appropriate quadrature.

$$p_F = \int_{-\infty}^{\infty} \left(1 - \left(1 - \Phi \left(\frac{-\beta - \sqrt{0.5}z}{\sqrt{1 - 0.5}} \right) \right)^{n_c} \right) f_Z(z) dz \tag{30}$$

From the above equation, recall that $f_Z(\cdot)$ and $\Phi(\cdot)$ represent the probability density function and cumulative density function of a standard Gaussian random variable, respectively.

The problem described above is solved by means of both the above integral and multidomain Line Sampling. Different combinations of the number of components n_c (10¹, 10², 10³, 10⁴) and of the threshold β (3, 4, 5) are investigated. For all these combinations, multidomain Line Sampling is implemented considering N = 200 lines (hence, $2 \times 200 = 400$ system analyses are performed). The results obtained for the failure probability estimates as well as their coefficient of variation are shown in Figures 7 and 8, respectively.

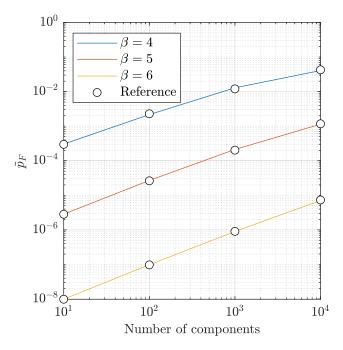


Figure 7: Example 1 - failure probability with respect to the number of components n_c and threshold level β . Solid line: estimates with multidomain Line Sampling. Circle: reference result.

Figure 7 illustrates the estimates of the failure probability generated with multidomain Line Sampling with solid line. In addition, the reference results provided by solving numerically the integral in eq. (30) are shown with circles. It is seen that there is an excellent agreement between the results, irrespective of the number of components and the threshold level. This is quite remarkable, as the failure probabilities involved in the figure span about six orders of magnitude. Figure 8 shows the coefficient of variation associated with the estimates of the failure probability associated with multidomain Line Sampling. It is observed that all coefficients of variation are relatively low, which is quite desirable from a practical viewpoint. Furthermore, it is observed that the coefficients of variation are quite small for small values of failure probabilities. Such

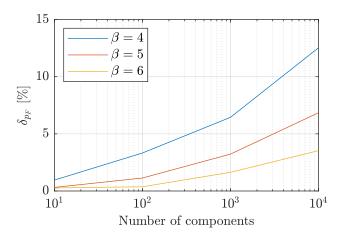


Figure 8: Example 1 - coefficient of variation of the failure probability with respect to the number of components n_c and threshold level β .

behavior is similar to the one observed in [29] and is explained by the fact that for small failure probabilities, interactions between components in the failure region become less relevant.

As a summary of this test example, it is observed that multidomain Line Sampling allows coping with a large number of random variables and components for estimating small failure probabilities with high precision and a reduced number of samples.

o2 4.2. Test Example 2: Truss Structure

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This test example involves a statically determined truss structure subject to three point loads P_k , k = 1, ..., 3, as depicted in Figure 9. These three point loads are modeled as Gaussian random variables with expected value 10 [kN], standard deviation 1 [kN] and constant correlation between them equal to 0.5. The maximum axial load that can be supported by the bars of the truss is as follows: bars 1, 4, 5, 6, 7, 8, 9 support a maximum load of 19 [kN]; bars 2 and 3 support a maximum load of 25 [kN]; bars 10 and 13 support a maximum load of 27 [kN]; and bars 11 and 12 support a maximum load of 10 [kN]. For simplicity, it is assumed that the bars are capable to support this maximum load either in tension or in compression.

The objective is determining the probability that the maximum allowable axial load due to the external loading is exceeded in one or more bars of the truss. As the truss possesses 13 bars and failure of any of these bars leads to failure of the system, this can be interpreted a series system with $n_c = 13$ components. The response of each component is its axial load and the allowable threshold is given by the maximum load supported by each bar.

The probability of failure associated with the series event is estimated by means of multidomain Line Sampling (mLS), considering a total of $N = 5 \times 10^5$ samples. As each sample comprises

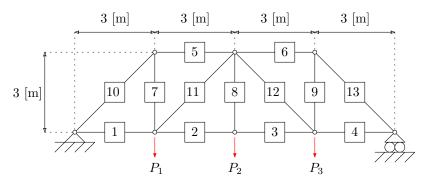


Figure 9: Example 2 - truss structure.

a total of two evaluations of the response of the components, a total of 10⁶ system analyses are
carried out. Note that this large number of samples for mLS is considered in order to carry
out comparisons with Monte Carlo simulation, which provides reference results. In this case,
Monte Carlo simulation is applied considering a total of 10⁶ samples. The results obtained for the
estimates of the failure probability and its coefficient of variation are shown in figures 10 and 11,
respectively.

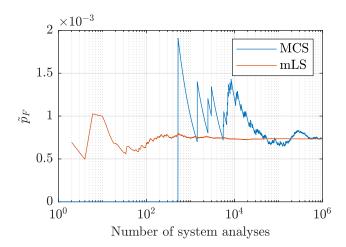


Figure 10: Example 2 - evolution of estimate of failure probability with respect to the number of system analyses (MCS: Monte Carlo simulation, mLS: multidomain Line Sampling).

An examination of Figure 10 indicates that both multidomain Line Sampling and Monte Carlo simulation provide similar estimates of the failure probability for a large number of system analyses. However, the estimator produced with multidomain Line Sampling stabilizes extremely fast:
in fact, with about only 100 system analyses, it provides an excellent estimator of the failure probability. This is quite remarkable, considering that the failure probability is relatively small, that is, about $p_F \approx 7 \times 10^{-4}$.

The evolution of the coefficient of variation as shown in Figure 11 reinforces the conclusions

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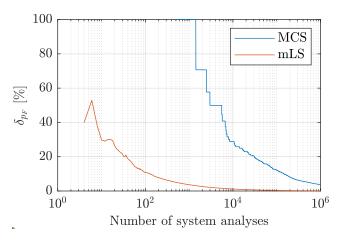


Figure 11: Example 2 - evolution of estimate of the coefficient of variation of the failure probability with respect to the number of system analyses (MCS: Monte Carlo simulation, mLS: multidomain Line Sampling).

drawn from Figure 10. It is seen that the coefficient of variation of the probability estimate produced with mLS is about 10% with only 100 system analyses. In order to produce an estimate with comparable coefficient of variation, Monte Carlo simulation demands 10⁵ system analyses. Such result highlights the benefits of mLS for estimating failure probabilities associated with a series event.

4.3. Application Example 3: Six-Story Building Subject to Stochastic Gaussian Ground Acceleration

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This application example involves a six-story reinforced concrete building subject to a stochastic Gaussian ground acceleration, as illustrated in Figure 12. The objective is estimating the first excursion probability that the interstory drifts of the building exceed a prescribed threshold within the duration of the acceleration. It is assumed that the building experiences small vibrations and hence, its behavior can be modeled as linear elastic. In fact, as discussed in the sequence, this problem can be modeled as a series system with a large number of components arising due to the discretization of time.

The assumption of linear elastic behavior of the building is appropriate for analyzing serviceability conditions, see e.g. [44, 45, 46], and allows conducting a reliability analysis by means of multidomain Line Sampling. For those cases where the assumption of a linear elastic behavior does not hold (e.g. progressive collapse or collapse), other more general methods should be applied, see e.g. [47, 48, 49, 50, 51].

Each floor of the building is composed of a square slab of side 32 [m] and thickness 0.2 [m], and is supported by 16 columns of square cross section of 0.4 [cm] and a shear wall of 0.2 [m] thickness.

The building is modeled as linear elastic and classically damped, with Young's modulus is equal to 2.3×10^{10} [Pa]. It is assumed that the building experiences small displacements and hence, its elements remain within the linear elastic range. The finite element model, which is taken from [52], involves about 9500 shell and beam elements and more than 50×10^3 degrees-of-freedom. Classical damping of 5% is considered for all modes retained in the analysis.

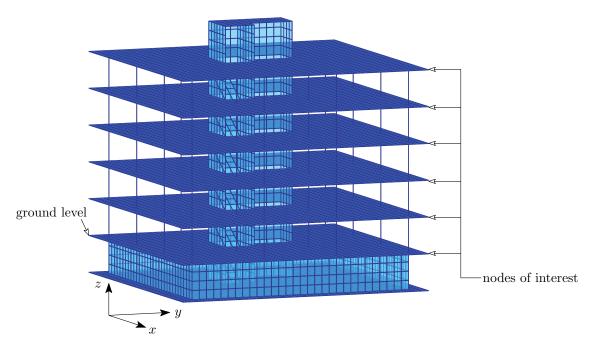


Figure 12: Example 3 - building model.

The building is excited by a stochastic Gaussian ground acceleration along the y direction (see 457 Figure 12), which is modeled by means of the Clough-Penzien power spectrum (see, e.g. [53, 54, 458 55]), and which is in turn modulated by the Shinozuka-Sato envelope function [56]. The stochastic 459 ground acceleration possesses a duration of 20 [s], with discrete representation considering a time 460 step of 0.01 [s]. The associated discrete white noise process possesses spectral intensity of $3 \times$ 461 10^{-4} [m²/s³] and the properties of the primary and secondary Clough-Penzien filters are circular 462 frequency $\omega_1 = 4\pi$ [rad/s] and $\omega_2 = 0.4\pi$ [rad/s] and damping ratios $\zeta_1 = \zeta_2 = 0.7$, respectively. 463 The parameters for the Shinozuka-Sato envelope are selected as $c_1=0.14$ and $c_2=0.16$. The 464 stochastic ground acceleration is represented with the help of the Karhunen-Loève expansion 465 considering 1466 terms (see, e.g. [57]). For additional details on the representation of the stochastic 466 ground acceleration model, it is referred to [53, 54, 55, 56]. 467 For design purposes, the interstory drifts along the y direction should not exceed a threshold level

of 2×10^{-3} times the story height within the duration of the stochastic loading. This condition

is verified considering the six nodes indicated in Figure 12, which implies that a total of five interstory drifts must be controlled. Appendix B provides a brief description on the procedure 471 for calculating these interstory drifts. The chances that any of these interstory drifts exceed its prescribed threshold within the duration of the stochastic ground acceleration can be interpreted 473 as the probability of failure of a series event, see e.g. [29]. In this case, the components are each 474 of the interstory drifts responses at each time instant. As there is a total of five interstory drifts 475 and 2001 discrete time instants, the total number of components is $n_c = 10005$. On the other hand, the total number of random variables involved in the problem is n = 1466, which is equal to the number of terms associated with the Karhunen-Loève expansion. Hence, the problem under 478 consideration corresponds to a case with a large number of random variables (over 10³ of them) 479 and a large number of components (over 10^4 components). 480

The probability of failure associated with the system event is calculated by means of multidomain Line Sampling, considering a total of $N = 5 \times 10^3$ samples. This implies that the system's response is calculated a total of 10^4 times. That is, 10^4 dynamic analyses are carried out. In addition and in order to provide a basis for comparison, the failure probability is also assessed by means of Efficient Importance Sampling [29] and Directional Importance Sampling [31, 32], which are simulation techniques specially developed for calculating failure probabilities of linear structural systems subject to Gaussian excitation. These two simulation techniques are implemented considering a total of 10^4 samples (which imply performing a total of 10^4 dynamic analyses). The results obtained are shown in Figures 13 and 14 as well as in Table 2.

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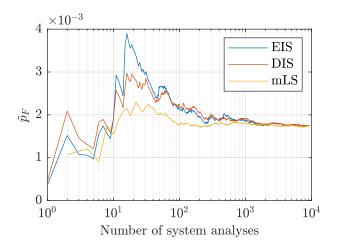


Figure 13: Example 3 - evolution of estimate of failure probability with respect to the number of system analyses (EIS: Efficient Importance Sampling, DIS: Directional importance Sampling, mLS: multidomain Line Sampling).

As noted from Figure 13, the three simulation techniques under consideration can produce good

estimates of the failure probability with a reduced number of samples. However, it is seen that
the estimator associated with multidomain Line Sampling stabilizes quicker than the estimators
associated with the other two simulation techniques. In fact, with about N = 50 samples (that is,
100 system analyses), multidomain Line Sampling already provides an excellent estimate of the
failure probability. This is confirmed by examining the results in table 2, where it is seen that the
associated coefficient of variation is already below 10%.

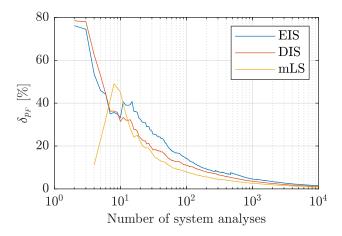


Figure 14: Example 3 - evolution of estimate of the coefficient of variation of the failure probability with respect to the number of system analyses (EIS: Efficient Importance Sampling, DIS: Directional importance Sampling, mLS: multidomain Line Sampling).

The results presented in Figure 14 support the observations already drawn from Figure 13. That is, the coefficient of variation associated with the probability estimates of all simulation techniques decreases quickly with the number of system analysis. Furthermore, the estimate associated with multidomain Line Sampling is the one presenting the smallest coefficient of variation (excluding the region of about 10 system analyses).

No. system analyses	$ ilde{p}_F^{ m EIS}$	$\delta_{p_F^{ m EIS}}$ [%]	$ ilde{p}_F^{ m DIS}$	$\delta_{p_F^{ m DIS}}$ [%]	$ ilde{p}_F^{ ext{mLS}}$	$\delta_{p_F^{ m mLS}}$ [%]
10^{2}	2.2×10^{-3}	14.2%	2.2×10^{-3}	11.1%	1.8×10^{-3}	7.9%
10^{3}	1.9×10^{-3}	4.6%	1.9×10^{-3}	3.6%	1.8×10^{-3}	2.6%
10^{4}	1.7×10^{-3}	1.5%	1.7×10^{-3}	1.1%	1.7×10^{-3}	0.9%

Table 2: Example 3 - Estimates of the failure probability \tilde{p}_F and its coefficient of variation δ_{p_F} (EIS: Efficient Importance Sampling, DIS: Directional importance Sampling, mLS: multidomain Line Sampling).

Table 2 reports the probability estimates and their coefficient of variation for the three simulation techniques under consideration for different number of system analyses. It is seen that all three techniques are quite successful in estimating the failure probability, as there is good agreement between the different estimates, with relatively low coefficient of variation. However, for

all cases reported in the table, multidomain Line Sampling provides probability estimates which are closer to the reference solution (with 10⁴ system analyses) and with the smallest coefficient of variation.

The (small) differences in performance between the three simulation techniques analyzed in this 509 example as presented in Table 2 can be understood as follows. Efficient Importance Sampling 510 (EIS) is a specially-designed variant of Importance Sampling that estimates the failure proba-511 bility by generating samples (realizations of \mathbf{Z}) exclusively in the failure domain. Directional 512 Importance Sampling (DIS) operates in a similar way as Efficient Importance Sampling, but it explores directions instead of samples; in other words, it explores an infinite number of samples 514 along a given ray starting at the origin of the standard normal space. Multidomain Line Sampling 515 (mLS) shares some common aspects with Directional Importance Sampling in the sense that an 516 infinite number of samples is explored. However, these samples fall in a line whose orientation is 517 different from the aforementioned ray. 518

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As a further comparison between the performance of the three simulation techniques discussed above, Table 3 presents both the number of system analyses and the relative execution time for attaining an estimate of the failure probability with coefficient of variation $\delta_{p_F} = 10\%$. It is observed that the smallest number of system analyses and relative execution time are associated with mLS. The other two simulation approaches, that is EIS and DIS, demand more samples to attain the prescribed coefficient of variation and more relative execution time than those associated with mLS. An an additional observation from Table 3, it should be noted that the relation between number of system analyses and relative execution time for the different simulation techniques is not proportional, as the specific implementation steps of EIS [29], DIS [31, 32] and mLS differ between them.

Simulation	No. system	Relative
technique	analyses	Execution time
EIS	171	249%
DIS	130	234%
mLS	62	100%

Table 3: Example 3 - Number of system analyses and relative execution time for achieving probability estimate with coefficient of variation of 10% (EIS: Efficient Importance Sampling, DIS: Directional importance Sampling, mLS: multidomain Line Sampling).

5. Conclusions and Outlook

This contribution has presented an approach for estimating the probability of occurrence of 530 a series system event by means of multidomain Line Sampling. In particular, multidomain Line 531 Sampling is applied in order to determine the effective contribution of a single component to the 532 overall failure probability. In this context, effective means that proper consideration is given to 533 the failure event associated with a particular component and its interaction with other components. The overall failure probability is then determined by randomly sampling among different 535 components. The examples presented in this contribution suggest that multidomain Line Sam-536 pling is applicable to problems involving both a small and a large number of random variables 537 and components, respectively. Much of the success of the multidomain Line Sampling strategy as reported herein can be attributed to the way each failure domain associated with an individual component is examined. By exploring lines, one can analyze an infinite number of realizations instead of a single one. In this way, each line provides a considerable amount of information. Moreover, in view of the linearity of the response with respect to the unknown random parameters, it is possible to solve the integral associated with that line by means of a closed-form, analytic formula. While the results presented are encouraging, several issues deserve further research. One line of 545 possible development involves extending the capabilities of multidomain Line Sampling in order 546 to account for problems that involve either non Gaussian random variables or responses which are non linear with respect to the unknown random parameters. Preliminary research efforts conducted by the authors suggest that such an extension is feasible applying a so-called smooth indicator function, as suggested in [58, 59, 60]. Another path for development considers the ex-550 tension of multidomain Line Sampling for the analysis of problems involving parallel systems or 551 more general configurations. In particular, for the case of parallel systems, the results reported in

4 6. Acknowledgments

[19] could serve as a basis.

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Appendix A. Calculation of $c_{2k-1}^{(j)},\,c_{2k}^{(j)},\,\xi_{2k-1}^{(j)}$ and $\xi_{2k}^{(j)}$

The scheme for evaluating the line integral associated with the implementation of multidomain Line Sampling as presented in Section 3.3.2 requires few modifications for its implementation in case $\boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}} = 0$. To motivate the discussion, consider the schematic representation in Figure A.15, that depicts a particular case where $n = n_c = 2$, N = 3 and $i^{(1)} = i^{(2)} = i^{(3)} = 2$.

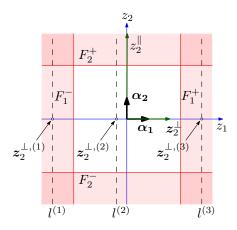


Figure A.15: Schematic representation of line associated with the application of multidomain Line Sampling in case $\alpha_k^T \alpha_{i^{(j)}} = 0$ $(n = n_c = 2)$.

As noted from Figure A.15, it is seen that component 1 fails for all points that belong to lines $l^{(1)}$ and $l^{(3)}$; on the contrary, it is seen that line $l^{(2)}$ never intersects the failure domain associated with component 1. Taking into account these observations and recalling the structure of eq. (21), it is concluded that for the case where $\boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}} = 0$, the unit step functions do not depend on $z_{i^{(j)}}^{\parallel}$ and that the Euclidean distances $c_{2k-1}^{i^{(j)}}$ and $c_{2k-1}^{i^{(j)}}$ must tend to either minus infinity or plus infinity in order to reflect that failure occurs or not with respect to the k-th component along line $l^{(j)}$. It is straightforward to demonstrate that the definitions for the distances $c_{2k-1}^{i^{(j)}}$ and $c_{2k-1}^{i^{(j)}}$ as presented in eqs. (22) and (23) must be extended as shown below in order to accommodate the

case where $\boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}} = 0$.

$$c_{2k-1}^{(j)} = \begin{cases} -\infty & \text{if } \boldsymbol{\alpha}_{k}^{T} \boldsymbol{\alpha}_{i(j)} = 0 \land \beta_{k}^{L} + \boldsymbol{\alpha}_{k}^{T} \boldsymbol{R}_{i(j)} \boldsymbol{z}_{i(j)}^{\perp,(j)} \leq 0 \\ -\frac{\beta_{k}^{L} + \boldsymbol{\alpha}_{k}^{T} \boldsymbol{R}_{i(j)} \boldsymbol{z}_{i(j)}^{\perp,(j)}}{\boldsymbol{\alpha}_{k}^{T} \boldsymbol{\alpha}_{i(j)}} & \text{if } \boldsymbol{\alpha}_{k}^{T} \boldsymbol{\alpha}_{i(j)} \neq 0 \\ \infty & \text{if } \boldsymbol{\alpha}_{k}^{T} \boldsymbol{\alpha}_{i(j)} = 0 \land \beta_{k}^{L} + \boldsymbol{\alpha}_{k}^{T} \boldsymbol{R}_{i(j)} \boldsymbol{z}_{i(j)}^{\perp,(j)} > 0 \end{cases}, k = 1, \dots, n_{c}, j = 1, \dots, N$$
(A.1)

$$c_{2k}^{(j)} = \begin{cases} -\infty & \text{if } \boldsymbol{\alpha}_{k}^{T} \boldsymbol{\alpha}_{i^{(j)}} = 0 \wedge \beta_{k}^{U} - \boldsymbol{\alpha}_{k}^{T} \boldsymbol{R}_{i^{(j)}} \boldsymbol{z}_{i^{(j)}}^{\perp,(j)} \leq 0 \\ \frac{\beta_{k}^{U} - \boldsymbol{\alpha}_{l}^{T} \boldsymbol{R}_{i^{(j)}} \boldsymbol{z}_{i^{(j)}}^{\perp,(j)}}{\boldsymbol{\alpha}_{k}^{T} \boldsymbol{\alpha}_{i^{(j)}}} & \text{if } \boldsymbol{\alpha}_{k}^{T} \boldsymbol{\alpha}_{i^{(j)}} \neq 0 \\ \infty & \text{if } \boldsymbol{\alpha}_{k}^{T} \boldsymbol{\alpha}_{i^{(j)}} = 0 \wedge \beta_{k}^{U} - \boldsymbol{\alpha}_{k}^{T} \boldsymbol{R}_{i^{(j)}} \boldsymbol{z}_{i^{(j)}}^{\perp,(j)} > 0 \end{cases}$$

$$(A.2)$$

The definition of the variables $\xi_{2k-1}^{(j)}$ and $\xi_{2k}^{(j)}$ as presented in eqs. (24) and (25) must be modified 576 well in order to accommodate the case where $\boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}} = 0$. It can be demonstrated that such 577 modification leads to: 578

$$\xi_{2k-1}^{(j)} = \begin{cases} 1 & \text{if } \boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}} \leq 0 \\ -1 & \text{otherwise} \end{cases}, \quad k = 1, \dots, n_c, \quad j = 1, \dots, N$$

$$\xi_{2k}^{(j)} = \begin{cases} 1 & \text{if } \boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}} \geq 0 \\ -1 & \text{otherwise} \end{cases}, \quad k = 1, \dots, n_c, \quad j = 1, \dots, N$$

$$(A.3)$$

$$\xi_{2k}^{(j)} = \begin{cases} 1 & \text{if } \boldsymbol{\alpha}_k^T \boldsymbol{\alpha}_{i^{(j)}} \ge 0 \\ -1 & \text{otherwise} \end{cases}, \ k = 1, \dots, n_c, \ j = 1, \dots, N$$
 (A.4)

Equations (A.1) to (A.4) as presented above allow calculating the line integral associated with 579 multidomain Line Sampling following the steps described in Section 3.3.2. In this sense, it is noted 580 that no modifications are required for eqs. (26) and (27) for the case where $\alpha_k^T \alpha_{i^{(j)}} = 0$.

Appendix B. Calculation of Interstory Drifts

The interstory drifts are calculated by means of the convolution integral, taking advantage of 583 the linearity of the response with respect to the ground acceleration. 584

$$\eta_i(t, \mathbf{z}) = \int_0^t h_i(t - \tau) p(\tau, \mathbf{z}) d\tau, \ i = 1, \dots, n_{\eta}$$
(B.1)

In the above equation, η_i represents the *i*-th interstory drift, h_i is the corresponding unit impulse response function and p represents the ground acceleration, which depends both on time t and a realization z of the standard Gaussian distribution. For the case considered in this contribution, the unit impulse response function is:

$$h_i(t) = \sum_{v=1}^{n_{\phi}} \frac{\kappa_v^T \phi_v \phi_v^T \rho}{\phi_v^T M \phi_v} \frac{1}{\omega_{d,v}} e^{-\zeta_v \omega_{n,v} t} \sin(\omega_{d,v} t), \quad i = 1, \dots, n_{\eta}$$
(B.2)

where κ_v , $i=1,\ldots,n_{\eta}$ is a vector that couples the degrees-of-freedom of the model for calculating the intestory drifts; $\boldsymbol{\rho}$ is a vector that couples the ground acceleration with the degrees-of-freedom of the model; $\boldsymbol{\phi}_v$, $v=1,\ldots,n_D$ are the eigenvectors associated with the eigenproblem of the undamped equation of motion involving the mass \boldsymbol{M} and stiffness \boldsymbol{K} matrices of the model; $\omega_{n,v}$, $v=1,\ldots,n_D$ are the natural frequencies of the system; ζ_v , $v=1,\ldots,n_D$ are the corresponding damping ratios; $\omega_{d,v}=\omega_{n,v}\sqrt{(1-\zeta_v^2)}$, $v=1,\ldots,n_D$ are the damped frequencies; and n_{ϕ} is the number of modes retained for modal analysis [61]. As the stochastic ground acceleration is represented by means of the Karhunen-Loève expansion, the interstory drift evaluated at discrete time instant t_k is approximated as:

$$\eta_i(t_k, \mathbf{z}) = \sum_{l=1}^k \Delta t \epsilon_l h_i(t_k - t_l) \mathbf{B}_{l:} \mathbf{z}, \ i = 1, \dots, n_{\eta}, \ k = 1, \dots, n_T$$
(B.3)

where Δt is the time discretization; n_T is the total number of discrete time instants; \boldsymbol{B}_l : denotes the l-th row of matrix \boldsymbol{B} (see eq. (1)), and ϵ_l is a coefficient depending on the numerical integration scheme used in the evaluation of the convolution integral. In this particular case, ϵ_l is chosen according to the trapezoidal integration rule [62], yielding $\epsilon_l = 1/2$ if l = 1 or l = k; otherwise, $\epsilon_l = 1$. From eq. (B.3), it is straightforward to see that the i-th interstory drift at the k-th time instant can be represented in the form $\eta_i(t_k, \boldsymbol{z}) = \boldsymbol{a}_{i,k}^T \boldsymbol{z}$, where $\boldsymbol{a}_{i,k}$ is a vector of constant coefficients. Such representation matches with the type of problems considered in this contribution, where the response of interest is a linear combination of a number of parameters following a Gaussian distribution (see Section 2.1).

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