### 1 Manuscript, October 2015

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# 4 A discretization procedure for rare events in Bayesian networks

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# 8 Abstract

9 Discrete Bayesian networks (BNs) can be effective for risk- and reliability assessments, in 10 which probability estimates of (rare) failure events are frequently updated with new information. To solve such reliability problems accurately in BNs, the discretization of 11 12 continuous random variables must be performed carefully. To this end, we develop an efficient discretization scheme, which is based on finding an optimal discretization for the 13 linear approximation of the reliability problem obtained from the First-Order Reliability 14 15 Method (FORM). Because the probability estimate should be accurate under all possible 16 future information scenarios, the discretization scheme is optimized with respected to the 17 expected posterior error. To simplify application of the method, we establish parametric formulations for efficient discretization of random variables in BNs for reliability problems 18 19 based on numerical investigations. The procedure is implemented into a software prototype. 20 Finally, it is applied to a verification example and an application example, the prediction of 21 runway overrun of a landing aircraft.

# 22 Keywords

23 Bayesian networks; discretization; near-real-time; structural reliability; updating

## 24 1 Introduction

25 For operational risk and reliability management, it is often desirable to compute the probability of a rare event F under potentially evolving information. Examples include 26 27 warning systems for natural and technical hazards, or the planning of inspection and 28 intervention actions in infrastructure systems. Ideally, this is achieved through Bayesian 29 updating of Pr(F) with the new information Z to the posterior probability Pr(F|Z). When 30 physically-based or empirical models for predicting the rare event exist, such updating is 31 possible with structural reliability methods (SRM) (Sindel and Rackwitz, 1998, Straub, 2011, 32 Straub et al., 2016). However, it is often difficult to perform the required computations in 33 near-real-time, due to a lack of efficiency or robustness. A modeling and computational 34 framework that does facilitate efficient Bayesian updating is the discrete Bayesian network 35 (BN). Hence it was proposed to combine SRMs with discrete Bayesian networks for near-36 real-time computations (Friis-Hansen, 2000, Straub and Der Kiureghian, 2010a, Straub and 37 Der Kiureghian, 2010b).

BNs are based on directed acyclic graphs (DAGs), to efficiently define a joint probability distribution  $p(\mathbf{Y})$  over a random vector  $\mathbf{Y}$  (Jensen and Nielsen, 2007, Kjaerulff and Madsen, 2013). The DAG of a BN, which is often referred to as the qualitative part of a BN, consists of a node for each variable in  $\mathbf{Y}$  and a set of directed links among nodes representing dependence among the variables. In the case of discrete BNs, conditional probability tables (CPTs) quantitatively define the type and strength of the dependence among the variables. The entries of the CPT of a variable  $Y_i$  are the probabilities for each state of  $Y_i$  conditional on

45 all possible combinations of states of its parents.

46 For hybrid BNs, which include both discrete and continuous variables, exact inference is 47 available only for two special cases, which are BNs with Gaussian nodes, whose means are 48 linear functions of their parents, and BNs, whose nodes are defined as a mixture of truncated 49 basic functions (MoTBFs) (Langseth et al., 2009, Langseth et al., 2012). Otherwise, approximate inference algorithms are available for hybrid BNs based on sampling techniques, 50 51 e.g. (Lerner, 2002, Hanea et al., 2006). However, these are computationally demanding and 52 not generally suitable for near-real-time decision support (Hanea et al., 2015). As an 53 alternative, the continuous random variables can be discretized, which enables the use of 54 exact inference algorithms that exist for general discrete BNs. These include the variable elimination algorithm (Zhang and Poole, 1994) and the junction tree algorithm (Lauritzen and 55 56 Spiegelhalter, 1988, Jensen et al., 1990).

57 The size of discrete BNs, and the associated computational effort, increases approximately 58 exponentially with the number of discrete states of its nodes, which motivates the 59 development of efficient discretization algorithms. While efficient discretization in the 60 context of machine learning and BNs in general has been investigated by multiple researchers 61 (Dougherty et al., 1995, Kotsiantis and Kanellopoulos, 2006), research on efficient 62 discretization in the context of engineering risk analysis or structural reliability has been 63 limited. In general, it is to be distinguished between static and dynamic discretization. While 64 the former discretizes the BN a-priori before entering evidence (offline), the latter is based on 65 an iterative scheme that updates the discretization scheme in function of the evidence (online).

66 Dynamic discretization for risk analysis applications has been developed mainly by (Neil et al., 2008), based on the work by (Kozlov and Koller, 1997). The procedure starts with an 67 68 initial discretization of a hybrid BN, for which an approximate entropy error is calculated. If 69 the error complies with a convergence criterion, the current discretization is accepted. 70 Otherwise the discretization is iteratively altered, by splitting the intervals with the highest 71 entropy error, until the convergence criterion is fulfilled. The approach is implemented in the 72 software AgenaRisk (Agena, 2005). Other dynamic discretization algorithms for reliability analysis have been proposed, e.g. in (Zhu and Collette, 2015) for dynamic BNs. The 73 74 advantage of dynamic discretization is its flexibility when evidence is entered in the BN, i.e. 75 when the model is updated with new observation.

76 Static discretization has the advantage of being computationally faster and simple to 77 implement. Some considerations for static discretization of BNs in reliability applications 78 have been presented in (Friis-Hansen, 2000, Straub, 2009, Straub and Der Kiureghian, 2010a). 79 As pointed out by (Friis-Hansen, 2000), for applications in which extreme events are 80 important, discretization of the distribution tails should be performed with care. Static 81 discretization facilitates a careful representation of these tails. However, the accuracy of the 82 static discretization varies with the available evidence. The difficulty is thus to find a 83 discretization scheme that is optimal under a wide variety of posterior distributions.

84 In this paper we derive a procedure for efficiently performing static discretization of 85 continuous reliability problems. An optimal discretization scheme is sought, which minimizes 86 the expected approximation error with respect to possible future observations (evidence). To 87 solve this optimization problem, we propose to approximate the reliability problem by the 88 First-Order Reliability Method (FORM). Section 2 of the paper describes the proposed 89 methodology. Section 3 presents numerical parameter studies, and simple parametric relations 90 for defining an efficient discretization scheme are derived. In Section 4, the procedure is 91 applied to a set of verification examples and to the computation of the probability of runway 92 overrun of a landing aircraft. While the theory is introduced for problems with only one 93 design point, considerations regarding problems with multiple design points are given in the 94 last verification example and in the discussion.

## 95 **2 Methodology**

#### 96 **2.1 Structural reliability**

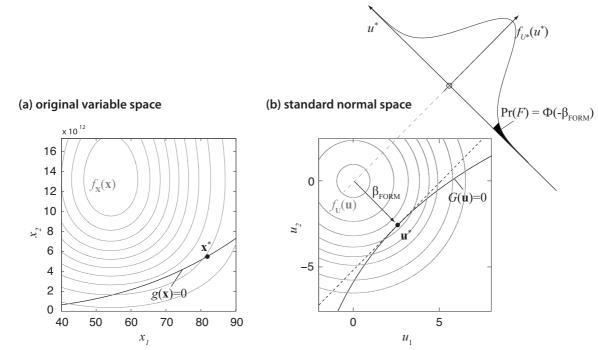
97 Since the 1970s structural reliability methods have been developed and applied in the 98 engineering community to estimate failure probabilities Pr(F) of components or systems, 99 based on physical or empirical models. The performance of engineering components is 100 described by a limit state function (LSF)  $g(\mathbf{x})$ , where  $\mathbf{X} = [X_1; ...; X_n]$  is a vector of basic 101 random variables influencing the performance of the component. By definition, failure 102 corresponds to  $g(\mathbf{x})$  taking non-positive values, i.e. the failure event is  $F = \{g(\mathbf{X}) \le 0\}$ .  $g(\mathbf{x})$ 103 includes the physical or engineering model, which is often computationally demanding. The 104 probability of failure is calculated by integrating the probability density function (PDF) of  $\mathbf{X}$ , 105  $f_{\mathbf{X}}(\mathbf{x})$ , over the failure domain:

$$\Pr(F) = \int_{g(\mathbf{x}) \le 0} f_{\mathbf{X}}(\mathbf{x}) d\mathbf{x}$$
(2)

106 The formulation can be extended to the reliability of general systems by defining the failure 107 domain as a combination of series and parallel systems (Ditlevsen and Madsen, 2007). In the general case, there is no analytical solution to Eq. 2 and the integral is potentially high-108 109 dimensional. For this reason, structural reliability methods (SRMs) are applied to approximate 110 it. These include the first- and the second order reliability method (FORM and SORM) as well as a large variety of sampling methods, including importance sampling methods such as 111 directional importance sampling, and sequential sampling methods such as subset simulation. 112 These methods are well-documented in the literature (Au and Beck, 2001, Rackwitz, 2001, 113 114 Der Kiureghian, 2005, Ditlevsen and Madsen, 2007).

## 115 **2.2** First order reliability method (FORM)

116 To obtain an approximation of the probability of failure through FORM, the LSF  $g(\mathbf{X})$  is 117 transformed to an equivalent LSF  $G(\mathbf{U})$  in the space of uncorrelated standard normal random 118 variables  $\mathbf{U} = [U_1; ...; U_n]$  (Fig. 1). The transformation is probability conserving, so that 119  $\Pr[g(\mathbf{X}) \leq 0] = \Pr[G(\mathbf{U}) \leq 0] = \Pr(F)$ . A suitable transformation for this purpose, which is 120 consistent with the BN, is the Rosenblatt transformation (Hohenbichler and Rackwitz, 1981). In case all basic random variables are independent, this transformation reduces to the 121 marginal transformations:  $U_i = \Phi^{-1}[F_{X_i}(X_i)]$ , with  $\Phi^{-1}$  being the inverse standard normal 122 123 CDF.



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Figure 1. Design point and linear approximation of the limit state surface. Left side: original random variable space; right side: standard normal space (from (Straub, 2014a)).

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The FORM approximation of Pr(F) is obtained by substituting the LSF in U-space  $G(\mathbf{U})$  by a linear function  $G_L(\mathbf{U})$ , i.e. a first-order Taylor expansion of  $G(\mathbf{U})$ . The key idea of FORM is to choose as the expansion point the so-called design point  $\mathbf{u}^*$ , which is the point that minimizes  $\|\mathbf{u}^*\|$  subject to  $G_L(\mathbf{U}) \leq 0$ .  $\mathbf{u}^*$  also known as the most likely failure point, as it is the point in the failure domain with the highest probability density. Since all marginal distributions of the standard uncorrelated multinormal distribution are standard normal, it can be shown that the FORM probability of failure  $Pr[G_L(\mathbf{U}) \leq 0]$  is:

$$\Pr[G_L(\mathbf{U}) \le 0] = \Phi(-\beta_{FORM}) \tag{3}$$

where  $\Phi$  is the standard normal CDF and  $\beta_{FORM}$  is the distance from the origin to the design point, i.e.  $\beta_{FORM} = ||\mathbf{u}^*||$ . The problem thus reduces to finding the design point  $\mathbf{u}^*$ . If  $G(\mathbf{U})$  is linear, the FORM solution of the probability of failure is exact, otherwise it is an approximation, which however is sufficiently accurate in most practical applications with limited numbers of random variables (Rackwitz, 2001).

140 The linearized LSF  $G_L(\mathbf{U})$  can be written as:

$$G_L(\mathbf{U}) = \beta_{FORM} - \alpha^{\mathrm{T}} \mathbf{U}$$
(4)

141 where  $\alpha = [\alpha_1, ..., \alpha_n]$  is the vector of FORM importance measures. These importance 142 measures are defined as:

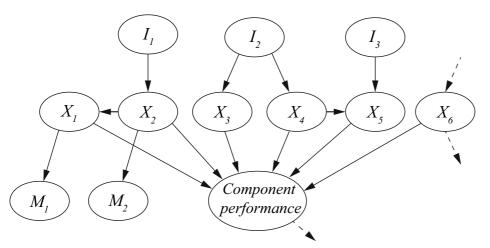
$$\alpha_i = \frac{u_i^*}{\beta_{FORM}} \tag{5}$$

where  $u_i^*$  is the *i*-th component of the design point coordinates. The  $\alpha_i$ 's take values between -1 and 1, and it is  $\|\boldsymbol{\alpha}\| = 1$ .  $\alpha_i$  is 0, if the uncertainty on  $U_i$  has no influence on  $\Pr(G_L(\mathbf{U}) \leq 0)$ , and it is 1 or -1, if  $U_i$  is the only random variable affecting  $\Pr(g_L(\mathbf{U}) \leq 0)$ . When the original random variables  $X_i$  are mutually independent, the  $\alpha_i$ 's are readily applicable also in the original space, otherwise the  $\alpha_i$ 's can be transformed as described in (Der Kiureghian, 2005).

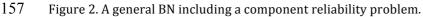
## 149 2.3 Treatment of a reliability problem in a BN

We combine discrete BNs and structural reliability concepts to facilitate updating of rare event (failure) probabilities under new observations. The general problem setting is illustrated in the BN of Fig. 2. We here limit the presentation to component reliability problems; system problems are considered later. The binary random variable 'Component performance' is described by the LSF  $g(\mathbf{X})$ .

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The basic random variables **X** of the model are included in the BN as parents of 'Component performance'. The nodes  $M_i$  represent measurements of individual random variables  $X_i$ , and nodes  $I_j$  represent factors influencing the basic random variables. Dependence between the variables in **X** is modeled either directly by links among them (here  $X_2 \rightarrow X_1$  and  $X_4 \rightarrow X_5$ ) or through common influencing factors (here  $I_2 \rightarrow X_3$  and  $I_2 \rightarrow X_4$ ). The component performance node can have (multiple) child nodes, which, however, does not impact the discretization of the reliability problem.

- 165 Ultimately, the goal is to predict the component performance, i.e. Pr(F), conditional on other 166 model parameters, such as measurements  $M_i$  or influencing variables  $I_j$ . Whenever new 167 evidence on these variables is available, the BN should be evaluated (in near-real time) 168 utilizing exact BN inference algorithms.
- 169 To enable exact inference algorithms, all continuous random variables are discretized. These
- include the **X**, and possibily the  $M_i$  and  $I_j$ . In the general case, the computational effort for
- solving the BN is a direct function of the CPT size of 'Component performance'. The size of

172 this CPT is  $2\prod_{i=1}^{n} n_i$ , where n is the number of random variables in **X**, and  $n_i$  is the number

173 of states used for discretizing  $X_i$ . In this paper we do not describe the discretization of random

174 variables  $M_i$  and  $I_i$ , since it is typically straightforward and does not contribute significantly

175 to computational performance. The key parameter for computational efficiency and accuracy

176 is the discretization scheme for **X**, which is described in sections 2.5 and 2.6.

#### 177 2.4 Simplification of BNs through node removal

178 Removing random variables from a BN is one possibility to reduce the computational effort 179 associated with a model. A formal approach for removing nodes from a BN is described in 180 (Straub and Der Kiureghian, 2010b). In order to decide which nodes to remove from the BN 181 the following questions should be considered:

- 182 • Which random variables are relevant for prediction? (These include 'Component 183 performance'.)
- 184 • Which random variables can potentially be observed? (These include the measurement 185 variables.)
- Which random variables simplify the modeling of dependencies? (These are e.g. 186 common influencing factors such as  $I_2$  in Fig. 2.) 187
- 188 • For which random variables is it desirable to explicitly show their influence on 189 component performance?

190 If a random variable does not belong to any of these categories, the corresponding node in the 191 BN can be removed. Since the computational efficiency of the model is governed by the size 192 of the CPT of the 'Component state' node, the primary interest is in removing basic random 193 variables  $X_i$  from the BN. As a measure for the relevance of a basic random variable, 194 importance measures  $\alpha_i$  from a FORM analysis may be used. To better understand the relation between  $\alpha_i$  and  $X_i$ 's relevance for prediction, consider a linearized LSF  $G_L(\mathbf{U})$ . 195 196 Following (Der Kiureghian, 2005), the variance of  $G_L(\mathbf{U})$  can be decomposed as:

$$\sigma_{G_L}^2 = \|\nabla G\|^2 (\alpha_1^2 + \alpha_2^2 + \dots + \alpha_n^2)$$
(6)

197 where  $\nabla G$  denotes the gradient vector of the non-linearized LSF  $G(\mathbf{U})$ . From Eq. 6 it is seen

that a random variable  $X_i$  with corresponding  $\alpha_i$  accounts for  $\alpha_i^2 \cdot 100\%$  of the variance  $\sigma_{G_i}^2$ . 198

Therefore, observing a random variable  $X_i$  with  $\alpha_i = 0.1$  will reduce the variance  $\sigma_{G_L}^2$  by 1%, 199 200 whereas observing  $X_i$  with  $\alpha_i = 0.5$  will reduce  $\sigma_{G_L}^2$  by 25%.

#### 201 2.5 Discretization of basic random variables

202 For ease of presentation, we first consider discretization of statistically independent basic 203 random variables **X**, i.e. the special case of the BN in Fig. 2 in which the  $X_i$ 's have no parents<sup>1</sup>.

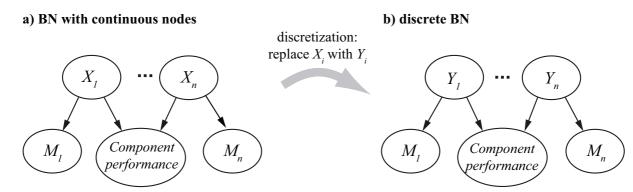
<sup>&</sup>lt;sup>1</sup> In a BN, basic random variables  $\mathbf{X}$  are independent if they are not connected through links, if they have no common (unknown) ancestors and if no evidence is available on any of their joint descendants.

The proposed procedure is extended to the general case of dependent basic random variables thereafter.

## 206 2.5.1 Independent basic random variables

207 The situation is illustrated in Fig. 3. The performance of the component depends on n208 statistically independent random variables and is described by a LSF  $g(\mathbf{X}) = g(X_1, \dots, X_n)$ . 209 For all basic random variables  $X_i$ , corresponding measurements  $M_i$  can be performed. To obtain an equivalent discrete BN, the continuous  $X_i$  are replaced by the discrete random 210 211 variables  $Y_i$ , and the LSF is replaced by the CPT of component performance conditional on 212  $\mathbf{Y} = [Y_1; ...; Y_n]$ . For each discrete random variable  $Y_i$  with  $n_i$  states 1,2, ...,  $n_i$ , we define a discretization scheme  $D_i = [d_0, d_1, ..., d_{n_i-1}, d_{n_i}]$  consisting of  $n_i + 1$  interval boundaries. 213 The first and the last interval boundaries are given by the boundaries of  $X_i$ 's outcome space. 214

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Figure 3. Representation of a basic reliability problem with *n* independent basic random variables in a BN. Left: original problem with continuous basic random variables  $X_i$ , right: discrete BN, in which  $X_i$ s are substituted with discrete nodes  $Y_i$ .

220 Since here the  $X_i$ , and thus the  $Y_i$ , have no parents, the PMF of  $Y_i$  is defined as:

$$p_{Y_i}(j) = F_{X_i}(d_j) - F_{X_i}(d_{j-1}); \quad \text{with } j \in [1, \dots, n_i]$$
(7)

where  $F_{X_i}$  denotes the cumulative distribution function (CDF) of  $X_i$ . The probability of failure corresponding to the discrete BN in Fig. 3b can be calculated as:

$$\Pr(F) = \sum_{y_1=1}^{n_1} \dots \sum_{y_n=1}^{n_n} p_{Y_1}(y_1) \cdot \dots \cdot p_{Y_n}(y_n) \cdot \Pr(F|Y_1 = y_1 \cap \dots \cap Y_n = y_n)$$
(8)

223 Note that the discretization does not introduce any approximation error here, as long as the

224 conditional  $\Pr(F|Y_1 = y_1 \cap ... \cap Y_n = y_n)$  is computed exactly.

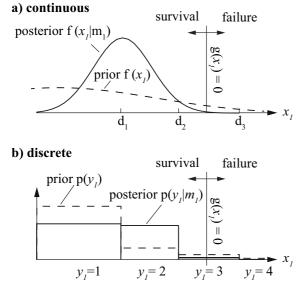
Once measurements from the nodes  $\mathbf{M} = [M_1; ...; M_n]$  are available, the conditional failure probability can be calculated as:

$$\Pr(F|\mathbf{M} = \mathbf{m}) \approx \frac{1}{p_{\mathbf{M}}(\mathbf{m})} \sum_{y_1=1}^{n_1} \dots \sum_{y_n=1}^{n_n} p_{Y_1}(y_1) \cdot p_{M_1|Y_1}(m_1|y_1) \cdot \dots \cdot p_{Y_n}(y_n) \cdot p_{M_n|Y_n}(m_n|y_n)$$
(9)

where  $\Pr(F|Y_1 = y_1, ..., Y_n = y_n)$  is the conditional probability of component failure given  $y_1, ..., y_n$ . If no measurements are available for some of the basic random variables, the corresponding likelihood terms  $p_{M_i|Y_i}(m_j|y_j)$  are simply omitted in Eq. 9.

While the computation of the unconditional failure probability following Eq. 8 is exact, the 230 231 computation of the conditional failure probability through Eq. 9 is only an approximation. 232 The reason is that the dependence between the measurement variable  $M_i$  and the 'Component 233 performance' variable is not fully captured in the discrete BN (see also Straub and Der 234 Kiureghian, 2010b). In Fig. 4, this is illustrated for a reliability problem with one basic 235 random variable  $X_i$ . Both the continuous distribution (Fig. 4a) and the corresponding discretized distribution (Fig. 4b) are updated correctly after observing  $M_1$ . However, for Eq. 9 236 to be exact, also the conditional failure probabilities  $Pr(F|Y_1 = y_1)$  would need to be updated. 237 This can be observed in Fig. 4a: in interval  $Y_1 = 3$ , which is the one cut by the limit state 238 239 surface, the ratio of the probability mass in the failure domain to that in the safe domain changes from the prior to the posterior case, i.e.  $\Pr(F|Y_1 = 3) \neq \Pr(F|Y_1 = 3, M_1 = m_1)$ . 240 241 Since the computation of the conditional failure probability following Eq. 9 is based on the 242 prior probability  $Pr(F|Y_1 = 3)$ , the discretization introduces an approximation in this case. 243 The error occurs only in the intervals that are cut by the limit state surface.

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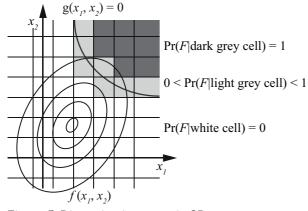


245 246

6 Figure 4. Discretization error in 1D.

In the simple one-dimensional case of Fig. 4, an optimal discretization approach would be to discretize the whole outcome space in two intervals, one capturing the survival and one the failure domain. This discretization would have zero approximation error. However, already in a two-dimensional case, such a solution is not possible. This is illustrated in Fig. 5, where the cells cut by the limit state surface are indicated in grey. The failure probability conditional on measurements calculated according to Eq. 9 will necessarily be an approximation. The approximation error will be small, if the contribution of the cells cut by the limit state surface

- 254 (the grey cells in Fig. 5) to the total failure probability is small. An efficient discretization will
- thus limit this contribution with as few intervals as possible.



- 256  $f(x_1, x_2)$ 257 Figure 5. Discretization error in 2D.
- 258 2.5.2 Dependent basic random variables

Eqs. (7-9) must be adjusted when dependence among the  $X_i$ 's is present, in accordance with the case-specific BN structure. However, the principles outlined above for independent  $X_1, ..., X_n$  hold equally for dependent basic random variables: The discretization error is a function of the cells cut by the limit state function.

When determining an optimal discretization, we propose in the following to find the FORM approximation of the reliability problem, which can readily account for the dependence among the random variables. Hence, there is no need to distinguish between cases with independent or dependent random variables.

# 267 2.6 Efficient discretization

268 2.6.1 Optimal discretization of linear problems in standard normal space

To find an efficient discretization of **X**, we consider the FORM solution to the reliability problem. Evaluating the linearized FORM LSF  $G_L(\mathbf{U})$  is computationally inexpensive once the design point  $\mathbf{u}^*$  is known. Therefore, it is feasible to find a discretization of **U** that is optimal for the event { $G_L(\mathbf{U}) \le 0$ } through optimization. If  $G(\mathbf{U})$  is not strongly non-linear, this solution will be an efficient discretization for { $G(\mathbf{U}) \le 0$ } and, after a transformation to the original space, also for { $g(\mathbf{X}) \le 0$ }.

- As discussed in Section 2.5.1, the approximation error of the discretization is associated with the change from the prior to the posterior distribution of the basic random variables. A measure of optimality must thus consider possible measurements of **X** or **U**. We consider hypothetical measurements  $\tilde{M}_i$  (this notation is used to distinguish hypothetical measurements from actual measurements  $M_i = m_i$ ) of all standard normal random variables  $U_i$  with additive measurement error  $\varepsilon_i \sim N(0, \sigma_{\varepsilon_i})$ . Since both the prior distribution and the measurement error
- are normal distributed, the likelihood is also normal distributed:

$$\widetilde{M}_i | \{ U_i = u_i \} \sim N(u_i, \sigma_{\varepsilon_i})$$
(10)

282 The posterior, i.e. the conditional distribution of  $U_i$  given a measurement outcome  $\widetilde{M}_i = \widetilde{m}_i$ ,

283 is the normal distribution with mean  $\frac{1}{1+\sigma_{\varepsilon_i}^2}\widetilde{m}_i$  and standard deviation  $\sqrt{\left(1-\frac{1}{1+\sigma_{\varepsilon_i}^2}\right)}$ .

We define an error measure based on comparing the true posterior probability of failure  $P_{F|\tilde{M}}(\tilde{m})$  with the posterior probability of failure calculated from the discretized **U**, denoted by  $\hat{P}_{F|\tilde{M}}(\mathbf{d}; \tilde{m})$ . Here, **d** are the parameters defining the discretization. The proposed error measure is:

$$e(\mathbf{d}, \widetilde{\mathbf{m}}) = \left| \frac{\log_{10} \widehat{P}_{F|\widetilde{M}}(\mathbf{d}; \widetilde{\mathbf{m}}) - \log_{10} P_{F|\widetilde{M}}(\widetilde{\mathbf{m}})}{\log_{10} P_{F|\widetilde{M}}(\widetilde{\mathbf{m}})} \right|$$
(11)

The error measure of Eq. 11 represents a tradeoff between the absolute and the relative error.
It weights the (logarithmic) relative error by the magnitude of the posterior failure probability.
This ensures that the same relative error is considered worse at a higher probability level

291 compared to an error at a lower probability level.

A-priori, the measurement outcomes are not known. Hence we define the optimal discretization as the one that minimizes the expected preposterior error  $E_{\tilde{M}}[e(\mathbf{d}, \tilde{M})]$ :

$$\mathbf{d}^{opt} = \arg\min_{\mathbf{d}} \mathbb{E}_{\widetilde{M}}[e(\mathbf{d}, \widetilde{M})] = \arg\min_{\mathbf{d}} \int_{\widetilde{M}} e(\mathbf{d}, \widetilde{\mathbf{m}}) f_{\widetilde{M}}(\widetilde{\mathbf{m}}) d\widetilde{\mathbf{m}}$$
(12)

The optimization is thus based on the computation of an expected value with respect to the possible measurements outcomes  $\tilde{M}$  before having taken any measurements. This is analogous to a preposterior analysis (Raiffa and Schlaifer, 1961, Straub, 2014b). However, unlike in traditional preposterior analysis, the objective is not to identify an optimal action under future available information, but to find the optimal discretization parameters  $\mathbf{d}^{opt}$ . The integral in Eq. 12 is evaluated through a simple Monte Carlo approach. All  $\tilde{M}_i$  have the normal distribution with zero mean and variance  $1 + \sigma_{\epsilon}^2$ .

301 The parameters in **d** describing the discretization scheme are:

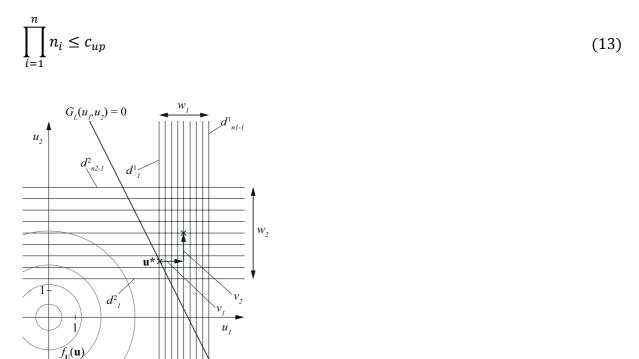
302 –  $n_i$ : number of intervals used to discretize each random variable  $U_i$ ,

- 303  $w_i$ : width of the discretization frame in the dimension of  $U_i$ , and
- $v_i$ : position of the midpoint of the discretization frame relative to the design point

These parameters are illustrated in Fig. 6. For a problem with *n* basic random variables, the full set of optimization parameters is  $\mathbf{d} = [w_1, \dots, w_n, n_1, \dots, n_{n-1}, v_1, \dots, v_n]$ .

307 Clearly, the discretization error reduces with increasing  $n_i$ . Because the computational 308 efficiency of the final BN is a direct function of the size of the CPT associated with 309 component performance, which is  $\prod_{i=1}^{n} n_i$ , we constrain its size. To this end, we define  $c_{up}$  as

- 310 the maximum allowed number of parameters of the CPT of the component state node. This
- 311 puts a constraint on the optimization of Eq. 12:



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Figure 6. Schematic representation of a discretization of a linear 2D reliability problem.  $w_i$  is the distance between interval boundaries  $d_1^i$  and  $d_{n_i-1}^i$ . All intervals between these boundaries are equi-spaced.  $v_i$  is the position of the midpoint of the discretization frame relative to the design point  $\mathbf{u}^*$  in dimension *i*.

The optimization is implemented through a two-level approach. The optimization of the continuous parameters width  $w_i$  and position of the discretization frame  $v_i$  for all i = 1, ..., nis carried out using unconstrained nonlinear optimization for fixed values of  $n_i$ . The optimization of the discrete  $n_i$  is performed through a local search algorithm. Note that the optimization is performed offline, i.e. prior to running the BN, hence it does not affect the goal of near-real time performance of the BN. Furthermore, in section 3 a heuristic is derived that can replace the time-consuming solution of the optimization problem.

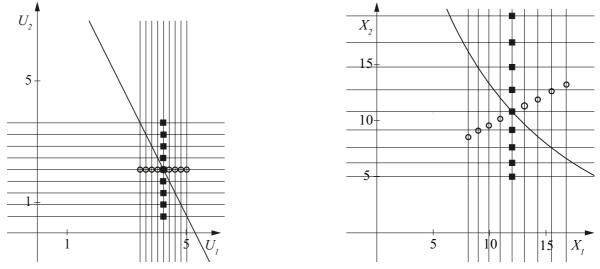
# 323 2.6.2 Efficient discretization of the original random variables **X**

324 Since the nodes in the BN represent random variables X in their original outcome space, the 325 discretization schemes, which are derived for the corresponding standard normal random 326 variables U, need to be transformed to the X-space. In the case of mutually independent random variables X<sub>i</sub>, any point on the i-th interval boundary in U-space – if transformed – will 327 328 result in the same corresponding i-th interval boundary in X-space. This is not the case for 329 dependent random variables X<sub>i</sub>, where a mapping of the interval boundaries in U-space to X-330 space will not lead to an orthogonal discretization scheme in X-space. To preserve orthogonality throughout the transformation, we propose to represent each interval boundary 331 332 through a characteristic point and determine the boundary in X-space through a 333 transformation of this point. For transforming the interval boundary of X<sub>i</sub>, the characteristic

- 334 point is selected as the design point u<sup>\*</sup>, where the i-th element is substituted by the coordinate
- of the interval boundary. In Fig. 7 this is shown for an example with n = 2 random variables.

a) Discretization in U-space

### b) Transformed discretization



336

Figure 7. Transformation of a discretization scheme from U-space to X-space. To preserve orthogonality each interval boundary in U-space is represented by a characteristic point. The random variables  $X_1$  and  $X_2$  are Weibull distributed with scale and shape parameter 1 and their correlation is 0.5.

# **340 3 Development of an efficient discretization procedure**

# 341 **3.1 Optimization of the FORM approximation**

We present the optimal discretization for the FORM approximation  $G_L(\mathbf{U})$  for n = 2 and n = 3 dimensions. Extension to higher numbers of random variables is discussed. Because the linear LSF employed in FORM is described only by the reliability index  $\beta_{FORM}$  and the vector  $\boldsymbol{\alpha}$  of FORM sensitives (Eq. 4), it facilitates parametric studies.

Initially, we consider a reliability index  $\beta_{FORM} = 4.26$ , corresponding to a probability of 346 failure of  $10^{-5}$ . The standard deviation of the additive measurement error is set to either 347  $\sigma_{\varepsilon} = 0.5$  or  $\sigma_{\varepsilon} = 1.0$ . Different combinations of FORM sensitivity values  $\alpha_i$  are selected, to 348 investigate their effect on the optimal discretization. In all investigated cases, we find that the 349 position of the midpoint of the optimal discretization frame coincides with the design point, 350 i.e.  $v_i^{opt} = 0$ . The optimal discretization widths  $w_i^{opt}$  vary significantly with the importance 351 measures  $\alpha_i$ , as shown in section 3.1.2. However, the optimal number of intervals  $n_i^{opt}$  is 352 approximately the same for all random variables in all investigated cases, independent of the 353  $\alpha_i$  values, i.e.  $n_i^{opt} = c_{up}^{1/n}$ . 354

355 3.1.1 On why the optimal number of intervals  $n_i^{opt}$  is independent of  $\alpha_i$ 

To better understand why the number of intervals does not depend on  $|\alpha_i|$  (for  $|\alpha_i|$  that are significantly larger than 0), recall that an efficient discretization scheme should focus on the area around the limit state surface. More precisely, the discretization error in the posterior 359 case is induced by the cells that are cut by the limit state surface. Exemplarily, Fig. 8 shows a linear problem in standard normal space with two basic random variables  $U_1$  and  $U_2$ , where 360  $\alpha_1 = 0.45$  and  $\alpha_2 = 0.89$ . While Fig. 8a shows an optimal discretization with 5 intervals per 361 362 dimension, Fig. 8b shows a discretization scheme, where the more important random variable  $U_2$  is discretized with 6 intervals and  $U_1$  with 4 intervals. In both cases, the discretization 363 364 frame is centered at the design point. It is observed that the probability mass of the (grey) cells, associated with the discretization error in the posterior case, is higher in Fig. 8b than for 365 the optimal discretion scheme in 8a. In the example shown here it is  $1.4 \cdot 10^{-3}$  compared to 366  $2.5 \cdot 10^{-4}$ . 367

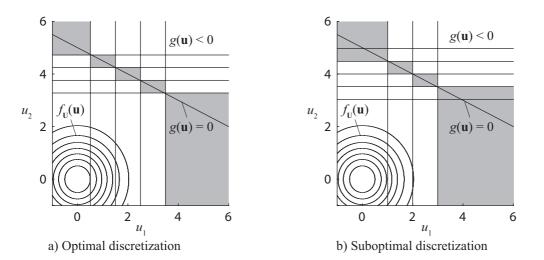


Figure 8. Linear problem in standard normal space, with two random variables  $U_1$  and  $U_2$ , where  $\alpha_1 = 0.45$  and  $\alpha_1 = 0.89$ . The intervals cut by the limit state surface, i.e. those which potentially lead to a posterior discretization error are marked in grey.

For input random variables  $X_i$  with a value of  $|\alpha_i|$  close to zero, the above observations do not hold. Following Section 2.4, these variables should be removed from the BN prior to discretization.

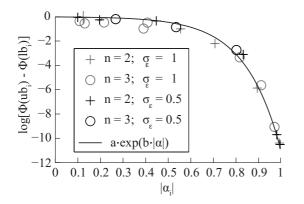
375 3.1.2 Dependence of the optimal discretization width on  $\alpha_i$ 

368

In the optimization, it is found that the optimal discretization width  $w_i^{opt}$  varies strongly with the random variable's importance, expressed through  $\alpha_i$ . It is reminded that the width  $w_i$ describes the domain in which a fine discretization mesh is applied (Fig. 6). In general,  $w_i^{opt}$ decreases with increasing  $\alpha_i$ . This effect can be observed in Fig. 8a, where  $U_2$  is the more important input random variable and it is  $w_2^{opt} < w_1^{opt}$ .

381 A clear relation between  $w_i^{opt}$  and  $\alpha_i$  can be observed by plotting the probability mass 382 enclosed by  $w_i^{opt}$  against  $\alpha_i$ , as shown in Fig. 9. The results of Fig. 9 indicate that the 383 probability mass contained within this interval should be a direct function of  $\alpha_i$ . The more 384 important the variable, the finer the discretization around the design point should become. 385 The observed relationship between this probability mass and  $\alpha_i$  follows a clear trend, and a 386 function can be fitted (Fig. 9). Neither the dimensionality of the problem nor the standard 387 deviation of the measurement error appear to have an influence on this relation. However, as

shown in the following section, it is found that the relation does depend on the prior failure probability of the problem (i.e. on  $\beta_{FORM}$ ) and on the number of intervals  $n_i$  used to discretize the domain.



391

392 Figure 9. Logarithm of the probability mass enclosed by the discretization frame plotted against  $\alpha_i$ .  $\Phi$ 393 denotes the standard normal CDF and  $ub_i$  respectively  $lb_i$  the last (upper) and the first (lower) interval 394 bound in dimension *i*.

To facilitate the application in practice and extending the results to larger numbers of random variables, in section 3.2 parametric functions are fitted to the optimization results to capture the dependency between the optimal discretization width  $w_i^{opt}$  and the FORM importance measures  $\alpha_i$ .

399 3.1.3 Dependence of the optimal discretization on the reliability index  $\beta$  and the

400 number of discretization cells  $c_{up}$ 

401 The influence of the prior failure probability and the maximum size of the CPT,  $c_{up}$ , on the 402 optimal discretization is investigated through 10 problems with n = 2 random variables in 403 standard normal space. The FORM importance measures of the random variables are selected 404 between 0.1 to 0.995 and the standard deviation of the measurement error is fixed to  $\sigma_{\varepsilon} = 1.0$ . 405 We find that the optimal discretization frame is generally centered at the design point, i.e. 406  $v_i^{opt} = 0$ , and that the intervals are distributed uniformly among the dimensions.

Firstly, we vary the maximum CPT size  $c_{up}$ , i.e. the total number of discretization cells. The reliability index is  $\beta_{FORM} = 5.2$ . Fig. 10 shows the influence of  $c_{up}$  on the resulting width of the discretization frame  $w_i$ . Three cases are considered:  $c_{up} = 25$ ,  $c_{up} = 100$  and  $c_{up} = 400$ . These choices correspond to 5, 10 and 20 intervals for each random variable. The left side of Fig. 10 shows the relation between the optimal  $w_i$  and  $|\alpha_i|$  The right side of Fig. 10 shows the same relation, where the  $w_i$ 's are scaled as in Fig. 9, i.e. the logarithm of the probability mass

- 413 enclosed by the outer interval boundaries is depicted. As in Fig. 9, there is a clear dependence
- 414 between the scaled  $w_i$  values and the  $|\alpha_i|$ 's. The interval frames increase with increasing
- 415 number of random variables.

416 Secondly, we vary the prior failure probability from  $10^{-3}$  ( $\beta = 3.1$ ) to  $10^{-7}$  ( $\beta = 5.2$ ). The 417 results are shown in Fig. 11. Again, a distinct dependence between the scaled  $w_i$  values and 418 the  $|\alpha_i|$ 's is found. The interval frames decrease with increasing reliability index (with 419 decreasing failure probability).

#### 420 **3.2** Parametric function of optimal discretion frame

421 As evident from Fig. 10 and Fig. 11, there is a clear dependence of the probability mass

- 422 enclosed by the optimal discretization frame (with width  $w_i$ ) on the FORM sensitivity values
- 423  $|\alpha_i|$ . The following parameteric function captures this dependence:

$$\log(\Phi(ub_i) - \Phi(lb_i)) = a \cdot \exp(b \cdot |\alpha_i|) \tag{14}$$

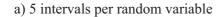
424  $ub_i$  is the upper and  $lb_i$  the lower interval boundary in dimension *i*, such that  $w_i = ub_i - lb_i$ . 425 *a* and *b* are the parameters of the exponential function. This function is depicted in Figs. 10 426 and 11. Tab. 1 shows the parameter values *a* and *b* for the different combinations of the prior 427 reliability index  $\beta$  and number of intervals per dimension

428 Table 1. Parameters *a* and *b* of Eq. 14 for  $\beta = 3.1$ ,  $\beta = 4.3$  and  $\beta = 5.2$  as well as 5, 10 and 20 intervals per dimension.

$\begin{bmatrix} a, \\ b \end{bmatrix}$	<i>n</i> <sub><i>i</i></sub> = 5	$n_{i} = 10$	$n_i = 20$
		$\begin{bmatrix} -1.6 \cdot 10^{-2}, \\ 5.8 \end{bmatrix}$	
$\beta = 4.3$	$\begin{bmatrix} -0.15, \\ 4.3 \end{bmatrix}$	$\begin{bmatrix} -2.4 \cdot 10^{-2}, \\ 6.1 \end{bmatrix}$	$\begin{bmatrix} -2.1 \cdot 10^{-2}, \\ 6.2 \end{bmatrix}$
$\beta = 5.2$	$\begin{bmatrix} -0.36, \\ 3.7 \end{bmatrix}$	$\begin{bmatrix} -0.11, \\ 5.0 \end{bmatrix}$	$\begin{bmatrix} -3.7 \cdot 10^{-2}, \\ 6.0 \end{bmatrix}$

From the left sides of Fig. 10 and Fig. 11, it can be observed that the relation between  $\alpha_i^2$  and the optimal  $w_i$  is fairly diffuse for random variables with  $|\alpha_i| < 0.6$ . Here, the parametric relationship of Eq. 14 is less accurate. However, these random variables by definition have lower importance on the reliability estimate. Hence, the inaccuracy of Eq. 14 for random variables with  $|\alpha_i| < 0.6$  is not critical, as is confirmed by the numerical investigations performed in the remainder of the paper.

The parameter values of Tab. 1 are derived from two-dimensional problems. In Fig. 9 it is shown that there are no notable differences between two and three dimensions. On this basis, it is hypothesized that the heuristics are applicable also to problems with higher dimensions. This assumption is furthermore supported by the verification examples presented in chapter 4, where the heuristics are applied also to four-dimensional problems without any notable deterioration in the results.



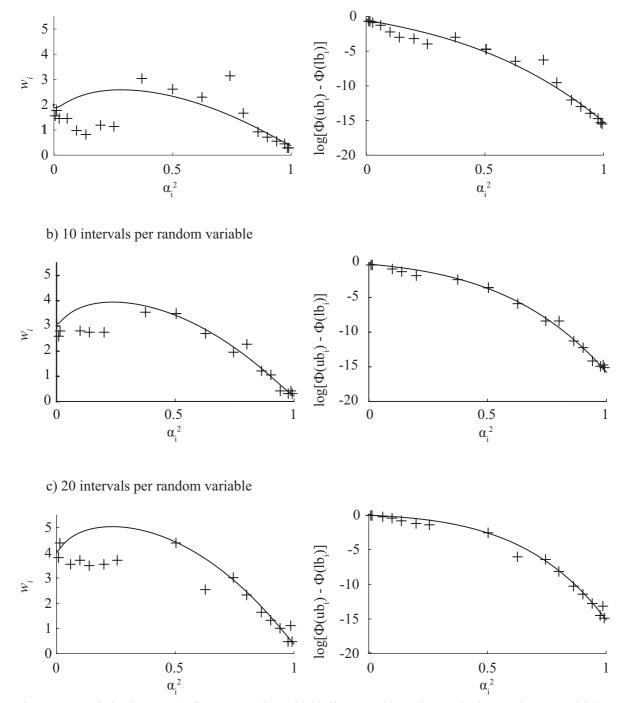


Figure 10. Optimization results for 10 two-dimensional, linear problems in standard normal space, which are discretized with 5, 10 and 20 intervals per dimension. In all cases the prior failure probability is  $10^{-7}$  ( $\beta = 5.2$ ). The crosses represent the optimization results. The solid lines are the fitted parametric functions (Eq. 14). The left-hand side shows the relation between the width of a discretization frame  $w_i$  and  $|\alpha_i|$  and the right-hand side shows the relation between the probability mass enclosed by the discretization frame with width  $w_i$  and  $|\alpha_i|$ .

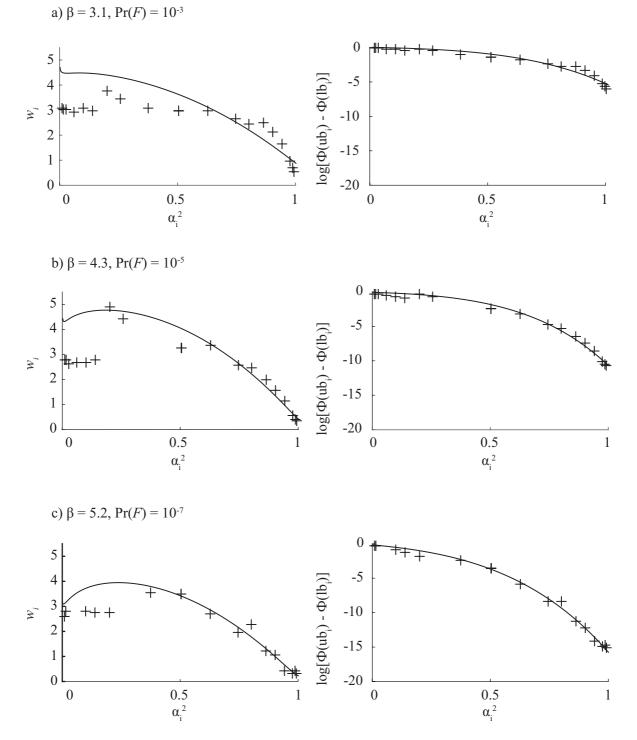




Figure 11. Optimization results for 10 two-dimensional, linear problems in standard normal space, which are discretized with 10 intervals per dimension. The prior failure probabilities are  $10^{-3}(\beta = 3.1)$ ,  $10^{-5}(\beta = 4.3)$ and  $10^{-7}(\beta = 5.2)$ . The crosses represent the optimization results. The solid lines are the fitted parametric functions (Eq. 14). The left-hand side shows the relation between the width of a discretization frame  $w_i$  and  $|\alpha_i|$ and the right-hand side shows the relation between the probability mass enclosed by the discretization frame with width  $w_i$  and  $|\alpha_i|$ .

## 455 **3.3** Summary of the proposed procedure

- 456 The steps of the proposed procedure are:
- 457 1. Formulate the reliability problem

458	2.	Set up the corresponding BN
459	3.	Perform a FORM analysis for the reliability problem
460	4.	Simplify the BN by removing nodes based on:
461		a. their importance for prediction
462		b. their observability
463		c. whether or not a node simplifies modeling of dependencies
464		d. whether or not it is desired to explicitly show a node in the BN for
465		communication purposes
466	5.	Find the discretization scheme in U-space based on the proposed heuristics i.e.:
467		a. the discretization scheme is centered at the design point from the FORM
468		analysis
469		b. the same number of intervals is used for each random variable
470		c. the width of the discretization frame follows Eq. 14
471	6.	Transform the discretization scheme to X-space
472	7.	Compute the CPTs of the component state node and the basic random variables using
473		Monte Carlo simulation or Latin hypercube sampling

A MATLAB-based software tool performing these steps is available for download under
 www.era.bgu.tum.de/software.

# 476 4 Applications

## 477 4.1 Verification example I

For verification purposes, we apply the proposed methodology to the discretization of a
general limit state with non-normal dependent random variables. The approximation error
made by this discretization is investigated for different measurement outcomes.

481 Failure is defined through the LSF  $g(\mathbf{x})$ :

$$g(\mathbf{x}) = a - \prod_{i=1}^{n} X_i \tag{15}$$

482 i.e., failure corresponds to the event  $\{\prod_{i=1}^{n} X_i \ge a\}$ .

The basic random variables are distributed as  $X_1 \sim LN(0,0.5)$  and  $X_2, ..., X_n \sim LN(1,0.3)$ (values in parenthesis are the parameters of the lognormal distribution). The statistical dependence among the  $X_i$  is described through a Gaussian copula model, with pairwise correlation coefficients  $\rho_{ij}$ . The parameters *a* and  $\rho_{ij}$  determine the prior failure probability  $P_F$ . Measurements  $M_i = m_i$  are available for all basic random variables; they are associated with multiplicative measurement errors  $\varepsilon_i \sim LN(0,0.71)$ . In Tabs. 2 and 3, different cases with 3 and 4 random variables are shown. These cases differ with respect to the prior failure 490 probability  $P_F$ , the correlation between the random variables  $\rho_{ij}$  and the observed 491 measurements **m**. For each case, a reference solutions  $P_{F|\mathbf{M}}$  is calculated analytically.

492

Table 2. Evaluation of the discretization error for different measurement outcomes **m**, for problems with n = 3random variables. *a* is the constant in the LSF, Eq. 15;  $\rho_{ij}$  is the correlation coefficient between  $X_i$  and  $X_j$  for all  $i \neq j$ ;  $P_F$  and  $P_{F|M}$  denote the analytically calculated prior and posterior failure probabilities;  $\hat{P}_{F|M}$  is the conditional failure probability calculated with the discrete BN.

а	c <sub>up</sub>	$ ho_{ij}$	P <sub>F</sub>	m	P <sub>F <i>M</i></sub>	$\widehat{P}_{F M}$		Relative error [%]
100	10 <sup>3</sup>	0	3.6 <i>E</i> — 5	[3.0,2.9,2,9]	4.3 <i>E</i> – 5	4.5 <i>E</i> — 5	3 <i>E</i> – 6	6
100	10 <sup>3</sup>	0	3.6 <i>E</i> – 5	[2.3,1.1,2,1]	4.6E - 6	5.3 <i>E</i> – 6	7 <i>E</i> — 7	14
100	10 <sup>3</sup>	0	3.6 <i>E</i> – 5	[0.9,2.4,0.9]	2.8 <i>E</i> – 7	3.5 <i>E</i> – 7	7 <i>E</i> – 8	25
200	15 <sup>3</sup>	0.5	1.6 <i>E</i> – 4	[1.6,2.0,1.2]	1.4E - 6	1.4 <i>E</i> – 6	1 <i>E</i> - 7	4
400	8 <sup>3</sup>	0.5	6.4 <i>E</i> – 6	[2.6,3.0,3.2]	8.2 <i>E</i> – 7	8.9 <i>E —</i> 7	7E - 8	9
400	12 <sup>3</sup>	0.5	6.4 <i>E</i> – 6	[3.6,3.3,4,3]	4.9 <i>E</i> – 6	5.0E - 6	1 <i>E</i> – 9	3

<sup>497</sup> 

Table 3. Evaluation of the discretization error for different measurement outcomes **m**. The number of random variables n = 4; *a* is the constant in the LSF, Eq. 15;  $\rho_{ij}$  is the correlation coefficient between  $X_i$  and  $X_j$  for all  $i \neq j$ ;  $P_F$  and  $P_{F|M}$  denote the analytically calculated prior and posterior failure probabilities;  $\hat{P}_{F|M}$  is the conditional failure probability calculated with the discrete BN.

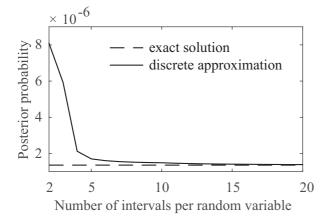
	а	C <sub>up</sub>	$ ho_{ij}$	$P_{\rm F}$	m	P <sub>F <i>M</i></sub>	₽ <sub>F <i>M</i></sub>	Absolute	Relative
_								error	error [%]
	400	10 <sup>4</sup>	0	1.7 <i>E</i> — 5	[2.2,3.2,2.4,3.4]	9.5 <i>E</i> — 6	1.0 <i>E</i> – 5	9 <i>E</i> — 7	9
	400	10 <sup>4</sup>	0	1.7 <i>E</i> — 5	[1.6,1.6,1.6,2.0]	6.5 <i>E</i> — 7	7.9E - 7	1 <i>E</i> – 7	21
	400	104	0	1.7 <i>E</i> — 5	[1.1,2.3,1.9,1.2]	2.4 <i>E</i> — 7	3.0 <i>E</i> - 7	6E - 8	26
	600	104	0.5	1.3 <i>E</i> – 3	[3.3,1.7,2.8,2.6]	4.2 <i>E</i> – 4	4.3 <i>E</i> – 4	2 <i>E</i> — 5	4
-	800	84	0.5	5.3 <i>E</i> – 4	[1.9,2.0,1.9,2.4]	1.8 <i>E</i> – 5	1.9 <i>E</i> – 5	1 <i>E</i> – 6	8

502

The results in Tables 2 and 3 show that the proposed methodology for discretization leads to errors in the posterior probability estimate that are acceptably small for most engineering applications. (It is reminded that discretization does not lead to a discretization error in the prior case.) As expected, the relative error is larger when the posterior probability is low, and the absolute error is larger when the posterior probability is high. This follows from the error measure defined in Eq. 11, which balances the relative with the absolute error. In addition, the results do not display any apparent effect of correlation on the accuracy. 510 To assess the effect of the choice of the number of discretization intervals, the failure

511 probability  $\hat{P}_{F|M}$  was calculated for a discretization scheme with up to 20 intervals per RV for 512 the fourth measurement case in Tab. 2. The estimated failure probabilities  $\hat{P}_{F|M}$  are plotted

513 together with the exact solution in Fig. 12.



515 Figure 12. Posterior probability  $\hat{P}_{F|M}$  as a function of the number of intervals per random variable together with 516 the exact (analytical) solution  $P_{F|M}$  for the fourth measurement case [1.6, 2.0, 1.2] in Tab 2.

#### 517 **4.2 Verification example II**

*i*=1

514

518 The failure criterion applied in verification example I (Eq. 15) leads to a linear LSF in U-

space. To verify the accuracy of the proposed method for problems with non-linear LSFs inU-space, we additionally investigate the following LSF:

$$g(\mathbf{x}) = a - \sum_{i=1}^{n} X_i \tag{16}$$

Again the basic random variables  $X_1$  to  $X_n$  are distributed as  $X_1 \sim LN(0,0.5)$ 521 and  $X_2, ..., X_n \sim LN(1, 0.3)$ . Different cases with n = 2, 3 and 4 random variables are 522 523 investigated. Measurements  $M_i = m_i$  are available for all basic random variables; associated 524 to these measurement are multiplicative measurement errors  $\varepsilon_i \sim LN(0,0.71)$ . For 525 independent random variables  $X_i$  it is possible to determine posterior distributions  $f_{X_i|M_i}(x_i|m_i)$  analytically. The posterior failure probabilities  $P_{F|M}$ , which are used as 526 reference solutions, are calculated through importance sampling with 107 samples. The 527 528 results are presented in Tab. 4.

Table 4. Evaluation of the discretization error for different measurement outcomes **m**. The problems have n = 2, 3 or 4 random variables; *a* is the constant in the LSF, Eq. 16;  $\rho_{ij}$  is the correlation coefficient between  $X_i$ and  $X_j$  for all  $i \neq j$ ;  $P_F$  and  $P_{F|M}$  denote the prior respectively posterior failure probabilities, which are calculated through importance sampling with 10<sup>7</sup> samples;  $\hat{P}_{F|M}$  is the conditional failure probability calculated with the discrete BN. Since for correlated basic random variables there is no analytical solution, in these cases the updating of the basic random variables was performed through rejection sampling with > 5*E*7 accepted samples. 535

а	C <sub>up</sub>	$ ho_{ij}$	P <sub>F</sub>	m	P <sub>F <i>M</i></sub>	₽ <sub><b>F</b> <i>M</i></sub>	Absolute error	Rel. error [%]
<u>n =</u>	2:							
12	10 <sup>2</sup>	0	1.3 <i>E</i> — 5	[2.8,4.5]	1.4E - 5	1.2E - 5	2E - 6	15
12	10 <sup>2</sup>	0	1.3E - 5	[2.3,2.4]	3.3E - 6	3.5E - 6	2 <i>E</i> – 7	6
10	12 <sup>2</sup>	0	1.7E - 4	[4.0,3.2]	4.0E - 4	3.7E - 4	3 <i>E</i> – 5	7
12	10 <sup>2</sup>	0.5	1.7E - 4	[2.3,2.4]	4.8E - 5	5.0E - 5	2E - 6	4
<u>n =</u>	3:							
15	10 <sup>3</sup>	0	3.7E - 5	[2.1,5.6,5.0]	4.8E - 5	4.5E - 5	4E - 6	7
15	10 <sup>3</sup>	0	3.7 <i>E</i> – 5	[1.1,3.7,3.4]	1.5 <i>E</i> – 5	1.8E - 5	3E - 6	20
13	12 <sup>3</sup>	0	5.0E - 4	[3.0,3.0,3.0]	5.4E - 4	5.4 <i>E</i> – 5	3 <i>E</i> – 6	1
16	10 <sup>3</sup>	0.5	9.1E - 4	[3.0,6.0,5.0]	1.8E - 3	1.9 <i>E</i> – 3	3 <i>E</i> – 5	2
<u>n =</u>	4:							
20	84	0	7.4E - 6	[2.0,4.0,3.4,3.0]	4.9E - 6	5.6E - 6	6E - 7	13
17	84	0	3.1E - 4	[1.0,1.4,1.2,2.0]	4.5E - 5	5.2E - 5	7 <i>E</i> – 6	15
17	12 <sup>4</sup>	0	3.1E - 4	[3.1,2.0,3.3,2.4]	2.5E - 4	2.5E - 4	7E - 6	3
24	84	0.5	3.7 <i>E</i> – 4	[1.0,1.4,1.2,2.0]	1.1E - 5	1.1E - 5	9E — 7	8

<sup>536</sup> 

537 The results in Tab. 4 do not differ substantially from Tabs. 2 and 3. This indicates that the 538 (weak) non-linearity of the LSF function describing failure does not affect the accuracy 539 significantly.

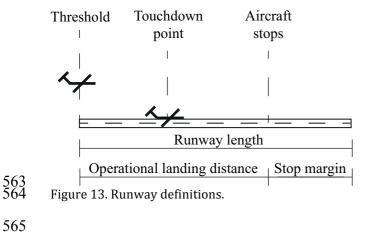
### 540 4.3 Runway overrun

541 Runway overrun (RWO) of a landing aircraft is one of the most critical accidents types in 542 civil aviation (IATA, 2013). A conceptual RWO warning system is developed with the

- 543 proposed discretization procedure. It provides RWO probabilities conditional on observations 544 of the landing-weight, the headwind and the approach speed for different aircraft types and 545 different airports. For a detailed description of how this problem can be treated in BN 546 framework we refer to (Zwirglmaier and Straub, 2015).
- 547 RWO is the event of the operational landing distance exceeding the available runway length
- 548 (Fig. 13). Correspondingly, a LSF for runway overrun can be defined as:

$$g(\mathbf{X}) = \text{Runway length} - \text{Operational landing distance}(\mathbf{X})$$
(17)

- 549 with **X** representing the basic random variables of the problem.
- 550 (Drees and Holzapfel, 2012) proposed a model for the operational landing distance required
- by a landing aircraft, which is applied here. The model, as well as the basic random variables
- 552 **X**, are presented in (Zwirglmaier et al., 2014), which also includes a detailed description of
- 553 the reliability and sensitivity analysis.
- 554 We consider two different airports (AP I and AP II) and two different aircraft types (AC A
- and AC B). While the aircraft type affects the landing-weight, the airport affects both the
- be headwind and the approach speed. The distribution models for landing-weight, headwind and
- 557 approach speed deviation at the different airports and with the different aircraft types are
- 558 given in Tabs. 5–7. All other basic random variables of the problem are not affected by the
- airport and aircraft type and are as in (Zwirglmaier et al., 2014).
- 560 Tab. 8 summarizes the FORM importance measures of all random variables **X** computed for
- 561 the four combinations of aircrafts and airports.
- 562



566

7 Table 5. Distribution models for landing weight conditional on the aircraft.

Landing weight [t]						
Aircraft	Distribution	Mean	Std. deviation			
А	Weibull (min)	59.25	1.69			
В	Weibull (min)	64.25	1.69			

569 Table 6. Distribution models for head wind conditional on the airport.

	Head wind [kts]						
Airport	Distribution	Mean	Std. deviation				
Ι	Normal	5.42	5.75				
II	Normal	6.51	5.75				

Table 7. Distribution models for approach speed deviation conditional on the airport.

Approach speed deviation [kts]							
Airport	Distribution	Mean	Std. deviation				
Ι	Gumbel (max)	4.69	4.21				
II	Gumbel (max)	5.63	4.21				

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574 Table 8. FORM importance measures  $\alpha_i$  for each aircraft-airport combination and every basic random variable 575 in the RWO application.

	$lpha_i$				
Random variable	(I/A)	(I/B)	(II/A)	(II/B)	Annotation
Landing weight [t]	0.09	0.10	0.11	0.09	Modeled
Headwind [kts]	-0.65	-0.61	-0.67	-0.60	Modeled
Temperature [°C]	0.03	-0.00	-0.03	-0.03	Not important
Air pressure [hPa]	0.01	-0.01	-0.01	-0.00	Not important
Touchdown point [m]	0.20	0.16	0.18	0.20	Modeled
Approach speed deviation [kts]	0.20	0.21	0.20	0.24	Not observable
Time of spoiler deployment [s]	-0.00	-0.00	0.01	0.01	Not important
Time of breaking initiation [s]	0.70	0.74	0.68	0.73	Not observable
Time of reverser deployment [s]	0.03	0.04	0.06	0.05	Not important
Time of breaking end [s]	-0.02	-0.01	0.01	0.02	Not important

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<sup>570</sup> 571 572

#### 576 4.3.1 Selection of relevant random variables

- 577 The applied RWO model includes 10 basic random variables. However, it is sufficient to 578 include only a selection of these explicitly in the BN. Random variables that are not relevant 579 for the prediction of RWO in the considered scenarios can be excluded. This is the case for 580 random variables with a low FORM importance, whose value does not depend significantly 581 on airport and aircraft type. Here, all random variables, whose absolute value of the FORM 582 importance measure  $|\alpha_i|'s$  is smaller than 0.1, are excluded (see Tab. 8). The one exception is
- 583 landing weight, since its mean value is substantially influenced by the aircraft type.
- 584 One can additionally exclude random variables that cannot be measured before the decision 585 on whether to land or not is made. This holds for Touchdown point and the time at which the 586 pilot initiates breaking. Since these basic random variables are also not needed to simplify the 587 modeling of dependencies, it is not necessary to explicitly model them in the BN, as indicated 588 in Tab. 8.
- 589 4.3.2 BN model

590 The resulting BN of the RWO warning system is shown in Fig. 14. During the aircraft 591 approach, measurements can be obtained for the three basic random variables included in the 592 BN.

The random variables are discretized separately for each aircraft-airport combination (joint states of discrete parents) with 8 intervals each, following the proposed discretization procedure. In a second step, the discretization schemes are merged, i.e. the regions of the outcome space, which are discretized with fine intervals for at least one of the aircraft-airport combinations, are discretized with the respective fine intervals also in the merged discretization scheme. In the end 15 (landing-weight), 10 (headwind) and 9 (approach speed deviation) intervals are used to discretize the three basic random variables.

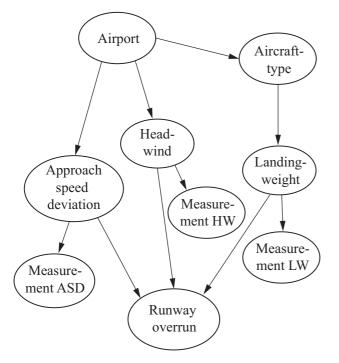
For all observable quantities, the measurements  $m_i$  are modeled with an additive observation error:

$$m_i = x_i + \varepsilon_i \tag{18}$$

602  $\varepsilon_i$  is modeled by a normal distribution with zero mean and standard deviation  $\sigma_{\varepsilon_i}$ .

For the random variable landing weight (at landing time), the standard deviation of the measurement error is  $\sigma_{\varepsilon_{LW}} = 0.34$  tons. Due to turbulences governing wind speeds, the measurement of the head wind speed at the time of the measurement is only an uncertain indicator for the head wind speed at landing time; we model the measurement error with a standard deviation  $\sigma_{\varepsilon_{HW}} = 2.88$  kts. The measurement uncertainty associated with the approach speed deviation at landing has standard deviation  $\sigma_{\varepsilon_{ASD}} = 4.21$  kts.

49 (Measurement LW), 57 (Measurement HW) and 57 (Measurement ASD) intervals are usedto discretize the measurement nodes.





- 613 Figure 14. BN structure for a RWO warning system.
- 614 4.3.3 Results

615 In Tab. 9, RWO probabilities for the different airports and aircrafts obtained with the discrete

616 BN are compared to solutions, which were calculated by importance sampling around the 617 design point.

618 Table 9. RWO probabilities for the different airports and aircrafts calculated with the discrete BN  $p_{_{BN'}}$ 619 together with solutions calculated by importance sampling around the design point  $p_{DS}$ . The latter have a

620 621 sampling error with coefficient of variation in the order of 10%.

AP/AC	$p_{BN}$	$p_{DS}$
I/A	2.0 <i>e</i> – 7	1.9 <i>e</i> – 7
I/B	1.0 <i>e</i> – 6	9.2 <i>e</i> – 7
II/A	1.3 <i>e</i> – 7	1.3 <i>e</i> – 7
II/B	6.9 <i>e</i> – 7	6.5 <i>e</i> – 7

622 In Tab. 10, results obtained with the BN for different hypothetical cases of aircrafts 623 approaching an airport are presented. In each of these cases, measurements associated with 624 landing weight, headwind and the approach speed deviation are made. A threshold on the probability of RWO is used to decide, whether or not the pilot should continue landing or 625 626 cancel the landing attempt. Here we assume that up to a RWO probability of  $10^{-6}$  the pilot 627 should continue landing.

628 Table 10. Probabilities of RWO and corresponding decision on landing, computed with the BN for different 629 sets of observations.

Case	Airport	Aircraft	Meas. LW	Meas. HW	Meas. ASD [kts]	Pr (RWO)	Landing
a)	Ι	В	63	0	10.5	2.5 <i>e</i> – 8	Yes
b)	Ι	А	61	-10	5	4.8 <i>e</i> – 6	No
c)	II	В	67	3	0	6.5 <i>e</i> – 10	Yes
d)	II	А	57.5	-12	3	1.3 <i>e</i> – 6	No

## 630 5 Discussion

631 When modeling with BNs, it is often necessary or beneficial to discretize continuous random 632 variables. When the BN includes rare events that are a function of such random variables, the 633 choice of the discretization scheme is non-trivial. In this contribution, we investigate this 634 discretization based on FORM concepts, and propose a heuristic procedure for an efficient 635 discretization in these cases. This is based on importance measures  $\alpha_i$  obtained through a 636 FORM analysis, which represent the influence of the uncertainty associated with a random 637 variable  $X_i$ .

638 The most important finding is that discretization should focus on the area around the most 639 likely failure point (design point), identified by a FORM analysis. Furthermore, we find that 640 optimally all random variables should be discretized with approximately equal numbers of 641 intervals, independent of their importance, as long as  $|\alpha_i|$  is not close to zero. The widths of 642 the intervals should be selected based on the FORM importance  $\alpha_i$  of the random variables. 643 With increasing importance, the interval width should be reduced, leading to finer 644 discretization for larger  $|\alpha_i|$ . This relation is particularly evident for  $|\alpha_i| \ge 0.8$ . We show that 645 it is possible to fit a parametric function to approximate the relation between  $|\alpha_i|$  and the 646 optimal width of the region on which the discretization should focus.

- 647 This parametric function is used to derive a heuristic procedure for finding an efficient 648 discretization. This allows the extrapolation of the optimization results to problems with more 649 random variables. As demonstrated by the verification examples, the heuristic procedure leads 650 to accurate results.
- This paper is restricted to static discretization. Application of the proposed procedure within dynamic discretization (e.g. (Neil et al., 2008)) should be investigated. The results of the
- 653 procedure can serve as an initial discretization scheme, which is iteratively adjusted within
- 654 dynamic discretization. This might strongly enhance the convergence performance of these
- algorithms.
- 656 The gain in computational efficiency resulting from the proposed procedure over alternative
- 657 static discretization approaches is problem specific. Some insights can be gained from Figure
- 8. A discretization with intervals of equal width centered at the origin would require between
- approximately 2 to 5 times more intervals per random variable to achieve the same accuracy.

660 Because of the exponential increase of computational effort with number of random variables, 661 this leads to a considerable increase in efficiency. The gain compared to discretization with 662 equal-frequency intervals is expected to be even higher, since equal frequency intervals 663 focuses the fine intervals on the region of high probability density rather than on the tails of 664 the distribution.

665 This paper focuses on component reliability problems, which are characterized by a single 666 design point. Nevertheless the heuristics derived can also be applied to system reliability 667 problems. System reliability problems can in general be treated as combinations of 668 component reliability problems. Parallel and series systems are to be distinguished. For parallel systems discretization should be performed based on the joint design point of the 669 670 problem. For series systems, following the same line of thought as in the runway overrun 671 example, discretization can be performed separately for each component problem 672 (corresponds to the discrete cases i.e. airport- aircraft combinations in the RWO example). In 673 a second step the discretization schemes can be merged. In the same way it is possible to 674 apply the heuristic to multi state components. One can treat each limit state surface (LSF) 675 defining the boundary between two states separately and merge the discretization schemes 676 afterwards.

The number of basic random variables in a single LSF that can be modeled explicitly in a BN is limited to around 5 to 8. This is due to the exponential growth of the target nodes CPT with increasing number of parents and is independent of the discretization method. Despite this limitation, BNs are applicable to many practical problems – particularly if one considers that usually not all basic random variables need to be modeled explicitly as nodes, as demonstrated in the presented example.

683 While in this paper the focus was on the discretization of the basic random variables, it is 684 straightforward to incorporate the BNs discussed into larger models.

# 685 6 Conclusion

686 We investigate discretization of continuous reliability problems such that they can be treated 687 in a discrete Bayesian network framework. Reliability problems with linear LSF in standard 688 normal space are considered. These can be seen as FORM approximations of reliability 689 problems. For these linear LSFs, optimal discretization schemes are found, which are optimal 690 with respect to an error measure calculated through a preposterior analysis. Since FORM is 691 known to give good approximations also for most non-linear reliability problem, the resulting 692 discretization schemes are efficient also for non-linear LSFs. The main findings presented in 693 this paper are:

An optimal discretization scheme should discretize finely the area around the FORM design point.

- The size of the sub-region of the outcome space of a random variable  $X_i$  can be reduced significantly for random variables whose corresponding uncertainty is dominating the reliability problem
- The number of intervals used for discretization should be approximately equal for all
   basic random variables
- 701 On this basis, we propose a heuristic that can be used to find an efficient discretization
- scheme. In verification examples, this heuristic is found to give good accuracy and efficiency.

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