

# Error-informed parallel adaptive Kriging method for time-dependent reliability analysis

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## Abstract

Active learning single-loop Kriging methods have gained significant attention for time-dependent reliability analysis. However, it still remains a challenge to estimate the time-dependent failure probability efficiently and accurately in practical engineering problems. This study proposes a new method, called ‘Error-informed Parallel Adaptive Kriging’ (EPAK) for efficient time-dependent reliability analysis. First, a sequential variance-amplified importance sampling technique is developed to estimate the time-dependent failure probability based on the trained global response Kriging model of the true performance function. Then, the maximum relative error of the time-dependent failure probability is derived to facilitate the construction of stopping criterion. Finally, a parallel sampling strategy is proposed through combining the relative error contribution and an influence function, which enables parallel computing and reduces the unnecessary limit state function evaluations caused by excessive clustering. The superior performance of the proposed method is validated through several examples. Numerical results demonstrate that the method can accurately estimate the time-dependent failure probability with higher efficiency than several compared methods.

*Keywords:* Time-dependent reliability analysis; Active learning; Kriging model; Importance sampling; Parallel computing; Estimation error

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## 1. Introduction

Reliability analysis aims to assess the likelihood that a structural system or component will consistently perform the intended functions when considering multi-source uncertainties, such as material properties, applied loads, geometry, model uncertainty, and others [1]. The traditional time-invariant reliability analysis neglects the time-dependent factors and is limited to assessing the reliability at a specific time instant. Considering the fact that the performance of engineered component or systems usually degrades with the increase of service time, time-dependent reliability analysis (TRA) has drawn much attention in recent decades [2–4]. Incorporating the time dimension adds complexity to the problem, which makes the TRA more time-consuming than the time-invariant cases [5, 6]. The current TRA methods could be categorized into three following groups: (1) out-crossing rate methods; (2) composite limit state methods; (3) extreme value methods.

In the out-crossing rate methods, the time-dependent failure probability (TDFP) is approximated by integrating the instantaneous out-crossing rate over a specified time interval. The origin of this type of methods can be tracked into the 1940s when Rice introduced the famous Rice formula [7], laying the theoretical foundation for the development of the out-crossing rate methods for time-dependent reliability problems. The out-crossing rate methods can be further classified into the two following groups. The first group consists of the numerical methods, mainly based on the FORM or the method of moments. The representative methods include PHI2 [8], PHI2+ [9], and MPHI2 [10], etc. The second is the analytical methods, including but not limited to [11–13]. Although the performance of the out-crossing rate methods have been improved in the last several decades [14, 15], the large computation cost and the inherent assumption still restrict the applicability of the out-crossing rate methods in TRA.

The composite limit state methods discretize the time-dependent limit state function (LSF) into a series of instantaneous LSFs, thereby transforming the time-dependent issue into a time-invariant one with the series system reliability concept [16]. Some studies use FORM to calculate the instantaneous reliability, including but not limited to [17–19]. These methods may produce inaccurate results when the LSF is highly nonlinear or contains multiple most probable points (MPPs). Simulation-based methods have also

53 been developed for TRA, e.g., subset simulation [20], line sampling [21] and importance sampling [22],  
54 etc. Despite better accuracy, the simulation-based methods still suffer from low efficiency in engineering  
55 practices.

56 The extreme value methods transform the time-dependent problem into a time-invariant one, and the  
57 TDFP is estimated by solving the extreme value distribution [23, 24]. Recent advancements in artificial  
58 intelligence have accelerated the application of machine learning in predicting the extreme value distribu-  
59 tion, where adaptive surrogate models have gained significant attention for their effective balance between  
60 accuracy and efficiency [25–27]. The extreme response surrogate-based methods, as a type of double-loop  
61 methods, need to identify the extreme response in the inner loop and build a surrogate model for the ex-  
62 treme response in the outer loop [28]. The typical methods falling into this category include the parallel  
63 efficient global optimization [29], confidence-based adaptive extreme response surface method [30], impor-  
64 tance sampling-based double-loop Kriging [31], mixed EGO method [32] and so forth. The double-loop  
65 methods may suffer from low accuracy due to the fact the accuracy of searching extreme time instant would  
66 influence the accuracy of surrogate model. Besides, this kind of method requires a large amount LSF evalu-  
67 ations for the problems with stochastic process with a long time interval. Instead of a double-loop scheme,  
68 a single-loop scheme involving constructing the global response surrogate models has been extensively in-  
69 vestigated [33, 34]. Among the various kinds of surrogates, the Kriging model is particularly prominent for  
70 its capability to interpolate and provide a local measure of prediction uncertainty. In this regard, the most  
71 pioneering is the single-loop Kriging surrogate modeling (SILK) method [33]. Some other representative  
72 single-loop methods include variance reduction-guided adaptive Kriging (VARAK) method [35], real-time  
73 estimation error-guided active learning Kriging (REAL) method [36], single-loop Gaussian process regres-  
74 sion based-active learning (SL-GPR-AL) method [37], and several others [38, 39]. In the aforementioned  
75 single-loop methods, the estimation of the TDFP is all based on Monte Carlo Simulation (MCS) and is  
76 computationally challenging for small failure probability problems. To solve this problem, several meth-  
77 ods have been developed by combining the single-loop Kriging model with importance sampling technique  
78 [40, 41] and subset simulation [42–45], respectively. Recently, the first author and his co-authors [46] have

79 extended the Bayesian active learning originally developed for time-invariant reliability analysis [47–49]  
80 to the time-dependent counterpart, and proposed uncertainty-aware adaptive Bayesian inference combined  
81 with hyper-ring decomposition importance sampling for TRA. As mentioned in [36], the estimation error of  
82 TDFP is an important measure for assessing whether the TDFP is sufficiently accurate as the final result  
83 throughout the active learning process. To the best of authors’ knowledge, however, none of existing studies  
84 have attempted to quantify and reduce the estimation error of TDFP provided by Kriging model and im-  
85 portance sampling. Besides, these single-loop methods can only identify one point per iteration, hindering  
86 their availability of the parallel computing.

87 This study aims to propose a novel method termed ‘Error-informed Parallel Adaptive Kriging’ (EPAK)  
88 for efficient TRA. The primary contributions can be outlined as follows:

89 • The variance-amplified importance sampling (VAIS) proposed in [48] is adapted in a sequential way  
90 for estimating the small TDFPs. The resulting sequential VAIS can reduce the sample size and total  
91 computation time but also avoid the computer memory issue due to the one-shot Kriging prediction on the  
92 large amount of samples;

93 • The maximum relative error of the TDFP is derived under the combination of the single-loop Kriging  
94 model and VAIS. This allows the quantification of error in estimating TDFP, and facilitates the construction  
95 of an effective stopping criterion. In this study, the adaptive updating of Kriging model is terminated by  
96 judging the maximum relative error;

97 • A parallel sampling strategy is developed through combining the relative error contribution and an  
98 influence function that considers the correlation between the existing training points and the candidate  
99 points. This strategy can select multiple training points and overcome the problem of unnecessary LSF  
100 evaluations caused by excessive clustering.

101 The rest of this study is structured as follows. Section 2 introduces the estimation of TDFP based on  
102 Kriging and MCS. In section 3, the proposed EPAK method is presented in detail. Four examples are  
103 studied in Section 4 to validate the proposed method. Section 5 concludes the present study.

104 **2. Background of time-dependent reliability analysis**

105 In this section, we first give the definition of TDFP. The MCS-based TDFP estimation is then reviewed.  
 106 The Kriging-based global response surrogate method is finally introduced.

107 *2.1. Definition of time-dependent failure probability*

108 The key to TRA is to calculate the failure probability (denoted as  $P_f(0, t_e)$ ) of a structural system  
 109 or component within a predefined time interval  $[0, t_e]$ . A failure event is defined when the LSF is below  
 110 zero at any time instant within  $[0, t_e]$ . Let  $g(\mathbf{X}, \mathbf{Y}(t), t)$  denote the LSF with an  $n$ -dimensional input  
 111 vector of random variables  $\mathbf{X} = [X_1, X_2, \dots, X_n]$  and an  $m$ -dimensional input vector of stochastic processes  
 112  $\mathbf{Y}(t) = [Y_1(t), Y_2(t), \dots, Y_m(t)]$ , where  $t$  denotes the time parameter.

113 The TDFP  $P_f(0, t_e)$  is expressed as follows:

$$P_f(0, t_e) = \mathbb{P} \{g(\mathbf{X}, \mathbf{Y}(t), t) < 0, \exists t \in [0, t_e]\} \quad (1)$$

114 where  $\mathbb{P}$  denotes the operation of probability.

115 Assuming that the stochastic processes  $\mathbf{Y}(t)$  are represented by a function of a  $d_0$ -dimensional vector of  
 116 random variables  $\Xi$  and a time parameter  $t$ , the TDFP can be expressed as an integral given by:

$$P_f(0, t_e) = \int_{\mathbb{R}^n} \int_{\mathbb{R}^{d_0}} I(\mathbf{x}, \mathbf{y}(\xi, t), t) f_{\mathbf{X}}(\mathbf{x}) f_{\Xi}(\xi) d\mathbf{x} d\xi \quad (2)$$

117 where  $I(\mathbf{x}, \mathbf{y}(\xi, t), t)$  is the time-dependent indicator function;  $\mathbf{x}$  and  $\xi$  are the realizations of  $\mathbf{X}$  and  $\Xi$ ,  
 118 respectively;  $f_{\mathbf{X}}(\mathbf{x})$  and  $f_{\Xi}(\xi)$  are the joint probability density functions (PDFs) of the random variables  
 119  $\mathbf{X}$  and  $\xi$ , respectively;  $I(\mathbf{x}, \mathbf{y}(\xi, t), t)$  is written as:

$$I(\mathbf{x}, \mathbf{y}(\xi, t), t) = \begin{cases} 1, & g(\mathbf{x}, \mathbf{y}(\xi, t), t) < 0, \exists t \in [0, t_e] \\ 0, & \text{otherwise} \end{cases} \quad (3)$$

120 *2.2. Discretization of stochastic processes*

121 The stochastic processes  $\mathbf{Y}(t)$  are discretized into random variables for computation purposes. The  
 122 commonly used expansion optimal linear estimation (EOLE) [50] is adopted in this study due to its high  
 123 efficiency and accuracy. One should note that it is not straightforward to simulate non-Gaussian processes  
 124 with EOLE method. For general stochastic processes, some advanced simulation methods can be used, e.g.,  
 125 [51, 52]. In this study, only Gaussian processes are considered for convenience.  $n_t$  time instants are employed  
 126 to discretize the time interval  $[0, t_e]$ . Considering a Gaussian process  $Y(t)$  for the sake of illustration,  $Y(t)$   
 127 is written as:

$$Y(t) \approx \mu(t) + \sum_{i=1}^p \frac{\xi_i}{\sqrt{\lambda_i}} \phi_i^\top \boldsymbol{\rho}_Y(t) \quad (4)$$

128 where  $\mu(t)$  denotes the mean function;  $p$  denotes the number of dominated eigenvectors, which can be  
 129 determined according to [8].  $\xi_i$  ( $i = 1, \dots, p$ ) denote the expanded random variables; For Gaussian process  
 130  $Y(t)$  here,  $\xi_i$  ( $i = 1, \dots, p$ ) are the standard normal variables;  $\lambda_i$  and  $\phi_i$  represent the dominated eigenvalues  
 131 and eigenvectors, respectively.  $\boldsymbol{\rho}_Y(t) = [\sigma(t) \sigma(t_1) \rho(t, t_1), \dots, \sigma(t) \sigma(t_{n_t}) \rho(t, t_{n_t})]^\top$  denotes the vector of  
 132 covariance function;  $\sigma(t)$  denotes the standard deviation function;  $\rho_Y(t_i, t_j)$  is the autocorrelation function.

133 *2.3. TDFP estimation by MCS*

134 After the stochastic processes  $\mathbf{Y}(t)$  are discretized, the LSF is expressed as  $g(\mathbf{x}, \mathbf{y}((\boldsymbol{\xi}_1, \boldsymbol{\xi}_2, \dots, \boldsymbol{\xi}_m), t), t)$ ,  
 135 where  $\boldsymbol{\xi}_i$  ( $i = 1, \dots, m$ ) denote the vectors of random variables. Based on MCS, the TDFP in Eq. (2) is  
 136 estimated as:

$$\tilde{P}_f(0, t_e) = \frac{1}{N_{mcs}} \sum_{i=1}^{N_{mcs}} I_t(\mathbf{x}^{(i)}, (\boldsymbol{\xi}_1^{(i)}, \boldsymbol{\xi}_2^{(i)}, \dots, \boldsymbol{\xi}_m^{(i)})) \quad (5)$$

137 where  $N_{mcs}$  denotes the number of samples; The indicator function  $I_t(\mathbf{x}^{(i)}, (\boldsymbol{\xi}_1^{(i)}, \boldsymbol{\xi}_2^{(i)}, \dots, \boldsymbol{\xi}_m^{(i)}))$  is expressed  
 138 as:

$$I_t(\mathbf{x}^{(i)}, (\boldsymbol{\xi}_1^{(i)}, \boldsymbol{\xi}_2^{(i)}, \dots, \boldsymbol{\xi}_m^{(i)})) = \begin{cases} 1, & \text{if } g(\mathbf{x}^{(i)}, \mathbf{y}((\boldsymbol{\xi}_1^{(i)}, \boldsymbol{\xi}_2^{(i)}, \dots, \boldsymbol{\xi}_m^{(i)}), t_j), t_j) < 0, \exists j = 1, \dots, n_t \\ 0, & \text{otherwise} \end{cases} \quad (6)$$

139 The coefficient of variation (COV) of  $\tilde{P}_f(0, t_e)$  is written as:

$$\text{COV}(\tilde{P}_f(0, t_e)) = \sqrt{\frac{1 - \tilde{P}_f(0, t_e)}{(N_{mcs} - 1) \times \tilde{P}_f(0, t_e)}} \quad (7)$$

140 It should be noted that the MCS involves a double loop computation procedure. That is, the realizations  
 141  $\{\mathbf{x}^{(i)}, (\boldsymbol{\xi}_1^{(i)}, \boldsymbol{\xi}_2^{(i)}, \dots, \boldsymbol{\xi}_m^{(i)})\}_{i=1}^{N_{mcs}}$  are first generated in the outer loop. For any realization, the LSF is evaluated  
 142 at  $n_t$  time instants in the inner loop, i.e.,  $\{g(\mathbf{x}^{(i)}, \mathbf{y}((\boldsymbol{\xi}_1^{(i)}, \boldsymbol{\xi}_2^{(i)}, \dots, \boldsymbol{\xi}_m^{(i)}), t_j), t_j)\}_{j=1}^{n_t}$ . If the minimum  
 143 response is less than zero (i.e.,  $\min\left(\{g(\mathbf{x}^{(i)}, \mathbf{y}((\boldsymbol{\xi}_1^{(i)}, \boldsymbol{\xi}_2^{(i)}, \dots, \boldsymbol{\xi}_m^{(i)}), t_j), t_j)\}_{j=1}^{n_t}\right) < 0$ ), the realization is  
 144 regarded to be failed; otherwise, it is considered safe. A schematic representation of the MCS is given in  
 145 Fig. 1, where the failed time trajectories are denoted by the red lines. The TDFP is calculated by dividing  
 146 the number of failed time trajectories by the total number of time trajectories.

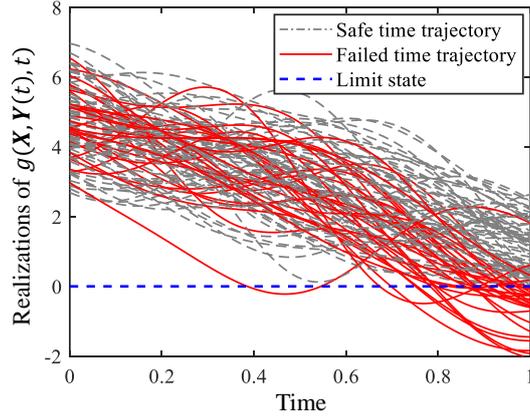


Figure 1: Illustrative diagram of MCS.

#### 147 2.4. Kriging-based global response surrogate method

148 The MCS-based TDFP estimation requires  $N_{mcs} \times n_t$  LSF evaluations, making it prohibitive in many  
 149 practical engineering problems. To address this problem, the Kriging model is adopted to build a global  
 150 response surrogate model for the LSF, enabling efficient estimation of the TDFP. The details of Kriging  
 151 model is presented in [Appendix A](#). The stochastic processes  $\mathbf{Y}(t)$  are expressed using  $\sum_{i=1}^m p^{(i)}$  random  
 152 variables. Hence, the input dimension of LSF is equal to  $n + \sum_{i=1}^m p^{(i)} + 1$ . In order to avoid dealing with  
 153 high dimensions, the stochastic processes are directly used as inputs of the Kriging model, instead of the

154 expanded random variables. The input dimension is thus reduced to  $n + m + 1$ . The transformation of input  
 155 can be given as:

$$\begin{bmatrix} \mathbf{x}^{(1)} & (\boldsymbol{\xi}_1^{(1)}, \dots, \boldsymbol{\xi}_m^{(1)}) & t^{(1)} \\ \mathbf{x}^{(2)} & (\boldsymbol{\xi}_1^{(2)}, \dots, \boldsymbol{\xi}_m^{(2)}) & t^{(2)} \\ \vdots & \vdots & \vdots \\ \mathbf{x}^{(n_0)} & (\boldsymbol{\xi}_1^{(n_0)}, \dots, \boldsymbol{\xi}_m^{(n_0)}) & t^{(n_0)} \end{bmatrix} \rightarrow \text{Eq. (4)} \rightarrow \begin{bmatrix} \mathbf{x}^{(1)} & \mathbf{y}_t^{(1)} & t^{(1)} \\ \mathbf{x}^{(2)} & \mathbf{y}_t^{(2)} & t^{(2)} \\ \vdots & \vdots & \vdots \\ \mathbf{x}^{(n_0)} & \mathbf{y}_t^{(n_0)} & t^{(n_0)} \end{bmatrix} \quad (8)$$

156 where  $n_0$  is the number of training points.

157 The adaptive Kriging based TRA methods starts with constructing a rough Kriging surrogate model  
 158 with a small number of initial training points. Then, new informative training points are sequentially  
 159 selected through a learning function and the Kriging model is updated. The procedure is terminated when  
 160 a predefined stopping criterion is fulfilled. Finally, the TDFP is estimated as:

$$\hat{P}_f(0, t_e) = \frac{1}{N_{mcs}} \sum_{i=1}^{N_{mcs}} \hat{I}_t(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}) \quad (9)$$

161 where  $\hat{I}_t(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)})$  is denoted as:

$$\hat{I}_t(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}) = \begin{cases} 1, & \text{if } \mu_{\hat{g}}(\mathbf{x}^{(i)}, \mathbf{y}_{t_j}^{(i)}, t_j) < 0, \exists j = 1, \dots, n_t \\ 0, & \text{otherwise} \end{cases} \quad (10)$$

162 where  $\mu_{\hat{g}}(\mathbf{x}^{(i)}, \mathbf{y}_{t_j}^{(i)}, t_j)$  is the mean prediction of the Kriging model.

### 163 3. Error-informed Parallel Adaptive Kriging

164 This section proposes a new method called EPAK, which can estimate small TDFPs and enable parallel  
 165 computing. First, the VAIS is adapted in a sequential way to reduce the sample size and computational  
 166 cost. Later, the maximum relative error of TDFP is derived under the combination of Kriging model and

167 VAIS. Finally, a stopping criterion and a parallel sampling strategy are developed to adaptively enrich the  
 168 training point set.

### 169 3.1. Sequential variance-amplified importance sampling

170 As mentioned in the last section, MCS involves a double loop computation procedure and requires  
 171  $N_{mcs} \times n_t$  Kriging model predictions. For problems with low TDFPs, a large  $N_{mcs}$  should be specified to  
 172 ensure the estimation accuracy, rendering the computation cumbersome. To address this problem, the VAIS  
 173 developed in [48] is introduced and adapted in a sequential way to replace the MCS in this study, so as to  
 174 reduce the sample size and total computation time.

175 The TDFP  $P_f(0, t_e)$  in Eq. (2) is rewritten as:

$$P_f(0, t_e) = \int_{\mathbb{R}^n} \int_{\mathbb{R}^{d_e}} I(\mathbf{x}, \mathbf{y}(\boldsymbol{\xi}, t), t) \frac{f_{\mathbf{X}}(\mathbf{x})}{h_{\mathbf{X}}(\mathbf{x})} h_{\mathbf{X}}(\mathbf{x}) f_{\Xi}(\boldsymbol{\xi}) d\mathbf{x} d\boldsymbol{\xi} \quad (11)$$

176 where  $d_e = \sum_{i=1}^m p^{(i)}$ ;  $h_{\mathbf{X}}(\mathbf{x}) = f_{\mathbf{X}}(\mathbf{x}; \mathbf{m}_{\mathbf{X}}, \gamma \cdot \boldsymbol{\sigma}_{\mathbf{X}})$  denotes the importance sampling density (ISD), which  
 177 is established by enlarging the vector of standard deviations  $\boldsymbol{\sigma}_{\mathbf{X}}$  (or enlarging the vector of variances  $\boldsymbol{\sigma}_{\mathbf{X}}^2$ )  
 178 of the PDF  $f_{\mathbf{X}}(\mathbf{x})$  (maintain the means  $\mathbf{m}_{\mathbf{X}}$  unchanged), where  $\gamma$  is the amplification factor. Note that  
 179 the stochastic processes are typically represented by many random variables, and amplifying the standard  
 180 deviations of these random variables greatly increases the computational complexity. For simplicity, only  
 181 the standard deviations of input random variables in LSF are amplified.

182 Then,  $\Delta N$  samples are generated from  $h_{\mathbf{X}}(\mathbf{x})$  and  $f_{\Xi}(\boldsymbol{\xi})$ .  $P_f(0, t_e)$  in Eq. (11) can be estimated as:

$$\hat{P}_f(0, t_e) = \frac{1}{\Delta N} \sum_{i=1}^{\Delta N} \hat{I}_t(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}) \frac{f_{\mathbf{X}}(\mathbf{x}^{(i)})}{h_{\mathbf{X}}(\mathbf{x}^{(i)})} \quad (12)$$

183 The variance and COV of  $\hat{P}_f(0, t_e)$  are given as:

$$\mathbb{V}[\hat{P}_f(0, t_e)] = \frac{1}{\Delta N - 1} \left( \frac{1}{\Delta N} \sum_{i=1}^{\Delta N} \hat{I}_t(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}) \left( \frac{f_{\mathbf{X}}(\mathbf{x}^{(i)})}{h_{\mathbf{X}}(\mathbf{x}^{(i)})} \right)^2 - \left( \hat{P}_f(0, t_e) \right)^2 \right) \quad (13)$$

184

$$\text{COV} \left[ \hat{P}_f(0, t_e) \right] = \frac{\sqrt{\mathbb{V} \left[ \hat{P}_f(0, t_e) \right]}}{\hat{P}_f(0, t_e)} \quad (14)$$

185 The samples are generated sequentially from the ISD and  $f_{\Xi}(\boldsymbol{\xi})$ , and then predicted on the Kriging  
 186 model, which can greatly save the computation time. First,  $\Delta N$  samples are generated. Let the number of  
 187 iteration  $s = 1$  and the total number of samples  $N_0 = s \times \Delta N$ . The TDFP is estimated by Eq. (12) and  
 188 expressed as  $\hat{P}_f^{(s)}$ . A quantity  $\varpi^{(s)}$  is introduced to efficiently store the Kriging prediction information for  
 189  $\Delta N$  samples, minimizing memory usage while enabling the calculation of the variance estimator.  $\varpi^{(s)}$  is  
 190 written as:

$$\varpi^{(s)} = \frac{1}{\Delta N} \sum_{i=1}^{\Delta N} \hat{I}_t \left( \mathbf{x}^{(i)}, \mathbf{y}_t^{(i)} \right) \left( \frac{f_{\mathbf{X}} \left( \mathbf{x}^{(i)} \right)}{h_{\mathbf{X}} \left( \mathbf{x}^{(i)} \right)} \right)^2 \quad (15)$$

191 Additional  $\Delta N$  samples are generated and let  $s = s + 1$ .  $\hat{P}_f^{(s)}$  and  $\varpi^{(s)}$  are estimated by Eq. (12) and  
 192 Eq. (15), respectively. The TDFP and its variance can be re-estimated as:

$$\hat{P}_f(0, t_e) = \frac{1}{s} \sum_{i=1}^s \hat{P}_f^{(s)} \quad (16)$$

193

$$\mathbb{V} \left[ \hat{P}_f(0, t_e) \right] = \frac{1}{N_0 - 1} \left( \frac{1}{s} \sum_{i=1}^s \varpi^{(s)} - \left( \hat{P}_f(0, t_e) \right)^2 \right) \quad (17)$$

194 The sampling process is executed until the COV of the TDFP is lower than the target threshold, i.e.,  
 195  $\text{COV} \left[ \hat{P}_f(0, t_e) \right] < \epsilon_p$ .

### 196 3.2. Relative error of TDFP

197 According to Eq. (10),  $\hat{I}_t \left( \mathbf{x}^{(i)}, \mathbf{y}_t^{(i)} \right)$  is estimated based on judging the sign of  $\mu_{\hat{g}} \left( \mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}, t \right)$ , which  
 198 is predicted by Kriging and may be wrongly estimated. The relative error of the predicted TDFP  $\hat{P}_f(0, t_e)$   
 199 with respect to the true result  $P_f(0, t_e)$  can be defined as:

$$\delta = \left| \frac{P_f(0, t_e) - \hat{P}_f(0, t_e)}{P_f(0, t_e)} \right| \quad (18)$$

200 The true result  $P_f(0, t_e)$  is expressed as:

$$P_f(0, t_e) = \frac{1}{N_0} \left[ \sum_{i=1}^{N_0} \hat{I}_t(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}) \frac{f_{\mathbf{X}}(\mathbf{x}^{(i)})}{h_{\mathbf{X}}(\mathbf{x}^{(i)})} + \sum_{h=1}^{\hat{N}_s^w} \frac{f_{\mathbf{X}}(\mathbf{x}^{(h)})}{h_{\mathbf{X}}(\mathbf{x}^{(h)})} - \sum_{k=1}^{\hat{N}_f^w} \frac{f_{\mathbf{X}}(\mathbf{x}^{(k)})}{h_{\mathbf{X}}(\mathbf{x}^{(k)})} \right] \quad (19)$$

201 where  $\hat{N}_s^w$  denotes the total number of time trajectories predicted to be safe by Kriging model but misclas-  
 202 sified;  $\hat{N}_f^w$  denotes the total number of time trajectories predicted to be failed but misclassified. Due to the  
 203 fact that the true number of misclassified time trajectories is unknown, the last two terms in Eq. (19) are  
 204 uncertain. Let  $\sum_{h=1}^{\hat{N}_s^w} \frac{f_{\mathbf{X}}(\mathbf{x}^{(h)})}{h_{\mathbf{X}}(\mathbf{x}^{(h)})} = \mathcal{N}_s$  and  $\sum_{k=1}^{\hat{N}_f^w} \frac{f_{\mathbf{X}}(\mathbf{x}^{(k)})}{h_{\mathbf{X}}(\mathbf{x}^{(k)})} = \mathcal{N}_f$ , the relative error  $\delta$  can thus be written as:

$$\delta = \left| 1 - \frac{\sum_{i=1}^{N_0} \hat{I}_t(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}) \frac{f_{\mathbf{X}}(\mathbf{x}^{(i)})}{h_{\mathbf{X}}(\mathbf{x}^{(i)})}}{\sum_{i=1}^{N_0} \hat{I}_t(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}) \frac{f_{\mathbf{X}}(\mathbf{x}^{(i)})}{h_{\mathbf{X}}(\mathbf{x}^{(i)})} + \mathcal{N}_s - \mathcal{N}_f} \right| \quad (20)$$

205 Although the exact values of  $\mathcal{N}_s$  and  $\mathcal{N}_f$  are unknown, it is possible to obtain the expectation and variance  
 206 of the two quantities. To achieve this goal,  $\hat{I}_t^s(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) = 1$  is first introduced to denote that the time  
 207 trajectory predicted to be safe by Kriging is actually in a failed status. Correspondingly,  $\hat{I}_t^s(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) = 0$   
 208 denotes that the time trajectory predicted to be safe is correctly classified.  $\mathcal{N}_s$  can thus be written as:

$$\mathcal{N}_s = \sum_{h=1}^{\hat{N}_s^w} \frac{f_{\mathbf{X}}(\mathbf{x}^{(h)})}{h_{\mathbf{X}}(\mathbf{x}^{(h)})} = \sum_{h=1}^{N_s} \hat{I}_t^s(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) \frac{f_{\mathbf{X}}(\mathbf{x}^{(h)})}{h_{\mathbf{X}}(\mathbf{x}^{(h)})} \quad (21)$$

209 where  $N_s$  is the total number of safe time trajectory predicted by Kriging.

210 The expectation and variance of  $\mathcal{N}_s$  can be expressed as:

$$\mathbb{E}[\mathcal{N}_s] = \mathbb{E} \left[ \sum_{h=1}^{N_s} \hat{I}_t^s(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) \frac{f_{\mathbf{X}}(\mathbf{x}^{(h)})}{h_{\mathbf{X}}(\mathbf{x}^{(h)})} \right] = \sum_{h=1}^{N_s} \mathbb{E} \left[ \hat{I}_t^s(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) \right] \frac{f_{\mathbf{X}}(\mathbf{x}^{(h)})}{h_{\mathbf{X}}(\mathbf{x}^{(h)})} \quad (22)$$

$$\mathbb{V}[\mathcal{N}_s] = \mathbb{V} \left[ \sum_{h=1}^{N_s} \hat{I}_t^s(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) \frac{f_{\mathbf{X}}(\mathbf{x}^{(h)})}{h_{\mathbf{X}}(\mathbf{x}^{(h)})} \right] = \sum_{h=1}^{N_s} \mathbb{V} \left[ \hat{I}_t^s(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) \right] \left( \frac{f_{\mathbf{X}}(\mathbf{x}^{(h)})}{h_{\mathbf{X}}(\mathbf{x}^{(h)})} \right)^2 \quad (23)$$

212 where the probability of  $\hat{I}_t^s(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) = 0$ , i.e., the probability of correct sign estimate of  $(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)})$ , can

213 be expressed as:

$$p_{s,c}(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) = \mathbb{P} \left\{ \hat{I}_t^s(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) = 0 \right\} = \mathbb{P} \left\{ \bigcap_{h=1}^{n_t} \hat{g}(\mathbf{x}^{(h)}, \mathbf{y}_{t_j}^{(h)}, t_j) > 0 \right\} \quad (24)$$

214 where the computation of  $p_{s,c}(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)})$  is quite cumbersome in practice. Instead of calculating the  
 215 exact value of the  $p_{s,c}(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)})$ , it is possible to obtain its sub-optimal estimate without sacrificing the  
 216 estimation accuracy according to our recent study [46]. The sub-optimal estimate  $\bar{p}_{s,c}(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)})$  can be  
 217 expressed as [46]:

$$\bar{p}_{s,c}(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) = \min_{j=1, \dots, n_t} \Phi \left( \frac{\mu_{\hat{g}}(\mathbf{x}^{(h)}, \mathbf{y}_{t_j}^{(h)}, t_j)}{\sigma_{\hat{g}}(\mathbf{x}^{(h)}, \mathbf{y}_{t_j}^{(h)}, t_j)} \right) \quad (25)$$

218 The sub-optimal estimate of misclassification probability is written as:

$$\bar{p}_{s,w}(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) = \mathbb{P} \left\{ \hat{I}_t^s(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) = 1 \right\} = 1 - \min_{j=1, \dots, n_t} \Phi \left( \frac{\mu_{\hat{g}}(\mathbf{x}^{(h)}, \mathbf{y}_{t_j}^{(h)}, t_j)}{\sigma_{\hat{g}}(\mathbf{x}^{(h)}, \mathbf{y}_{t_j}^{(h)}, t_j)} \right) \quad (26)$$

219 It is easy to find that  $\hat{I}_t^s(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)})$  follows the Bernoulli distribution. The expectation and variance of  
 220  $\mathcal{N}_s$  are rewritten as:

$$\mathbb{E}[\mathcal{N}_s] = \sum_{h=1}^{N_s} \bar{p}_{s,w}(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) \frac{f_{\mathbf{X}}(\mathbf{x}^{(h)})}{h_{\mathbf{X}}(\mathbf{x}^{(h)})} \quad (27)$$

$$\mathbb{V}[\mathcal{N}_s] = \sum_{h=1}^{N_s} \bar{p}_{s,w}(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)}) (1 - \bar{p}_{s,w}(\mathbf{x}^{(h)}, \mathbf{y}_t^{(h)})) \left( \frac{f_{\mathbf{X}}(\mathbf{x}^{(h)})}{h_{\mathbf{X}}(\mathbf{x}^{(h)})} \right)^2 \quad (28)$$

222 Similarly, let  $\hat{I}_t^f(\mathbf{x}^{(k)}, \mathbf{y}_t^{(k)}) = 1$  denote that the time trajectory predicted to be failed by Kriging is  
 223 actually in a safe status. Correspondingly,  $\hat{I}_t^s(\mathbf{x}^{(k)}, \mathbf{y}_t^{(k)}) = 0$  denotes the time trajectory predicted to be  
 224 failed is correctly classified.

225 The misclassification probability can be written as:

$$p_{f,w}(\mathbf{x}^{(k)}, \mathbf{y}_t^{(k)}) = \mathbb{P} \left\{ \hat{I}_t^f(\mathbf{x}^{(k)}, \mathbf{y}_t^{(k)}) = 1 \right\} = \mathbb{P} \left\{ \bigcap_{j=1}^{n_t} \hat{g}(\mathbf{x}^{(k)}, \mathbf{y}_{t_j}^{(k)}, t_j) > 0 \right\} \quad (29)$$

226 According to the aforementioned discussion, the corresponding sub-optimal estimate of  $p_{f,w}(\mathbf{x}^{(k)}, \mathbf{y}_t^{(k)})$

227 can be defined as:

$$\bar{p}_{f,w}(\mathbf{x}^{(k)}, \mathbf{y}_t^{(k)}) = \min_{j=1, \dots, n_t} \Phi \left( \frac{\mu_{\hat{g}}(\mathbf{x}^{(k)}, \mathbf{y}_{t_j}^{(k)}, t_j)}{\sigma_{\hat{g}}(\mathbf{x}^{(k)}, \mathbf{y}_{t_j}^{(k)}, t_j)} \right) \quad (30)$$

228 The expectation and variance of  $\mathcal{N}_f$  can be written as:

$$\mathbb{E}[\mathcal{N}_f] = \sum_{k=1}^{N_f} \bar{p}_{f,w}(\mathbf{x}^{(k)}, \mathbf{y}_t^{(k)}) \frac{f_{\mathbf{X}}(\mathbf{x}^{(k)})}{h_{\mathbf{X}}(\mathbf{x}^{(k)})} \quad (31)$$

$$\mathbb{V}[\mathcal{N}_f] = \sum_{k=1}^{N_f} \bar{p}_{f,w}(\mathbf{x}^{(k)}, \mathbf{y}_t^{(k)}) (1 - \bar{p}_{f,w}(\mathbf{x}^{(k)}, \mathbf{y}_t^{(k)})) \left( \frac{f_{\mathbf{X}}(\mathbf{x}^{(k)})}{h_{\mathbf{X}}(\mathbf{x}^{(k)})} \right)^2 \quad (32)$$

230 where  $N_f$  is the total number of failed points predicted by Kriging model.

231 For any time trajectory predicted to be either safe or failed, the sub-optimal estimate of the misclassifi-  
232 cation probability is expressed as:

$$\bar{p}_w(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}) = \Phi \left( - \left| \min_{j=1, \dots, n_t} \frac{\mu_{\hat{g}}(\mathbf{x}^{(i)}, \mathbf{y}_{t_j}^{(i)}, t_j)}{\sigma_{\hat{g}}(\mathbf{x}^{(i)}, \mathbf{y}_{t_j}^{(i)}, t_j)} \right| \right) \quad (33)$$

233 With the VAIS technique presented in Section 3.1, the dispersedly distributed samples are generated as  
234 the candidate samples. It is reasonable to consider that the number of safe time trajectories  $N_s$  and the  
235 number of failed time trajectories  $N_f$  are both large enough. In this case, the confidence intervals of  $\mathcal{N}_s$   
236 and  $\mathcal{N}_f$  can be approximately obtained using the central limit theorem. Besides,  $\mathcal{N}_s$  and  $\mathcal{N}_f$  approximately  
237 follow normal distributions, which are expressed as:

$$\mathcal{N}_s \sim N(\mathbb{E}[\mathcal{N}_s], \mathbb{V}[\mathcal{N}_s]) \quad (34)$$

$$\mathcal{N}_f \sim N(\mathbb{E}[\mathcal{N}_f], \mathbb{V}[\mathcal{N}_f]) \quad (35)$$

239 The confidence intervals of  $\mathcal{N}_s$  and  $\mathcal{N}_f$  can be obtained as:

$$\mathcal{N}_s \in [\mathcal{N}_s^l, \mathcal{N}_s^u] = [\Phi_{\mathcal{N}_s}^{-1}(\alpha/2), \Phi_{\mathcal{N}_s}^{-1}(1 - \alpha/2)] \quad (36)$$

240

$$\mathcal{N}_f \in [\mathcal{N}_f^l, \mathcal{N}_f^u] = [\Phi_{\mathcal{N}_f}^{-1}(\alpha/2), \Phi_{\mathcal{N}_f}^{-1}(1 - \alpha/2)] \quad (37)$$

241 where  $\Phi_{\mathcal{N}_s}^{-1}(\cdot)$  and  $\Phi_{\mathcal{N}_f}^{-1}(\cdot)$  denote inverse cumulative distribution functions (CDFs);  $\alpha$  is the confidence level  
 242 ( $\alpha = 0.05$  is used in this study).

243 The maximum relative error of TDFP can thus be obtained by:

$$\delta_{\max} = \max \left( \left| \frac{\mathcal{N}_s^l - \mathcal{N}_f^u}{\sum_{i=1}^{N_0} \hat{I}_t(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}) \frac{f_{\mathbf{x}}(\mathbf{x}^{(i)})}{h_{\mathbf{x}}(\mathbf{x}^{(i)})} + \mathcal{N}_s^l - \mathcal{N}_f^u} \right|, \left| \frac{\mathcal{N}_s^u - \mathcal{N}_f^l}{\sum_{i=1}^{N_0} \hat{I}_t(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}) \frac{f_{\mathbf{x}}(\mathbf{x}^{(i)})}{h_{\mathbf{x}}(\mathbf{x}^{(i)})} + \mathcal{N}_s^u - \mathcal{N}_f^l} \right| \right) \quad (38)$$

244 It should be noted that the quantification of the relative error of TDFP can be regarded as an extension  
 245 of the study in static reliability analysis [53] and the one in time-dependent reliability analysis [36]. The  
 246 significant difference is that the estimation of relative error is based on the VAIS in this study, whereas it is  
 247 based on MCS in the two existing studies.

### 248 3.3. Stopping criterion and parallel sampling strategy

249 In adaptive Kriging based TRA method, a suitable stopping criterion is required to determine whether  
 250 the obtained estimate of TDFP is accurate enough as the final result. Based on the Eq. (38), the stopping  
 251 criterion can be defined by judging whether the maximum relative error of TDFP is below a prescribed  
 252 threshold:

$$\delta_{\max} \leq \epsilon_r \quad (39)$$

253 where  $\epsilon_r$  is the specified threshold. Note that the adaptive updating of Kriging model is terminated only  
 254 when Eq. (39) is satisfied twice consecutively to prevent the potential fake convergence.

255 If the stopping criterion is not met, new training points need to be identified and evaluated on the true  
 256 LSF to update the Kriging model. In order to efficiently decrease the maximum relative error of TDFP,

257 the number of misclassified time trajectories should be as small as possible. In other words, the greater the  
 258 misclassification of a time trajectory, the greater the contribution to reducing the relative error of TDFP.  
 259 That is, the  $\bar{p}_w(\mathbf{x}, \mathbf{y}_t)$  in Eq. (33) measures the contribution of the time trajectory to the relative error  
 260 of TDFP. Therefore, the relative error can be minimized as much as possible through evaluating the point  
 261 that maximizes the misclassification probability in Eq. (33). However, directly selecting the point with the  
 262 highest misclassification probability overlooks the distance between the newly added training point and the  
 263 existing ones, which may cause excessive clustering and lead to unnecessary LSF evaluations. To address  
 264 this problem, this paper introduces a influence function by considering the correlation between the existing  
 265 training points and all candidate points as the distance measure, and then proposes a new learning function  
 266 called improved relative error contribution (IREC):

$$\text{IREC}(\mathbf{x}, \mathbf{y}_t) = IF(\mathbf{x}; \mathbf{x}^{(1)}, \mathbf{x}^{(2)}, \dots, \mathbf{x}^{(n_0)}) \times \bar{p}_w(\mathbf{x}, \mathbf{y}_t) \quad (40)$$

267 where  $IF(\mathbf{x}; \mathbf{x}^{(1)}, \mathbf{x}^{(2)}, \dots, \mathbf{x}^{(n_0)})$  is the introduced influence function and denoted as [54]:

$$IF(\mathbf{x}; \mathbf{x}^{(1)}, \mathbf{x}^{(2)}, \dots, \mathbf{x}^{(n_0)}) = \prod_{i=1}^{n_0} [1 - R(\mathbf{x}, \mathbf{x}^{(i)})] \quad (41)$$

268 where  $\mathbf{x}^{(1)}, \mathbf{x}^{(2)}, \dots, \mathbf{x}^{(n_0)}$  are the  $n_0$  existing training points;  $R(\cdot)$  is the Gaussian correlation function of  
 269 the Kriging model in this study.

270 The IREC function can be further extended to parallel sampling case by exploiting the advantages of  
 271 the introduced influence function. In an active learning iteration,  $q$  training points are sequentially selected  
 272 and simultaneously enriched to the dataset  $\mathcal{D}$ , which is elaborated as follows. First, the time trajectory  
 273 maximizing the IREC function is selected and then the time instant with the largest prediction uncertainty is  
 274 chosen. Herein, the commonly used U function is used to determine the optimal time instant [55]. Therefore,

275 the selection strategy of the first added training point  $\mathbf{v}_{new}^{(n_0+1)}$  is expressed as follows:

$$\mathbf{v}_{new}^{(n_0+1)} = [\mathbf{x}^{(i^*)}, \mathbf{y}_{t_{j^*}}^{(i^*)}, t_{j^*}]$$

$$i^* = \arg \max_{i=1, \dots, N_0} IF(\mathbf{x}; \mathbf{x}^{(1)}, \mathbf{x}^{(2)}, \dots, \mathbf{x}^{(n_0)}) \bar{p}_w(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}), j^* = \arg \min_{j=1, \dots, n_t} \frac{|\mu_{\hat{g}}(\mathbf{x}^{(i^*)}, \mathbf{y}_{t_{j^*}}^{(i^*)}, t_{j^*})|}{\sigma_{\hat{g}}(\mathbf{x}^{(i^*)}, \mathbf{y}_{t_{j^*}}^{(i^*)}, t_{j^*})} \quad (42)$$

276 Sequentially, after  $q - 1$  updated training points have been identified, the  $q$ -th training point can be  
 277 obtained as:

$$\mathbf{v}_{new}^{(n_0+q)} = [\mathbf{x}^{(i^*)}, \mathbf{y}_{t_{j^*}}^{(i^*)}, t_{j^*}]$$

$$i^* = \arg \max_{i=1, \dots, N_0} IF(\mathbf{x}; \mathbf{x}^{(1)}, \mathbf{x}^{(2)}, \dots, \mathbf{x}^{(n_0+q-1)}) \bar{p}_w(\mathbf{x}^{(i)}, \mathbf{y}_t^{(i)}), j^* = \arg \min_{j=1, \dots, n_t} \frac{|\mu_{\hat{g}}(\mathbf{x}^{(i^*)}, \mathbf{y}_{t_{j^*}}^{(i^*)}, t_{j^*})|}{\sigma_{\hat{g}}(\mathbf{x}^{(i^*)}, \mathbf{y}_{t_{j^*}}^{(i^*)}, t_{j^*})} \quad (43)$$

278 It is evident from Eqs. (42)-(43) that  $q$  training points are sequentially identified through considering  
 279 the relative error contribution and the correlation between all training points and candidate points. After  
 280  $q$  training points are selected within an iteration, the Kriging model is updated. Note that the developed  
 281 parallel sampling strategy may be similar to the idea that multiple informative points are sequentially gener-  
 282 ated by maximizing the product of the learning function and influence function, which is recently introduced  
 283 in the field of time-invariant reliability analysis [56, 57] and time-dependent system reliability analysis [58].  
 284 The difference lies mainly in two aspects. First, this study develops a new IREC learning function based on  
 285 the concept of minimizing the relative error of TDFP. Second, the several existing researches have only con-  
 286 sidered the correlation between the current and previously selected points within an iteration, ignoring the  
 287 correlation of the training points prior to the current iteration, whereas this study addresses this problem.

### 288 3.4. Implementation of the proposed method

289 The flowchart of the proposed method is shown in Fig. 2. The implementation procedures are summa-  
 290 rized as follows:

291 **Step 1:** Discretize the time interval  $[0, t_e]$  and the stochastic processes  $\mathbf{Y}(t)$  (if involved).

292 **Step 2:** Generate  $n_0$  initial training points by Sobol sequence. The sampling domain of these initial  
 293 points is set to  $[\boldsymbol{\mu} - 3\boldsymbol{\sigma}, \boldsymbol{\mu} + 3\boldsymbol{\sigma}]$ , where  $\boldsymbol{\mu}$  and  $\boldsymbol{\sigma}$  are the mean and standard deviation of the input random

294 variables, respectively. The corresponding responses are calculated by evaluating the LSF  $g(\cdot)$ . Establish  
 295 the initial dataset  $\mathcal{D} = \{[\mathbf{x}^{(i)}, \mathbf{y}_{t_i}^{(i)}, t_i], g^{(i)}\}_{i=1}^{n_0}$  with the EOLE method. Let the number of LSF evaluations  
 296  $N_{eva} = n_0$  and the number of iterations  $N_{ite} = 1$ .

297 **Step 3:** Construct the candidate sample pool  $\mathcal{S} = [\mathbf{x}^{(i)}, (\boldsymbol{\xi}_1^{(i)}, \dots, \boldsymbol{\xi}_m^{(i)})]_{i=1}^{\Delta N}$ , where  $\Delta N$  is the number of  
 298 initial candidate samples. Let  $s = 1$  and the the number of total candidate samples  $N_0 = s \times \Delta N$ .

299 **Step 4:** Build the Kriging model based on the dataset  $\mathcal{D}$ .

300 **Step 5:** Estimate the TDFP  $\hat{P}_f(0, t_e)$  and its variance  $\mathbb{V}[\hat{P}_f(0, t_e)]$  based on Eqs. (16)-(17). Calculate  
 301 the maximum relative error  $\delta_{\max}$  in Eq. (38).

302 **Step 6:** If the stopping criterion in Eq. (39) is fulfilled twice consecutively, turn to Step 8; else, turn to  
 303 Step 7.

304 **Step 7:** Identify  $q$  points  $\mathcal{D}_+ = [\mathbf{x}^{(i)}, \mathbf{y}_{t_i}^{(i)}, t_i]_{i=1}^q$  using the developed parallel sampling strategy. Evaluate  
 305 the LSF on the  $q$  selected points to obtain the responses  $\{g^{(i)}\}_{i=1}^q$ . Enrich the points and responses into the  
 306 dataset  $\mathcal{D}$ . Let  $N_{eva} = N_{eva} + q$  and  $N_{ite} = N_{ite} + 1$ , and return to Step 4.

307 **Step 8:** Check if the COV of the TDFP is below the target threshold, i.e,  $\text{COV}[\hat{P}_f(0, t_e)] < \epsilon_p$ . If  
 308 satisfied, go to Step 9; else, generate additional  $\Delta N$  candidate samples and enrich the candidate sample  
 309 pool  $\mathcal{S}$ , let  $s = s + 1$  and return to Step 5.

310 **Step 9:** Return the estimated TDFP  $\hat{P}_f(0, t_e)$  and end the proposed method.

## 311 4. Examples and results

312 The effectiveness of the proposed EPAK method is validated by comparing it with several existing non-  
 313 parallel TRA methods. The two compared methods, SILK [33] and REAL [36], are implemented in all  
 314 examples and the number of initial training points is set to be 12. Other implementation details follow the  
 315 default settings in the original studies. The results for the remaining compared methods, including eSPT  
 316 [59], STRA [60], SLK-co-SS [45], etc., are taken from the respective publications unless otherwise stated.  
 317 The parameter settings for the proposed method are as follows:  $n_0 = 12$ ,  $\gamma = 1.5$ ,  $\Delta N = 10^4$ ,  $\epsilon_p = 5\%$ ,  
 318  $\epsilon_r = 5\%$ . Different values are specified for  $q$  to investigate the impact on the results. The reported results are

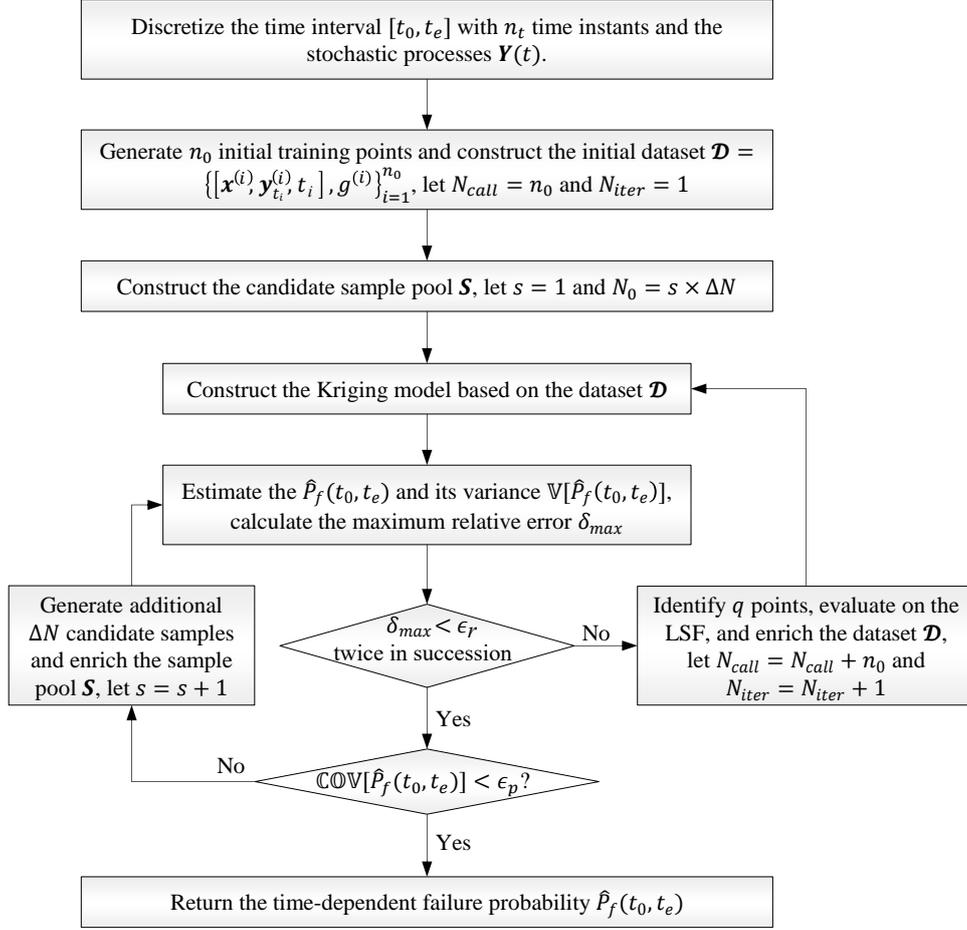


Figure 2: Flowchart of the proposed method.

319 averaged over 10 independent runs in MATLAB R2019(b) under the environment (CPU of Intel i5-13400F,  
 320 RAM of 16 GB) unless otherwise stated.

#### 321 4.1. Mathematical example

322 A mathematical model is investigated in this section [45, 59, 60]. The LSF is expressed as:

$$g(\mathbf{X}, Y(t), t) = X_1^2 X_2 + (X_2 + 1)t^2 - 5X_1(1 + Y(t))t - 20 \quad (44)$$

323 where the input variables  $X_1$  and  $X_2$  both follow a normal distribution with the mean of 3.5 and the standard  
 324 deviation of 0.25;  $Y(t)$  denotes a Gaussian process, where the mean and standard deviation are 0 and 1,

325 respectively. The autocorrelation function of the Gaussian process is defined as  $\kappa(t_1, t_2) = \exp[-(t_1 - t_2)^2]$ ;  
 326  $t \in [0, 1]$  represents the time parameter.

327 The time interval and Gaussian process  $Y(t)$  are discretized with 101 time instants and three independent  
 328 standard normal variables, respectively. Fig. 3 depicts one hundred realizations of  $Y(t)$ . The TDFP results  
 329 provided by different methods are listed in Table 1. The TDFP  $\hat{P}_f = 0.3079$  by MCS is considered as  
 330 the reference result. It is observed that all methods can produce results close to the reference result. In  
 331 terms of efficiency, the proposed method requires fewer iterations with the help of the developed parallel  
 332 sampling strategy. When  $q = 1$  (indicating that the parallel computing is unavailable), the proposed method  
 333 costs less LSF evaluations than other compared methods though it produces slightly larger COVs. Through  
 334 comparing the computation time between different methods, it can be found that the proposed method  
 335 typically requires less CPU time than other active learning TRA methods.

Table 1: TRA results of example 4.1.

Methods	$N_{ite}$	$N_{eva}$	$\hat{P}_f$	$\epsilon_{\hat{P}_f}$	$COV[\hat{P}_f]$	CPU Time (s)	
MCS	-	$101 \times 10^6$	0.3079	-	0.15%	1.9	
eSPT <sup>1</sup>	-	139	0.2986	3.02%	-	-	
STRA <sup>2</sup>	16.4	27.4	0.3041	1.23%	-	-	
SLK-co-SS <sup>3</sup>	35.8	46.8	0.3072	0.23%	-	-	
SILK	12.5	23.5	0.3075	0.13%	2.95%	31.1	
REAL	10.8	21.8	0.3070	0.29%	3.56%	29.5	
Proposed method	$q = 1$	6.6	17.6	0.3059	0.65%	6.58%	2.5
	$q = 2$	4.3	18.6	0.2995	2.73%	7.02%	2.0
	$q = 3$	4.1	21.3	0.3014	2.11%	5.16%	2.2
	$q = 4$	3.7	22.8	0.3062	0.55%	4.16%	2.2
	$q = 5$	3.8	23.2	0.3031	1.56%	4.95%	2.3
	$q = 6$	3.5	27.0	0.3093	0.45%	3.14%	2.6
	$q = 7$	2.8	24.6	0.3086	0.23%	2.57%	2.2
	$q = 8$	2.8	26.4	0.3081	0.07%	3.51%	2.4

<sup>1</sup> The results are taken from research [59];

<sup>2</sup> The results are taken from research [60];

<sup>3</sup> The results are taken from research [45].

#### 336 4.2. Corroded simply supported beam

337 In this section, we consider the TRA of a corroded beam [60] shown in Fig 4. The span of the beam  
 338 is 5m. The bending loads include the concentrated load  $F(t)$  and the uniformly distributed load  $p =$

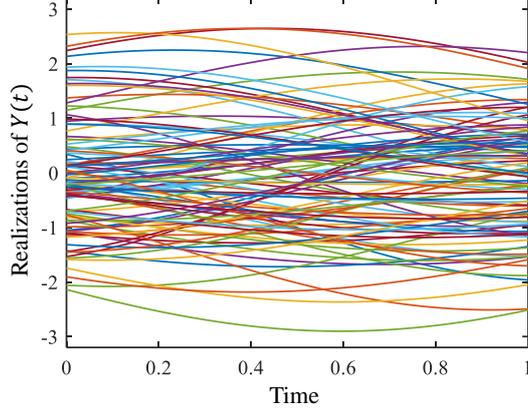


Figure 3: Realizations of  $Y(t)$  of example 4.1.

339  $7.85 \times 10^4 b_0 h_0 (N/m)$ . The LSF is defined as follows:

$$g(\mathbf{x}, F(t), t) = \frac{(b_0 - 2\kappa t)[h_0 - 2\kappa t]^2 f_y}{4} - \left( \frac{F(t)L}{4} + \frac{7.85 \times 10^4 b_0 h_0 L^2}{8} \right) \quad (45)$$

340 where  $\kappa = 3 \times 10^{-5} \text{m/year}$  control the corrosion rate and  $t \in [0, 20]$  years represents time parameter;  $b_0$ ,  $h_0$   
 341 and  $f_y$  are the initial width and height of the cross section, and the yield strength, respectively. Table 2  
 342 lists the details of the random parameters.

343 The time interval  $[0, 12]$  and Gaussian process  $F(t)$  are discretized with 241 time instants and six inde-  
 344 pendent standard normal variables, respectively. The TRA results provided by different methods are listed  
 345 in Table 3. One can observe that all investigated methods can yield the results close to the reference value  
 346 provided by MCS. In terms of the efficiency, however, the proposed method requires much less iterations and  
 347 LSF evaluations than other methods. The comparison of computation time shows that the proposed method  
 348 requires less CPU time than both REAL and SILK methods. When  $q$  varies from 1 to 8 in the proposed  
 349 method, the number of LSF evaluations gradually increases. Besides, the proposed method requires the  
 350 least number of iterations and computation time for the setting of  $q = 5$ . This observation indicates that  
 351 selecting too many training points per iteration does not necessarily result in a reduction in the the number  
 352 of iterations and the total computation time.

353 Fig. 5 schematically presents the TDFP within the time interval  $[0, 20]$  years, where the error bar indicate

354 the range of the mean  $\pm 2$  standard deviations of the TDFP. It is found that the TDFP estimates obtained  
 355 by the proposed method are similar to the reference values provided by MCS. These results confirm that  
 356 the proposed EPAK method is capable of estimating the TDFP across varied subintervals with satisfactory  
 357 accuracy. As shown in Fig. 5, the failure probability is rather low (e.g., around  $10^{-3}$ ) within small time  
 358 intervals. The proposed method still maintains high estimation accuracy, which indicates the effectiveness  
 359 of the proposed method in small TDFP problems.

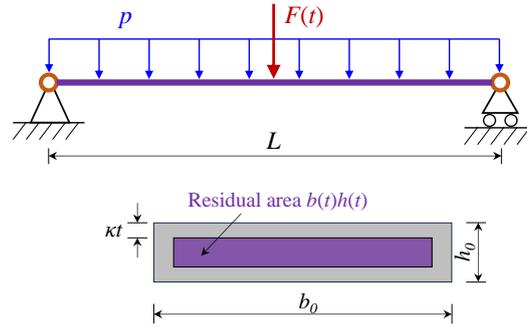


Figure 4: Schematic diagram of the corroded beam.

Table 2: Distributions of inputs of example 4.2.

Input variable	Distribution	Mean	Standard deviation	Autocorrelation function
$f_y$ (MPa)	Lognormal	240	24	-
$b_0$ (m)	Lognormal	0.2	0.01	-
$h_0$ (m)	Lognormal	0.03	0.003	-
$F(t)$ (kN)	Gaussian process	3.5	0.7	$\kappa(t_1, t_2) = \exp\left[-\frac{1}{25}(t_1 - t_2)^2\right]$

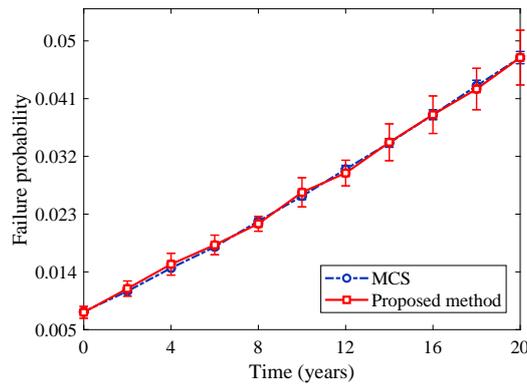


Figure 5: TDFP of example 4.2 ( $q = 1$ ).

Table 3: TRA results of example 4.2.

Methods	$N_{ite}$	$N_{eva}$	$\hat{P}_f$	$\epsilon_{\hat{P}_f}$	COV $[\hat{P}_f]$	CPU Time (s)	
MCS	-	$201 \times 10^6$	$4.74 \times 10^{-2}$	-	0.45%	2.0	
t-IRS <sup>1</sup>	-	171.4	$4.71 \times 10^{-2}$	0.63%	-	-	
eSPT <sup>1</sup>	-	113.4	$4.88 \times 10^{-2}$	2.95%	-	-	
STRA <sup>1</sup>	45.4	56.4	$4.78 \times 10^{-2}$	0.84%	-	-	
SILK	40.4	51.4	$4.73 \times 10^{-2}$	0.21%	2.51%	432.8	
REAL	31.1	42.1	$4.75 \times 10^{-2}$	0.21%	3.29%	183.9	
Proposed method	$q = 1$	8.7	19.7	$4.78 \times 10^{-2}$	0.84%	4.50%	6.1
	$q = 2$	7.3	24.6	$4.74 \times 10^{-2}$	0	8.12%	7.2
	$q = 3$	6.1	27.3	$4.69 \times 10^{-2}$	1.05%	6.63%	6.8
	$q = 4$	4.9	27.6	$4.62 \times 10^{-2}$	2.53%	4.91%	6.5
	$q = 5$	4.2	28.0	$4.74 \times 10^{-2}$	0	3.98%	6.0
	$q = 6$	5.2	37.2	$4.79 \times 10^{-2}$	1.05%	3.14%	6.9
	$q = 7$	4.6	37.2	$4.76 \times 10^{-2}$	0.42%	5.85%	6.3
	$q = 8$	4.5	40.0	$4.60 \times 10^{-2}$	2.95%	3.51%	7.1

<sup>1</sup> The results are taken from research [60].

### 360 4.3. Turbine blade

361 A turbine blade model adapted from Matlab PDE toolbox is investigated as the third example, whose  
362 finite element model (FEM) and von Mises stress distribution are shown in Fig. 6. The Young's modulus,  
363 Poisson's ratio, coefficient of thermal expansion and the thermal conductivity are denoted as  $E$ ,  $\lambda$ ,  $\alpha$  and  
364  $T_c$ , respectively. The temperature of the interior cooling air and the suction sides are denoted as  $T_1$  and  
365  $T_2$ , respectively. In this example, failure is defined as the maximum von Mises stress exceeding an allowable  
366 threshold. Taking into account that the allowable threshold decreases with the time  $t$ , the LSF is defined  
367 as:

$$g(\mathbf{x}, \mathbf{F}(t), t) = \sigma_{at} e^{-0.02t} - \sigma_m(\mathbf{x}, \mathbf{F}(t)) \quad (46)$$

368 where  $\sigma_{at} = 1.5\text{GPa}$  represents the initial allowable threshold;  $\sigma_m(\mathbf{x}, \mathbf{F}(t))$  denotes the maximum stress  
369 provided by FEM;  $F_1(t)$  and  $F_2(t)$  in  $\mathbf{F}(t)$  represent the pressure loads at the pressure side and suction  
370 side, respectively. Table 4 lists the details of the input variables.

371 The time interval  $[0, 12]$  and the Gaussian processes  $F_1(t)$  and  $F_2(t)$  are discretized with 121 time instants,  
372 five and eight independent normal variables, respectively. Table 5 presents the TRA results using different  
373 methods. For MCS,  $121 \times 10^5$  LSF evaluations are required to ensure the accuracy of TDFP estimate,  
374 which is computationally prohibitive. Instead, a simplified simulation with 12,100 LSF evaluations (taking

375 approximately 3.76 hours) is performed to provide an approximate assessment of the computational time.  
 376 The proposed method, SILK and REAL provide similar TDFP estimates, while the proposed method exhibits  
 377 higher efficiency in terms of  $N_{ite}$  and computation time than the counterparts. When  $q$  varies from 1 to  
 378 8, the proposed method requires minimal  $N_{ite}$  for the setting of  $q = 5$ . As for the computation time, the  
 379 setting of  $q = 4$  would minimize the computation time in this example.

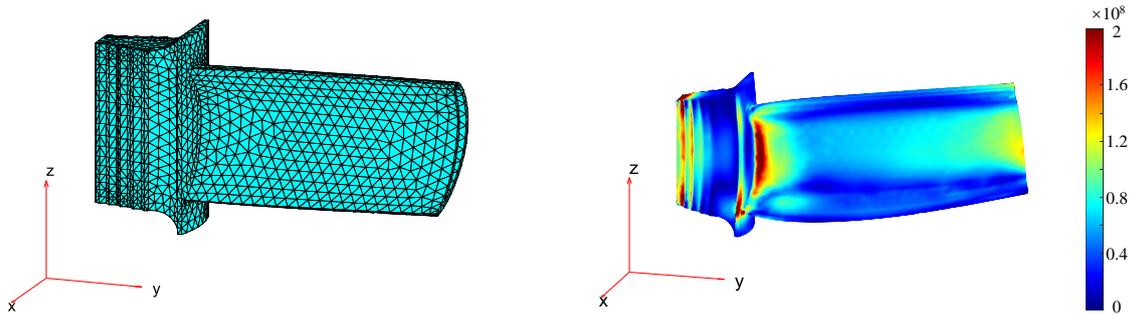


Figure 6: FEM and stress distribution of the turbine blade (unit: Pa).

Table 4: Distributions of inputs of example 4.3.

Variable	Distribution	Parameter 1	Parameter 2	Autocorrelation function
$E$	Normal	220	22	-
$\alpha$	Normal	$1.27 \times 10^{-5}$	$1.27 \times 10^{-6}$	-
$\lambda$	Lognormal	0.27	0.0216	-
$T_c$	Lognormal	11.5	1.38	-
$T_1$	Uniform	130	170	-
$T_2$	Uniform	950	1050	-
$F_1(t)$	Gaussian process	500	100	$\kappa(t_1, t_2) = \exp\left[-\frac{1}{16}(t_1 - t_2)^2\right]$
$F_2(t)$	Gaussian process	450	90	$\kappa(t_1, t_2) = \exp\left[-\frac{1}{4}(t_1 - t_2)^2\right]$

Note: For the uniform distributions, parameter 1 and 2 respectively represent the lower and upper-bounds; for other distributions, they denote the mean and standard deviation, respectively.

#### 380 4.4. Arch bridge

381 In this section, a lower-bearing arch bridge with the calculated span of 150m and the rise-to-span ratio  
 382 of 1:5 is considered [61], as shown in Fig. 7. The arch bridge features 34 suspenders, each spaced 6.8m  
 383 apart. A three-dimensional FEM model is built on the OpenSEES platform, consisting of 241 nodes and  
 384 439 elements, as shown in Fig. 8. The elastic beam-column element is adopted to simulate the suspenders,  
 385 main girder and arch ribs. A increasing heavy load  $F(t)$  applied in the mid-span is considered and modeled  
 386 as a non-stationary Gaussian process. Considering that the suspenders are subjected to the corrosion effect,

Table 5: TRA results of example 4.3.

Methods	$N_{ite}$	$N_{eva}$	$\hat{P}_f$	$\epsilon_{\hat{P}_f}$	$COV[\hat{P}_f]$	CPU Time	
MCS	-	$121 \times 10^5$	-	-	-	$\approx 157.0$ (days)	
SILK	369.5	380.5	$5.28 \times 10^{-3}$	-	2.65%	29,962.6 (s)	
REAL	191.3	202.3	$5.28 \times 10^{-3}$	0	6.86%	8382.3 (s)	
	$q = 1$	110.8	121.8	$5.19 \times 10^{-3}$	1.70%	5.41%	1121.4 (s)
	$q = 2$	58.7	127.4	$5.15 \times 10^{-3}$	2.46%	2.82%	1032.3 (s)
	$q = 3$	42.2	135.6	$5.36 \times 10^{-3}$	1.52%	4.49%	709.8 (s)
Proposed method	$q = 4$	33.3	141.2	$5.33 \times 10^{-3}$	0.95%	6.08%	432.4 (s)
	$q = 5$	28.9	151.5	$5.28 \times 10^{-3}$	0	5.24%	643.2 (s)
	$q = 6$	36.4	224.4	$5.26 \times 10^{-3}$	0.38%	5.90%	769.4 (s)
	$q = 7$	34.7	247.9	$5.17 \times 10^{-3}$	2.08%	4.59%	858.8 (s)
	$q = 8$	37.9	307.2	$5.38 \times 10^{-3}$	1.89%	4.90%	1180.9 (s)

387 the residual area is denoted as  $A(t) = A_1 \times (1 - 0.007t)$ , where  $A_1$  is the initial area of the suspenders. The  
388 failure is defined as the maximum deflection in excess of a safety threshold  $\Delta_s = 10$  cm. The LSF is thus  
389 written as:

$$g(\mathbf{X}, F(t), t) = \Delta_s - \Delta(A(t), E_1, A_2, E_2, I, F_1, F(t)) \quad (47)$$

390 where the initial area of the arch ribs is denoted as  $A_2$ .  $E_1$  and  $E_2$  represent the Young's modulus for the  
391 suspenders and arch ribs, respectively; The main girder's stiffness is characterized by its area moment of  
392 inertia  $I$ ; The total applied load, including both dead and live components, is represented as  $F_1$ . Table 6  
393 lists the details of input variables.

394 The time interval  $[0, 50]$  and the non-stationary Gaussian process  $F(t)$  are discretized with 501 time  
395 instants and sixteen independent normal variables, respectively. Fig. 9 shows one hundred realizations of  
396  $F(t)$ . Table 7 presents the obtained TDFP estimates by different methods. Similar to the third example,  
397 the CPU time required by MCS is approximately calculated with 10,200 LSF evaluations in this example.  
398 The TDFP estimates obtained by all methods are relatively close; however, the proposed EPAK method  
399 demonstrates significantly higher efficiency than the other methods. Specifically, the proposed method with  
400  $q = 1$  requires 66.6 LSF evaluations on average, while SILK and REAL requires 113.8 and 88.0 evaluations,  
401 respectively. Besides, the proposed method costs much less iterations than the counterparts when specifying  
402 a large value of  $q$ . As for the computation time, SILK and REAL takes 13,702.3 seconds and 6000.4 seconds,  
403 respectively, whereas the proposed EPAK method takes 270.1 to 456.2 seconds. When  $q$  varies from 1 to 8,

404 the required computation time is minimal for the setting of  $q = 5$ . Meanwhile, the corresponding  $N_{ite}$  and  
 405 COV are relatively small.

406 The time-dependent reliability results are schematically depicted in Fig. 10, where the reliability index  
 407 (denoted as  $\beta$ ) is obtained by  $\beta = \Phi^{-1}(1 - \hat{P}_f)$ . The error bars show the range of the mean  $\pm 2$  standard  
 408 deviations of the TDFP and the reliability index, respectively. It can be observed that as the service life  
 409 increases, the failure probability increases. Correspondingly, the reliability index gradually decreases, which  
 410 following an approximately linear trend. Besides, the reliability index at  $t = 50$  ( $\beta_{50} = 2.95$ ) decreases by  
 411 29.76% compared to the initial service status ( $\beta_0 = 4.20$ ), which reflects the necessity of performing TRA  
 412 for this arch bridge problem.

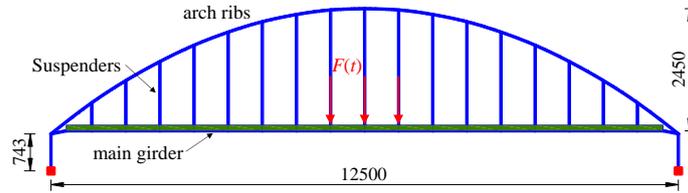


Figure 7: Illustrative diagram of the arch bridge (unit: cm).

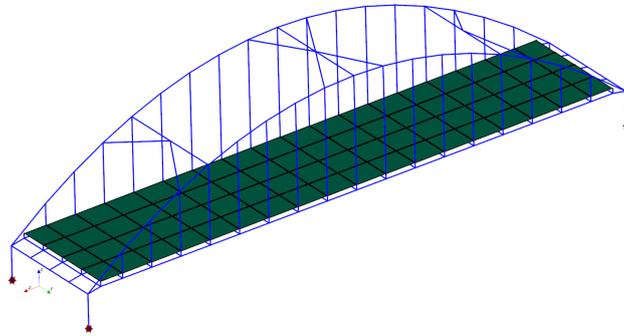


Figure 8: FEM of the arch bridge.

## 413 5. Conclusions

414 This study proposed a new method termed ‘Error-informed Parallel Adaptive Kriging’ (EPAK) for  
 415 efficient TRA. Specifically, the VAIS was adapted in a sequential way to estimate the small TDFP based  
 416 on the trained single-loop Kriging model, which could decrease the sample size and total computation

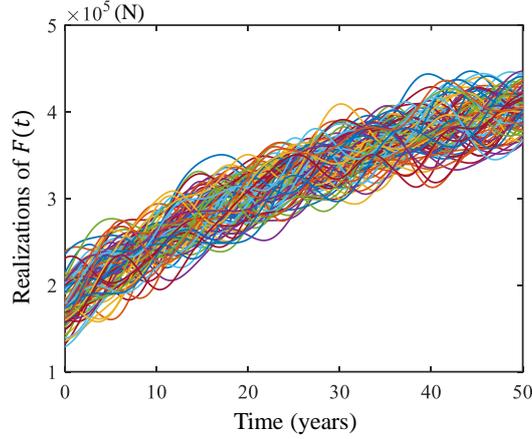


Figure 9: Realizations of the Gaussian process of example 4.4.

Table 6: Distributions of inputs of example 4.4.

Input variable	Distribution	Mean	Standard deviation	Autocorrelation function
$A_1(\text{m}^2)$	Lognormal	$3 \times 10^{-3}$	$3 \times 10^{-4}$	-
$E_1(\text{Pa})$	Normal	$2 \times 10^{11}$	$2 \times 10^{10}$	-
$A_2(\text{m}^2)$	Lognormal	2.8	0.28	-
$E_2(\text{Pa})$	Normal	$2.1 \times 10^{11}$	$2.1 \times 10^{10}$	-
$I(\text{m}^4)$	Lognormal	$5.6 \times 10^{-2}$	$8.4 \times 10^{-3}$	-
$F_1(\text{N/m})$	Gumbel	$5.5 \times 10^7$	$1.1 \times 10^7$	-
$F(t)(\text{kN})$	Gaussian process	$180 + 180 \ln \left(1 + \frac{t}{20}\right)$	$18 + 18 \ln \left(1 + \frac{t}{20}\right)$	$\kappa(t_1, t_2) = \exp \left[ -\frac{1}{16} (t_1 - t_2)^2 \right]$

417 time. Besides, the maximum relative error of TDFP estimation was derived, based on which a stopping  
418 criterion was developed by judging whether the maximum relative error was below a predefined threshold.  
419 Finally, a parallel sampling strategy was proposed through combining the relative error contribution and an  
420 introduced influence function, which could not only select multiple training points but also overcome the  
421 problem of unnecessary LSF evaluations caused by excessive clustering. Several examples were studied to  
422 validate the applicability of proposed EPAK method. Results demonstrated that the proposed method can  
423 estimate small TDFPs with satisfactory accuracy. More importantly, the proposed EPAK method required  
424 much less LSF evaluations, iterations and CPU time when compared to other TRA methods, demonstrating  
425 its superior efficiency. Besides, numerical results showed that selecting too many training points in each  
426 iteration did not necessarily result in a reduction in the number of iterations and the total computation  
427 time. According to the investigated examples, four or five points was sufficient and effective to select, and  
428 was therefore suggested for the proposed method.

Table 7: TRA results of example 4.4.

Methods	$N_{ite}$	$N_{eva}$	$\hat{P}_f$	$\epsilon_{\hat{P}_f}$	$COV[\hat{P}_f]$	CPU Time	
MCS	-	$501 \times 10^6$	-	-	-	$\approx 6543.1$ (days)	
SILK	102.8	113.8	$1.57 \times 10^{-3}$	-	3.55%	13,702.2 (s)	
REAL	77.0	88.0	$1.60 \times 10^{-3}$	1.91%	4.29%	6000.4 (s)	
Proposed method	$q = 1$	55.6	66.6	$1.61 \times 10^{-3}$	2.55%	2.84%	456.2 (s)
	$q = 2$	29.4	68.8	$1.60 \times 10^{-3}$	1.91%	6.05%	282.3 (s)
	$q = 3$	22.3	75.9	$1.61 \times 10^{-3}$	2.55%	4.12%	279.0 (s)
	$q = 4$	17.0	76.0	$1.58 \times 10^{-3}$	0.64%	4.14%	354.9 (s)
	$q = 5$	15.3	83.5	$1.59 \times 10^{-3}$	1.27%	3.33%	270.1 (s)
	$q = 6$	15.8	100.8	$1.62 \times 10^{-3}$	3.18%	4.56%	355.8 (s)
	$q = 7$	15.2	111.4	$1.61 \times 10^{-3}$	2.55%	5.91%	389.1 (s)
	$q = 8$	13.8	114.4	$1.59 \times 10^{-3}$	1.27%	4.96%	340.0 (s)

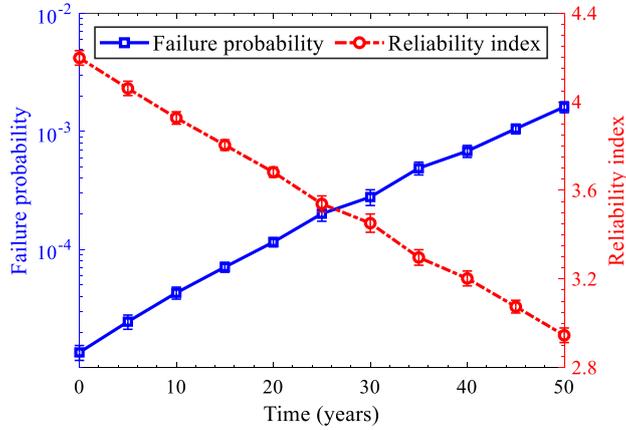


Figure 10: TDFP and reliability index of example 4.4.

429 The proposed EPAK method may perform weakly in high dimensions owing to the inherent limitations  
 430 of the Kriging model. Further research is still required to address this problem. Besides, the proposed  
 431 method could be adapted for time-dependent reliability-based design optimization.

#### 432 Declaration of competing interest

433 The authors declare that they have no known competing financial interests or personal relationships that  
 434 could have appeared to influence the work reported in this paper.

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440 **Appendix A. Kriging model**

441 Kriging model is an interpolation-based regression method including linear regression and stochastic  
 442 process, which is written as [55]:

$$\hat{g}(\mathbf{x}) = \mathbf{f}^T(\mathbf{x})\boldsymbol{\zeta} + \epsilon(\mathbf{x}) \quad (\text{A.1})$$

443 where  $\hat{g}(\mathbf{x})$  denotes the predicted response;  $\mathbf{f}(\mathbf{x})$  is the basis function vector;  $\boldsymbol{\zeta}$  represents the regression  
 444 coefficients vector;  $\epsilon(\mathbf{x})$  denotes a Gaussian process with the mean of zero and the covariance of  $\kappa(\mathbf{x}^{(i)}, \mathbf{x}^{(j)}) =$   
 445  $\sigma^2 R(\mathbf{x}^{(i)}, \mathbf{x}^{(j)})$ , where  $R(\mathbf{x}^{(i)}, \mathbf{x}^{(j)})$  is the correlation function and denoted as:

$$R(\mathbf{x}^{(i)}, \mathbf{x}^{(j)}) = \exp \left[ - \sum_{k=1}^n \theta_k \left( x_k^{(i)} - x_k^{(j)} \right)^2 \right] \quad (\text{A.2})$$

446 where  $\theta_k$  ( $k = 1, 2, \dots, n$ ) are correlation parameters and estimated using the maximum likelihood method  
 447 [55]. Given  $n_0$  training points  $\mathbf{X} = [\mathbf{x}^{(1)}, \mathbf{x}^{(2)}, \dots, \mathbf{x}^{(n_0)}]$  and the responses  $\mathbf{G} = [g(\mathbf{x}^{(1)}), g(\mathbf{x}^{(2)}), \dots, g(\mathbf{x}^{(n_0)})]$ ,  
 448  $\boldsymbol{\zeta}$  and  $\sigma^2$  are estimated as:

$$\hat{\boldsymbol{\zeta}} = (\mathbf{F}^T \mathbf{R}^{-1} \mathbf{F})^{-1} \mathbf{F}^T \mathbf{R}^{-1} \mathbf{G} \quad (\text{A.3})$$

$$\hat{\sigma}^2 = \frac{(\mathbf{G} - \hat{\boldsymbol{\zeta}})^T \mathbf{R}^{-1} (\mathbf{G} - \hat{\boldsymbol{\zeta}} \mathbf{F})}{n_0} \quad (\text{A.4})$$

450 where  $\mathbf{F}$  is the regression matrix with  $F_{i,j} = f_j(\mathbf{x}^{(i)})$ ,  $i = 1, \dots, n_0$ ,  $j = 1, \dots, n$ ;  $\mathbf{R} = [R(\mathbf{x}^{(i)}, \mathbf{x}^{(j)})]$  ( $i, j =$   
 451  $1, \dots, n_0$ ) is the correlation matrix.

452 The mean prediction  $\mu_{\hat{g}}(\mathbf{x})$  and the variance prediction  $\sigma_{\hat{g}}^2(\mathbf{x})$  are obtained as follows:

$$\mu_{\hat{g}}(\mathbf{x}) = \mathbf{f}^T(\mathbf{x})\hat{\boldsymbol{\zeta}} + \mathbf{r}^T(\mathbf{x})\mathbf{R}^{-1}(\mathbf{G} - \mathbf{F}\hat{\boldsymbol{\zeta}}) \quad (\text{A.5})$$

453

$$\sigma_{\hat{g}}^2(\mathbf{x}) = \hat{\sigma}^2 \left[ 1 + \mathbf{u}^T(\mathbf{x}) (\mathbf{F}^T \mathbf{R}^{-1} \mathbf{F}^T)^{-1} \mathbf{u}(\mathbf{x}) - \mathbf{r}^T(\mathbf{x}) \mathbf{R}^{-1} \mathbf{r}(\mathbf{x}) \right] \quad (\text{A.6})$$

454 where  $\mathbf{u}(\mathbf{x}) = \mathbf{F}^T \mathbf{R}^{-1} \mathbf{r}(\mathbf{x}) - \mathbf{f}(\mathbf{x})$ ;  $\mathbf{r}(\mathbf{x}) = [R(\mathbf{x}, \mathbf{x}^{(1)}), \dots, R(\mathbf{x}, \mathbf{x}^{(n_0)})]^T$  is the correlation coefficient vector  
455 between the predicted point and the training set. Note that the construction and prediction of the Kriging  
456 model can be easily performed with DACE toolbox. More details may refer to [62].

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