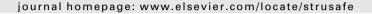
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Evaluation of multivariate normal integrals for general systems by sequential compounding

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ABSTRACT

System reliability analysis often requires efficient and accurate evaluation of a multivariate normal integral. Despite recent advances in system reliability analysis methods, it is still a challenging task especially when the definition of the system event is complex; the system has a large number of components; and/ or the component events have significant statistical dependence. This paper presents a new method developed for evaluating multivariate normal integrals defined for general system events including series, parallel, cut-set and link-set systems. The method compounds two components coupled by union or intersection sequentially until the system becomes a single compound event. Efficient numerical procedures are developed for obtaining the reliability index of the new compound event, and the correlation coefficients between the compound event and the remaining component events, at each step of the sequential compounding. The accuracy and efficiency of the proposed method, and its applicability to various types and sizes of multivariate normal integrals are demonstrated by numerical examples.

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1. Introduction

For risk management of an engineering system, it is essential to evaluate the likelihood of its failure event efficiently and accurately. Sometimes, the failure needs to be described by a *system* event, i.e. a logical (or Boolean) function of multiple *component* events, each of which representing the failure of a physical component, the occurrence of a failure mode, or the structural damage at a particular location. The probability of such a system event is often formulated as a multivariate normal integral

$$P(E_{\text{sys}}) = \int_{\Omega} \varphi_n(\mathbf{z}; \mathbf{R}) d\mathbf{z}$$

$$= \int_{\Omega} \frac{1}{(2\pi)^{n/2} \sqrt{\det \mathbf{R}}} \exp\left(-\frac{1}{2} \mathbf{z}^{\mathsf{T}} \mathbf{R}^{-1} \mathbf{z}\right) d\mathbf{z}$$
(1)

where E_{sys} is the system event of interest; Ω denotes the domain of a system event defined in the space of n standard normal random variables \mathbf{Z} ; $\varphi_n(.)$ is the joint probability density function (PDF) of \mathbf{Z} ; \mathbf{R} is the correlation coefficient matrix of \mathbf{Z} ; and det \mathbf{R} denotes the determinant of \mathbf{R} . For example, if the system failure event is defined as $E_{\text{sys}} = (E_1 \cap E_2) \cup E_3$, the domain Ω is determined as

$$\Omega = \{(z_1,z_2,z_3) | [(z_1\leqslant -\beta_1)\cap (z_2\leqslant -\beta_2)] \cup (z_3\leqslant -\beta_3)\} \tag{2}$$

where β_i is the reliability index of the *i*th component event, i = 1, 2, 3. If the first order reliability method (FORM) is used for component reliability analysis, the correlation coefficient matrix **R** can be determined from the inner products of negative normalized gradient vectors evaluated at the design points [1].

There exists no closed form solution for this multi-fold integration. A direct numerical integration is impractical in many cases because, as the number of component events increases, the computational cost increases rapidly and significant numerical errors are accumulated [2]. Furthermore, a complex system failure domain may render a numerical integration challenging. A number of methods have been developed to compute the multivariate normal integral in Eq. (1) mostly for parallel systems (i.e., intersection of component events) and series systems (i.e., union of component events): first order approaches to multivariate normal integration (FOMN) [3,4], sequentially conditioned importance sampling (SCIS) [5], the product of conditional marginal (PCM) method [6–8], and quasi Monte Carlo simulation methods employing transformation and conditional expectations [9–11].

One of the few methods developed for non-series or non-parallel system problems can be found in the software RELSYS [12]. The method estimates the system probability using the concept of "equivalent component" [13], which sequentially replaces sub-series or sub-parallel systems into equivalent normal components until the system is simplified to a single equivalent component. In this method, individual component reliability indexes and the correlation coefficients between components are first evaluated by

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FORM. The reliability indexes of the equivalent components are then estimated by use of a multivariate normal integral evaluation method while the equivalent directional cosines are approximately obtained by a finite difference method. It is known that this method provides accurate results for parallel systems with five or fewer components. However, it may result in significant error for series systems consisting of components with the same reliability indexes [12], and the errors may accumulate as the size of a general system increases.

The authors of this paper recently developed a matrix-based system reliability (MSR) method [14,15]. This method depends on neither sampling methods nor a finite difference based method, and computes the probability of a general system event including series, parallel, cut-set and link-set system and its parameter sensitivities efficiently and accurately using a convenient matrixbased framework. The method identifies the sources of statistical dependence between components as "common source random variables (CSRVs)" by using, e.g. a generalized Dunnett-Sobel class correlation coefficient matrix [15]. If the MSR method is used for evaluating multivariate normal integrals [16], the dimension of the multivariate integral is reduced from the number of components to that of the CSRVs. However, when it is required to have many CSRVs to describe the correlation coefficients between components accurately, this approach may require a time-consuming multi-fold integration, which prevents rapid evaluations.

In order to overcome these challenges and limitations, this paper presents a new systematic procedure developed for evaluating the multivariate normal integral of any general system efficiently and accurately. The proposed method sequentially compounds two components coupled by union or intersection operation in a general system event into a single equivalent component while the reliability index of the compound event and the correlation coefficients between the compound event and the remaining components are obtained through efficient numerical procedures. The accuracy and efficiency of the proposed method is demonstrated by various numerical examples.

2. Proposed method

The proposed method compounds two components coupled by a logical operation (e.g. union or intersection) sequentially until the system event of interest is simplified into a single compound event. Since it deals with the logical operation of only two components each time, the compounding process would not be encumbered by the complexity of the logical description of a given general system. This section introduces the compounding procedures developed for two components coupled by intersection or union, and explains the overall sequential compounding scheme toward the final solution.

2.1. Compounding two components coupled by intersection

Let us consider two components E_1 and E_2 coupled by intersection. The goal is to compound these into a single equivalent event $E_{1\text{and}2}$. For example, they can appear in a parallel system or a cutset system, and can be compounded as follows.

$$P(E_1 \cap E_2 \cap \dots \cap E_{10}) = P(E_{1\text{and}2} \cap E_3 \cap \dots \cap E_{10}) P(E_1 E_2 E_3 \cup E_4 E_5 E_6 \cup E_7 E_8 E_9) = P(E_{1\text{and}2} E_3 \cup E_4 E_5 E_6 \cup E_7 E_8 E_9)$$
(3)

In order to proceed to the next compounding, we need to find the reliability index of the compound event $E_{1\text{and2}} = E_1 \cap E_2$ (denoted by $\beta_{1\text{and2}}$), and the correlation coefficient between the new compound event $E_{1\text{and2}}$ and each of the remaining component events in the system.

First, the reliability index of the compound event, $\beta_{1\text{and}2}$ is obtained by use of the single-fold numerical integration of $P(E_1 \cap E_2)$ as follows.

$$\beta_{1\text{and2}} = -\Phi^{-1}[P(E_1 \cap E_2)] = -\Phi^{-1}[\Phi_2(-\beta_1, -\beta_2; \rho_{12})] \tag{4}$$

where $\Phi(\cdot)$ is the marginal cumulative distribution function (CDF) of the standard normal distribution; $\Phi_2(\cdot)$ denotes the joint CDF of the bi-variate standard normal distribution; β_1 and β_2 are the reliability indexes of E_1 and E_2 respectively; and $\rho_{1,2}$ is the correlation coefficient between the standard normal random variables Z_1 and Z_2 which respectively represent E_1 and E_2 by $Z_1 \leqslant -\beta_1$ and $Z_2 \leqslant -\beta_2$. For simplicity, the correlation coefficients between standard normal random variables will be hereinafter termed as the correlation coefficient between component *events*. The bi-variate CDF $\Phi_2(-\beta_1, -\beta_2; \rho_{1,2})$ in Eq. (4) can be computed by a single-fold numerical integral [17]

$$\Phi_2(-\beta_1,-\beta_2;\rho_{1,2}) = \Phi(-\beta_1)\Phi(-\beta_2) + \int_0^{\rho_{1,2}} \varphi_2(-\beta_1,-\beta_2;\rho)d\rho \tag{5}$$

where $\varphi_2(\cdot)$ is the bi-variate joint PDF of standard normal distribution. The efficiency of the evaluation of the bi-variate joint CDF can be further improved by using special algorithms. In all the numerical examples in this paper, we use the algorithm that was developed in [18] and later modified in [19].

Next, we obtain the correlation coefficients between the new compound event $E_{1\text{and2}}$ and each of the other remaining component events in the system. Consider a triple of E_1 , E_2 , and E_k , $k=3,\ldots,n$ where n is the current total number of components in the system during a sequential compounding process. We define the equivalent correlation coefficient $\rho_{(1\text{and2}),k}$, $k=3,\ldots,n$ as the one that provides the same estimate on the probability of the event $\Omega_p = [(Z_1 \leqslant -\beta_1) \cap (Z_2 \leqslant -\beta_2)] \cap (Z_k \leqslant -\beta_k)$ even after compounding i.e.

$$\Phi_3(-\beta_1, -\beta_2, -\beta_k; \rho_{1,2}, \rho_{1,k}, \rho_{2,k}) = \Phi_2(-\beta_{1and2}, -\beta_k; \rho_{(1and2),k})$$
 (6)

where $\Phi_3(\cdot)$ denotes the joint CDF of the tri-variate standard normal distribution. Since all the terms in Eq. (6) except $\rho_{(1\text{and2}),k}$ are already given, one can solve the equation for $\rho_{(1\text{and2}),k}$ numerically by finding $\rho_{(1\text{and2}),k}$ that minimizes the difference between the two CDFs by use of a nonlinear optimization algorithm with the constraint $-1\leqslant\rho_{(1\text{and2}),k}\leqslant 1$. However, this is a time-consuming task because the numerical analysis would require repeated evaluations of the multi-fold integrals.

In order to obtain the equivalent correlation coefficients efficiently, we first decompose the CDFs in Eq. (6) using the conditional probabilities, i.e.

$$P(Z_1 \leqslant -\beta_1 \cap Z_2 \leqslant -\beta_2 | Z_k \leqslant -\beta_k) \cdot \Phi(-\beta_k)$$

$$= P(Z_{1and2} \leqslant -\beta_{1and2} | Z_k \leqslant -\beta_k) \cdot \Phi(-\beta_k)$$
(7)

where $Z_{1\text{and2}}$ is the standard normal random variable representing the compound event. For exact evaluations, these conditional probabilities would be obtained from non-normal distributions [6], but the proposed method approximates them by the CDFs of normal random variables. By dividing both terms by $\Phi(-\beta_k)$, Eq. (7) is approximated to

$$\Phi_2(-\beta_{1|k}, -\beta_{2|k}; \rho_{1,2|k}) = \Phi(-\beta_{(1\text{and}2)|k})$$
(8a)

The conditional reliability indexes and conditional correlation coefficients given $Z_k \le -\beta_k$ in this equation are determined exactly by [20]

$$\begin{split} \beta_{1|k} &= (\beta_{1} - \rho_{1,k} A) \Big/ \sqrt{1 - \rho_{1,k}^{2} B} \\ \beta_{2|k} &= (\beta_{2} - \rho_{2,k} A) \Big/ \sqrt{1 - \rho_{2,k}^{2} B} \\ \rho_{1,2|k} &= (\rho_{1,2} - \rho_{1,k} \rho_{2,k} B) \Big/ \Big(\sqrt{1 - \rho_{1,k}^{2} B} \sqrt{1 - \rho_{2,k}^{2} B} \Big) \\ \beta_{(1\text{and}2)|k} &= (\beta_{1\text{and}2} - \rho_{(1\text{and}2),k} A) \Big/ \sqrt{1 - \rho_{(1\text{and}2),k}^{2} B} \end{split} \tag{8b}$$

where $A = \varphi(-\beta_k)/\Phi(-\beta_k)$; and $B = A(-\beta_k + A)$ in which $\varphi(\cdot)$ denotes the PDF of the standard normal distribution. Note that the bi-variate CDF in Eq. (8a) can be computed by a single-fold integral [17], i.e.

$$\begin{split} \Phi_{2}(-\beta_{1|k}, -\beta_{2|k}; \rho_{1,2|k}) &= \Phi(-\beta_{1|k}) \Phi(-\beta_{2|k}) \\ &+ \int_{0}^{\rho_{1,2|k}} \varphi_{2}(-\beta_{1|k}, -\beta_{2|k}, \rho) d\rho \end{split} \tag{9}$$

or by use of an efficient algorithm [18,19].

In summary, Eq. (8) has only one unknown, i.e. $\rho_{(1\text{and}2),k}$ because β_1 , β_2 , β_k , $\rho_{1,2}$, $\rho_{1,k}$ and $\rho_{2,k}$ are available from the original component analysis or previous compounding processes. The efficiency of numerical analysis for obtaining $\rho_{(1\text{and}2),k}$ is significantly increased by solving Eq. (8) instead of Eq. (6) because it does not require performing any multi-fold numerical integration. At each compounding, this numerical analysis is performed (n-2) times to obtain $\rho_{(1\text{and}2),k}$, $k=3,\ldots,n$. In all the numerical examples in the paper, we used a sequential quadratic programming method [21] to obtain the correlation coefficients. In order to check the accuracy of the approximate equation in Eq. (8), Fig. 1 compares the equivalent correlation coefficients obtained by Eqs. (6) and (8) when the three components are equi-correlated, i.e. $\rho_{1,2}=\rho_{1,k}=\rho_{2,k}=\rho$, and have equal reliability indexes, i.e. $\beta_1=\beta_2=\beta_k=\beta$. Close matches are observed at all levels of correlation and reliability indexes considered.

2.2. Compounding two components coupled by union

Let us consider compounding two components E_1 and E_2 coupled by union into a single equivalent event $E_{1 \text{or}2}$. For example, this compounding can appear in a series system or a link-set system, and can be compounded as follows.

$$P(E_1 \cup E_2 \cup \dots \cup E_{10}) = P(E_{1 \text{or} 2} \cup E_3 \cup \dots \cup E_{10})$$

$$P((E_1 \cup E_2 \cup E_3) \cap (E_4 \cup E_5 \cup E_6)) = P((E_{1 \text{or} 2} \cup E_3) \cap (E_4 \cup E_5 \cup E_6))$$
(10)

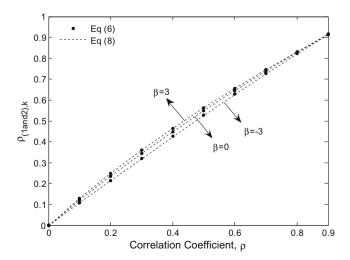


Fig. 1. Comparison of equivalent correlation coefficients $\rho_{(1\text{and}2),k}$ obtained by exact (Eq. (6)) and approximate (Eq. (8)) formulations.

First, using De Morgan's rule and the symmetry of the standard normal distribution, the reliability index of the compound event $E_{1 \text{or}2}$ is obtained by

$$\begin{split} \beta_{1\text{or2}} &= -\Phi^{-1}[P(E_1 \cup E_2)] = -\Phi^{-1}[1 - P(\overline{E}_1 \cap \overline{E}_2)] \\ &= \Phi^{-1}[P(\overline{E}_1 \cap \overline{E}_2)] = \Phi^{-1}[\Phi_2(\beta_1, \beta_2; \rho_{1,2})] \end{split} \tag{11}$$

Next, we obtain the correlation coefficients between the new compound event $E_{1\text{or}2}$ and each of the other remaining component events in the system. Consider a triple of E_1 , E_2 , and E_k , $k=3,\ldots,n$ where n is the current total number of components in the system during a sequential compounding process. We define the equivalent correlation coefficient $\rho_{(1\text{or}2),k}$ as the one that provides the same estimate on the probability of the event $\Omega_u = [(Z_1 \leqslant -\beta_1) \cup (Z_2 \leqslant -\beta_2)] \cap (Z_k \leqslant -\beta_k)$ after compounding, i.e.

$$\int_{\Omega_{-}} \varphi_3(z_1, z_2, z_3; \rho_{1,2}, \rho_{1,k}, \rho_{2,k}) d\mathbf{z} = \Phi_2(-\beta_{1\text{or}2}, -\beta_k; \rho_{(1\text{or}2),k})$$
(12)

Using the same decomposition and approximation used for the intersection case, Eq. (12) is approximated as

$$1 - \Phi_2(\beta_{1|k}, \beta_{2|k}; \rho_{1,2|k}) = \Phi(-\beta_{(1\text{or}2)|k}) \tag{13a}$$

where

$$\beta_{(1\text{or2})|k} = (\beta_{1\text{or2}} - \rho_{(1\text{or2}),k} A) / \sqrt{1 - \rho_{(1\text{or2}),k}^2 B} \tag{13b}$$

The bi-variate CDF in Eq. (13a) can be computed by performing the single-fold numerical integration in Eq. (9) with $-\beta_{1|k}$ and $-\beta_{2|k}$ replaced by $\beta_{1|k}$ and $\beta_{2|k}$, respectively or by use of a special algorithm [18,19]. At each compounding, Eq. (13) is solved numerically for $\rho_{(1\text{or}2),k}, k=3,\ldots,n$ with the constraint $-1\leqslant\rho_{(1\text{or}2),k}\leqslant 1$. Fig. 2 compares the equivalent correlation coefficients obtained by Eqs. (12) and (13) when the three components are equi-correlated, i.e. $\rho_{1,2}=\rho_{1,k}=\rho_{2,k}=\rho$, and have equal reliability indexes, i.e. $\beta_1=\beta_2=\beta_k=\beta$. Close matches are once again observed at all levels of correlation and reliability indexes considered.

2.3. Sequential compounding processes

For a parallel or series system, the procedures introduced in Section 2.1 or Section 2.2 can be applied to adjacent components sequentially until it becomes a single compound event. For a cut-set or link-set system, one can first compound components in each cut-set or link-set until the system becomes an equivalent series or parallel system, respectively. Then, the compound components in

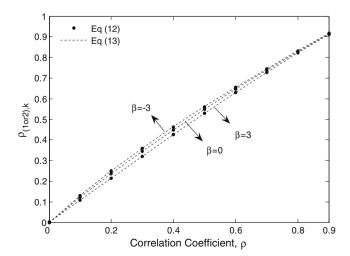


Fig. 2. Comparison of equivalent correlation coefficients $\rho_{(1\text{or2}),k}$ obtained by exact (Eq. (12)) and approximate (Eq. (13)) formulations.

the series or parallel system are compounded sequentially again. One can follow alternative orders of compounding as long as it is compatible with event operation rules (e.g. associative rule and commutative rule). Although each compounding process requires solving the nonlinear equation in Eq. (8) or Eq. (13) numerically for all remaining components, the proposed procedure is efficient because it does not involve sampling or expensive multi-fold numerical integrations. It is also noteworthy that the proposed approach can be used to quantify the statistical dependence between sub-systems, e.g., cut-sets or link-sets by the equivalent correlation coefficients between the compound events.

3. Numerical examples

3.1. Illustrative example: a link-set system with five components

Consider the following link-set system event consisting of five equally reliable components with, $\beta_i = -1$, i = 1, ..., 5:

$$E_{sys} = E_1 \cap E_2 \cap (E_3 \cup E_4) \cap E_5 \tag{14}$$

The correlation coefficient matrix of the standard normal random variables representing the five components is given as

$$\mathbf{R} = \begin{bmatrix} 1 & 0.8 & 0.6 & 0.4 & 0.2 \\ & 1 & 0.8 & 0.6 & 0.4 \\ & & 1 & 0.8 & 0.6 \\ sym. & & 1 & 0.8 \\ & & & & 1 \end{bmatrix}$$
(15)

In this example, each step of sequential compounding is shown along with the updated reliability index and correlation coefficients to demonstrate the proposed method. See Fig. 3 for an illustration of the procedure.

First, we compound the subsystem $E_1 \cap E_2$ into an equivalent component E_A . The reliability index of this compound component, β_A is computed as -0.773 by Eq. (4), and the system definition and correlation coefficient matrix are then updated to

$$E_{\text{sys}} = E_A \cap (E_3 \cup E_4) \cap E_5 \tag{16}$$

and

$$\mathbf{R} = \begin{bmatrix} 1 & \mathbf{0.725} & \mathbf{0.522} & \mathbf{0.319} \\ 1 & 0.8 & 0.6 \\ sym. & 1 & 0.8 \\ & & & 1 \end{bmatrix}$$
 (17)

where the correlation coefficients between the new compound component and the remaining components (in bold) are obtained by solving Eq. (8) numerically.

Next, we merge the subsystem $(E_3 \cup E_4)$ in Eq. (16) into a compound component E_B . Using Eq. (11) in Section 2.3, the reliability index of the compound component β_B is computed as -1.295, and the system definition and correlation coefficients are respectively updated to

$$E_{sys} = E_A \cap E_B \cap E_5 \tag{18}$$

and

$$\mathbf{R} = \begin{bmatrix} 1 & \mathbf{0.651} & 0.319 \\ & 1 & \mathbf{0.729} \\ sym. & 1 \end{bmatrix}$$
 (19)

where the correlation coefficients between the compound event E_B and the other events (in bold) are computed by solving Eq. (13) numerically. The original system event in Eq. (14) is now trans-

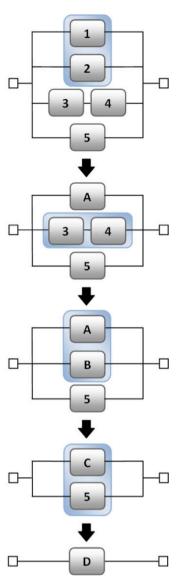


Fig. 3. Sequential compounding procedure for a general system with five components.

formed to a parallel system with three components E_A , E_B and E_5 . By applying the procedure in Section 2.1 two more times, the system in Eq. (18) becomes a single compound event (E_D in Fig. 3). As a result, the system failure probability $P(E_{sys})$ is computed as 0.671 which is fairly close to 0.673 by a Monte Carlo simulation with 10^8 samples.

3.2. Parallel system consisting of 10 components with equal reliability indexes and equal correlation coefficients

The proposed approach is applied to a parallel system consisting of 10 components with the equal reliability indexes, i.e. $\beta_i = \beta, i = 1, 2, \ldots, 10$ and equal correlation coefficients $\rho_{i,j} = 0.5, \ i \neq j$. The probabilities computed by the proposed method are compared to those by the PCM method [6] which is known as an efficient and reasonably accurate method for evaluating multivariate normal integrals for series and parallel systems. Due to the equal correlation coefficients, the exact solution can be obtained by a single-fold numerical integration employing the Dunnett–Sobel class correlation model [22], i.e.

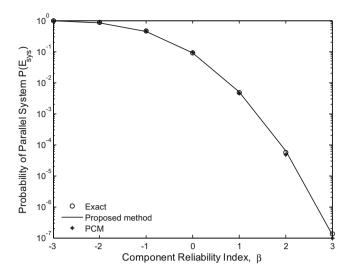


Fig. 4. Comparison of the proposed method and the PCM for parallel systems consisting of 10 components with equal reliability indexes and equal correlation coefficients.

$$P(E_{\rm sys}) = \int_{-\infty}^{\infty} \left[\Phi\left(\frac{-\beta - \sqrt{0.5} \cdot s}{\sqrt{1 - 0.5}}\right) \right]^{N} \varphi(s) ds \tag{20}$$

where N=10. Fig. 4 and Table 1 show the comparison as the equal component reliability index β varies from -3 to 3. Both methods provide quite accurate results while the proposed method shows slightly better accuracy overall.

3.3. Parallel system consisting of 10 components with equal reliability indexes but unequal correlation coefficients

Consider a parallel system with 10 components having equal reliability indexes but their unequal correlation coefficients are determined by

$$\rho_{ij} = 1 - \frac{|i-j|}{10-1}, i, j = 1, \dots, 10, i \neq j$$
 (21)

which satisfies the positive definite condition. The system probabilities are computed by the proposed method, the PCM method, and the MSR method using one, two and three CSRVs. The MSR method requires numerical integration whose dimension is the same as the number of CSRVs while the accuracy of the MSR method gradually enhances with the number of CSRVs increased. Since there is no closed form solution available for this example, we perform Monte Carlo simulation (MCS) with 10⁸ samples for each case. The results in Fig. 5 show that the proposed method has a good agreement with the MCS results for the entire range of the reliability index considered.

3.4. Cut-set system consisting of 10 components with equal reliability indexes but unequal correlation coefficients

Consider the following cut-set system consisting of the components used in the previous example:

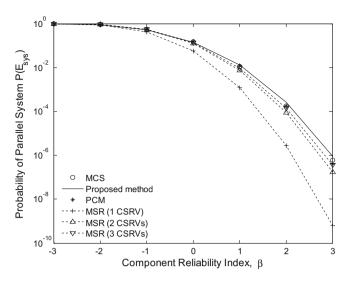


Fig. 5. Comparison of the proposed method, the PCM and MSR method for parallel systems consisting of 10 components with equal reliability indexes but unequal correlation coefficients.

$$E_{sys} = E_1 E_2 \cup E_3 E_4 \cup \ldots \cup E_9 E_{10} \tag{22}$$

Fig. 6 compares the results by the proposed method with those by the MSR method and MCS (10^8 samples). It is noted that most existing methods for multivariate normal integral including the PCM method are developed only for parallel or series systems, thus can not be used for this cut-set system. The proposed method provides the results that are close to those by MCS and its accuracy is even better than that of the MSR using three CSRVs.

3.5. Parallel system consisting of 10 components with unequal reliability indexes and unequal correlation coefficients

Consider a parallel system with 10 components having the equal correlation coefficients 0.5 except $\rho_{1.5}=\rho_{2.4}=0.4$ and $\rho_{1.7}=\rho_{2.6}=\rho_{3.5}=0.9$. The unequal reliability indexes of the 10 components are given by $\beta_i=(4\cdot i-22)/9,\ i=1,\ldots,10$. For this example, the MSR and PCM method show significant errors, as shown in Table 2. By contrast, the proposed method shows a good agreement with the MCS result (10^8 samples).

3.6. Parallel system consisting of 5–50 components with equal reliability indexes and equal correlation coefficients

In order to examine the potential effect of the system size on the performance of the proposed method, we consider parallel systems with 5–50 components having the equal reliability indexes ($\beta = 1, 2 \text{ or } 3$) and equal correlation coefficients 0.5. The exact solution can be obtained by Eq. (20) as N is varied from 5 to 50. As shown by the results in Fig. 7, in contrast to the PCM, the proposed method does not show significant error accumulation as the system size increases. Although the proposed method is not as efficient as the PCM method, it requires affordable computational cost. For example, the computations for 5, 10, ..., 50-component

Table 1Comparison of the proposed method with the PCM for parallel systems consisting of 10 components with equal reliability indexes and equal correlation coefficients.

β	-3	-2	-1	0	1	2	3
Exact Proposed method PCM	$\begin{array}{l} 9.890 \times 10^{-1} \\ 9.884 \times 10^{-1} \\ 9.878 \times 10^{-1} \end{array}$	$\begin{array}{l} 8.669 \times 10^{-1} \\ 8.658 \times 10^{-1} \\ 8.628 \times 10^{-1} \end{array}$	$\begin{array}{l} 4.606 \times 10^{-1} \\ 4.621 \times 10^{-1} \\ 4.697 \times 10^{-1} \end{array}$	9.091×10^{-2} 9.238×10^{-2} 9.332×10^{-2}	4.791×10^{-3} 5.013×10^{-3} 4.552×10^{-3}	5.658×10^{-5} 6.158×10^{-5} 4.735×10^{-5}	1.361×10^{-7} 1.549×10^{-7} 9.982×10^{-8}

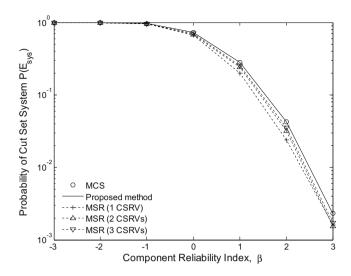


Fig. 6. Comparison of the proposed method and MSR method for cut-set systems consisting of 10 components with equal reliability indexes but unequal correlation coefficients.

Table 2Comparison of the existing methods and the proposed method for a parallel system with unequal reliability indexes and unequal correlation coefficients.

	Proposed method	MSR (3 CSRVs)	PCM	MCS
Probability (×10 ⁻³)	4.585	3.613	2.373	4.480

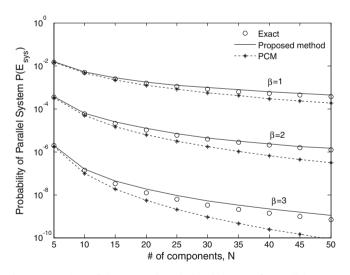


Fig. 7. Comparison of the proposed method and the PCM for parallel systems consisting of 5–50 components with equal reliability indexes and equal correlation coefficients.

systems took 0.16, 0.75, 1.64, 2.59, 3.69, 5.01, 6.38, 8.31, 9.82, and 11.90 seconds, respectively (Matlab® on a personal computer with AMD dual core 2.0 GHz.) The accuracy and efficiency of the proposed method can be further improved by using a numerical solver with higher-performance in computing equivalent correlation coefficients.

3.7. Series system consisting of 10–100 components with equal reliability indexes and equal correlation coefficients

The size effect is examined for series systems as well. Consider series systems having 10–100 components having the equal reli-

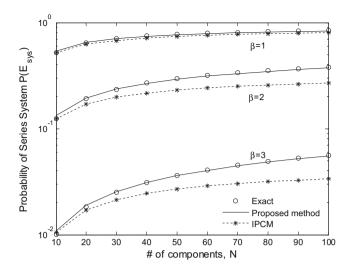


Fig. 8. Performance of the proposed method for series systems consisting of 10–100 components with equal reliability indexes and equal correlation coefficients.

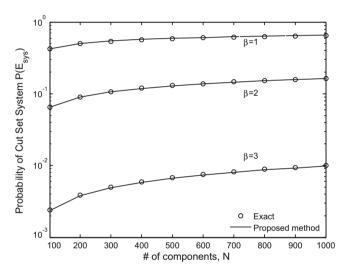


Fig. 9. Performance of the proposed method for cut-set systems consisting of 100–1000 components with equal reliability indexes and equal correlation coefficients.

ability indexes (β = 1, 2 or 3) and equal correlation coefficients 0.5. The exact solution of this problem can be obtained by

$$P(\textit{E}_{\textit{sys}}) = \int_{-\infty}^{\infty} \left\{ 1 - \left[1 - \varPhi \left(\frac{-\beta - \sqrt{0.5} \cdot \textit{s}}{\sqrt{1 - 0.5}} \right) \right]^{\textit{N}} \right\} \phi(\textit{s}) \textit{ds} \tag{23} \right\}$$

where *N* is varied from 10 to 100. Fig. 8 confirms that the size effect of the proposed method is negligible in contrast to the PCM method. This example uses an improved version of the PCM method (IPCM) [8] that was developed for more accurate evaluation of multivariate normal integrals of series systems.

3.8. Cut-set system consisting of 100–1000 components with equal reliability indexes and equal correlation coefficients

Consider the following cut-set system having the equal reliability indexes ($\beta = 1, 2 \text{ or } 3$) and equal correlation coefficients 0.5:

$$E_{\text{sys}} = E_1 E_2 \cup E_3 E_4 \cup \ldots \cup E_{N-1} E_N \tag{24}$$

The exact solution of this problem can be obtained by

$$P(E_{\text{sys}}) = \int_{-\infty}^{\infty} \left\{ 1 - \left[1 - \Phi\left(\frac{-\beta - \sqrt{0.5} \cdot s}{\sqrt{1 - 0.5}}\right)^{2} \right]^{N/2} \right\} \varphi(s) ds$$
 (25)

where *N* is varied from 100 to 1,000. As shown in Fig. 9, the results by the proposed method show good agreement with the exact solutions.

4. Summary and conclusion

A new method is developed for evaluating multivariate normal integrals defined for general systems efficiently and accurately. It sequentially compounds components in a general system event until the system becomes one compound event. For two components coupled by intersection and union, efficient procedures are developed to obtain the equivalent reliability index of the new compound event and the correlation coefficients between the compound event and the remaining components. In order to obtain the correlation coefficients efficiently, the numerical analysis employs approximate decomposition of bi- and tri-variate normal integrals. An important merit of the proposed method is that one can quantify the statistical dependence between sub-systems by the equivalent correlation coefficients between compound events. The proposed method's accuracy and efficiency for series, parallel, cut-set and link-set systems are demonstrated through comparison with other existing methods such as the PCM method and MSR method in various numerical examples. It is also confirmed that the accuracy of the method is not significantly affected by the large number of components in a system. Due to its wide applicability, accuracy and efficiency, the method is expected to enhance the computational capability in various system reliability analysis applications.

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