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**A NON-SIMULATION BASED METHOD FOR INDUCING PEARSON'S
CORRELATION BETWEEN INPUT RANDOM VARIABLES**

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by

Eric R. Druker, Richard L. Coleman and Peter J. Braxton

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A Non-simulation Based Method for Inducing Pearson's Correlation between Input Random Variables

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Abstract

Several previously published papers have cited the need to include correlation in risk-analysis models. In particular, a landmark paper published by Philip Lurie and Matthew Goldberg presented a methodology for inducing Pearson's correlation between input/independent random variables. The one subject, absent from the paper, was a methodology for finding the optimal applied correlation matrix given a desired outcome correlation. Since the publishing of the Lurie-Goldberg paper, there has been continuing discussion on its implementation; however, there has not been any presentation of an optimization algorithm that does not involve the use of computing-heavy simulations. This paper reviews the general methodology used by Lurie and Goldberg (along with its predecessor papers) and presents a non-simulation approach to finding the optimal input correlation matrix, given a set of marginal distributions and a desired correlation matrix.

Introduction

The Complete Correlation Algorithm (CCA) developed by Northrop Grumman and recently implemented in NG developed risk models is a product of more than two years of research and development. Several previously published papers have cited the need to include correlation in risk-analysis models; however, none present an optimization algorithm that does not involve the use of computing-heavy simulations. In particular, a landmark paper published by Philip Lurie and Matthew Goldberg (1998) presented a methodology for inducing Pearson's correlation between input random variables. This paper reviews the general methodology used by Lurie and Goldberg (along with its predecessor papers) and presents the Druker Algorithm: a non-simulation approach to finding the optimal input correlation matrix given a set of marginal distributions and a desired correlation matrix.

The CCA was deliberately created bearing in mind identified environmental factors that prevent easy implementation of commercially available models. No one on the team had any experience implementing correlation into Monte Carlo simulations beyond the use of COTS programs, such as @Risk™ and Crystal Ball™. To determine the best development method, the following factors were considered:

1. The Northrop Grumman risk models need to be of an easily transferable electronic size, as the models are often shared via email or network drives.
2. A diverse group of users must be able to run the software in a variety of work environments; Microsoft Office is the only platform that is transferable to all parties. Users include risk practitioners, program managers and members of pricing organizations; locations include unclassified and classified Northrop Grumman facilities, unclassified and classified customer facilities and home offices.



3. Custom implementations are frequent; much of NGIT-TASC risk work requires risk simulations to be built into pre-existing cost and price models. These models are generally limited to Microsoft Excel and Access; however, Web-based platforms are not unheard of.

The above concerns drove the decision to use Visual Basic source code to develop *the CCA*.

Initially, the development was focused on an algorithm that could induce Pearson's correlation between typical distributions in risk analysis: Bernoulli (discrete), Triangular, Normal and Log-Normal. By limiting the problem to the most-common applications, in theory, the solution should have been easier to find. While attempting to ascertain the maximum correlation between any two Bernoulli distributions, however, the general solution was uncovered. The resulting algorithm induces Pearson's correlation between any set of random variables (while still preserving the marginal distributions) using the Lurie-Goldberg Method and without the use of simulation to find the optimal applied correlation matrix.

The CCA is a compilation of multiple algorithms (each named for their main author(s)) from several sources: existing papers, public source code and internally-developed code. Most of the algorithms used were taken from a variety of existing papers. Although these papers all provided complete algorithms, they sometimes lacked details in how to accomplish key steps; in cases such as these, gaps were filled with open-source code solutions. The optimization of the applied correlation matrix, the last step in the correlation algorithm, was developed entirely by the Northrop Grumman Team.

Definitions and Assumptions

Matrix Definitions:

1. **Consistent Correlation Matrix**—Consistent Correlation matrices have diagonal entries equal to 1.0, all other entries between [-1, 1] are symmetric and positive definite. Consistency is necessary for a viable correlation matrix, but a Consistent Correlation Matrix may not necessarily be viable given the **Parent Distributions**.
2. **Input Correlation Matrix (I)**—The user-inputted correlation matrix. This matrix may or may not be a consistent correlation matrix.
3. **Adjusted Correlation Matrix (L)**—The **Input Correlation Matrix** adjusted to be a **Consistent Correlation Matrix**. This matrix will, by definition, be positive definite. Additionally, the adjusted matrix will be viable as correlations between various distributions of random variables will be achievable. When **(L)** is generated, the differences between **(I)** and **(L)** are minimized.
4. **Applied Correlation Matrix (A)**—The correlation matrix used by the *grand algorithm* to generate correlated random number draws. This matrix may be the same, or very different from, the **Adjusted Correlation Matrix**; the extent of the differences will depend on the random variables to be correlated.
5. **Optimal Applied Correlation Matrix (A')**—The Applied Correlation Matrix optimized using the Lurie-Goldberg Method.
6. **Outcome Correlation Matrix (O)**—The correlation matrix of the simulated variables following the simulation run. The goal of the *grand correlation algorithm* is for **(O)** to be identical to **(L)**.



Other Definitions:

1. **Parent Distribution**—The distributions correlated for use in the simulation. The distributions are simulated using the Inverse CDF technique. The goal is to induce a desired correlation between these distributions.
2. **Pearson's Correlation**—A parametric statistic that measures the *strength and direction of a linear relationship between two random variables* ("Correlation," 2008).
3. **Spearman's Rank Correlation**—A non-parametric statistic that measures the monotonicity of a function without making any assumptions as to the distribution of the variables.
4. **Eigenvalues**—A scalar (L) associated with a matrix such that if (A) is a matrix and (X) is a vector, $AX = LX$. The vector (X) is known as the **Eigenvector** that corresponds to the **Eigenvalue** (L).

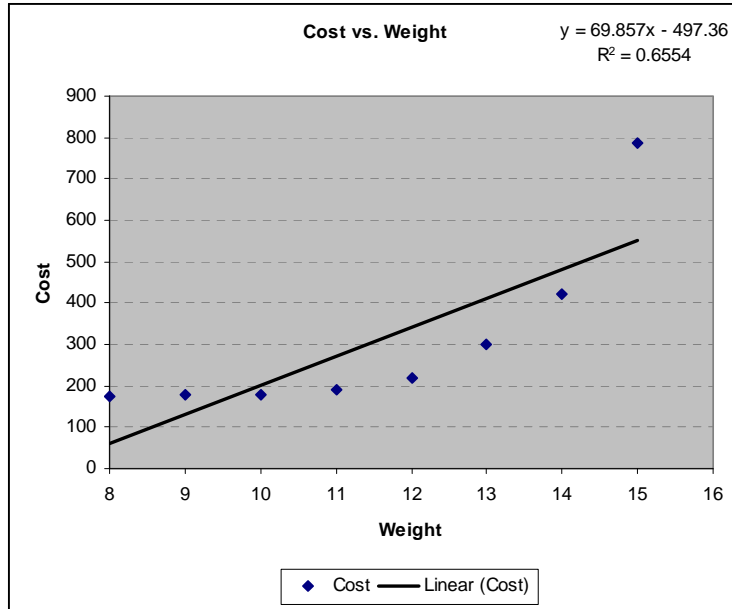
Assumptions:

1. **Normal Distributions**—Any reference to the normal distribution, whether in a univariate or bivariate case, is assumed to be the Standard Normal distribution (Mean of 0, Standard Deviation of 1).

Pearson's vs. Rank Correlation

Most COTS risk tools use Spearman's rank correlation as a substitute for Pearson's correlation between parent distributions. Spearman's rank correlation (a non-parametric statistic) differs from Pearson's correlation (a parametric statistic) in that it measures the monotony of a function, whereas Pearson's correlation measures the strength of the linear relationship between two functions (see Figure 1). Though studies have shown that, using the most common risk distributions, models using rank correlation yield similar results to those using Pearson's (Robinson & Salls, 2004), there is a distinct difference between the two. Although this paper will not detail all the differences between the two measures, a quick (and exaggerated) example is presented below. The *grand algorithm* supersedes the need to substitute for Pearson's correlation with Spearman's rank correlation.





Pearson's Rho	0.81
Spearman's Rho	1.00

Figure 1. Pearson's vs. Spearman's Rank Correlation

Algorithm Overview

There are three main steps behind the *grand algorithm*. An outline of these steps follows, and the upcoming sections of this paper will review each individual step in detail.

1. Correct the **User-Input Correlation Matrix (I)**
 - a. Correct **I** so that it is consistent—both in terms of a general correlation matrix and the properties of the parent distributions being correlated.
 - b. Through these corrections, the Adjusted Correlation Matrix (**L**) will be generated.
2. Optimize the **Applied Correlation Matrix**
 - a. Find the **Optimal Applied Correlation Matrix (A')** such that when **A'** is run through the Lurie-Goldberg Method, the **Outcome Correlation Matrix (O)** is identical to **L**.
3. Correlate the Input Random Variables
 - a. Using **A'**, apply the Lurie-Goldberg Method to correlate the parent distributions.

For purposes of presenting the methodology, it is necessary to show how the input random variables are to be correlated before discussing how to find **A'**.

Correcting the User-Input Correlation Matrix (Part I)

Giving users the ability to input their own correlation matrix allows for the possibility that the **User-Input Correlation Matrix (I)** may not be a viable correlation matrix. Correlation matrices, by definition, have diagonal entries of 1.0. All other entries between [-1, 1] are

symmetric and are positive definite. The first step in inducing correlation between input random variables is checking whether \mathbf{I} is a consistent correlation matrix. If it is not, it must be corrected that it is such.

The *Iman-Davenport Algorithm*, which is based on a paper by Ronald Iman and James Davenport (1982) is used to correct \mathbf{I} in order to make it a consistent correlation matrix. While numerous other papers have been published describing methods to correct \mathbf{I} such that it is altered as little as possible (Higham, 2002), the *Iman-Davenport Algorithm* is the most computationally efficient method the authors uncovered. Given that additional adjustment may be required based on the parent distributions being correlated; the resulting matrix is close enough to \mathbf{I} to satisfy this requirement.

The algorithm corrects \mathbf{I} in three main phases. First, the algorithm checks whether \mathbf{I} is symmetric with diagonal entries of 1.0 and off-diagonal entries between $[-1, 1]$. If it is not, the user is prompted to re-input the matrix, correcting for the discrepancies.

Second, once the above conditions are satisfied, the algorithm checks whether \mathbf{I} is positive-definite. One way to test this is to find the eigenvalues for \mathbf{I} (positive-definite matrices have all positive eigenvalues). The paper referenced did not describe an approach for finding the eigenvalues of the matrix. After further research, the Jacobi Eigenvalue Algorithm was determined to be a sufficiently efficient way to evaluate a matrix's eigenvalues. As a result, the eigenvalues are produced as the diagonals of an otherwise zero-matrix. The Jacobi Eigenvalue Algorithm is computationally inexpensive and pre-existing source code was used in its implementation.

If all eigenvalues for \mathbf{I} are positive and the other conditions have been satisfied, then \mathbf{I} is a consistent correlation matrix. Otherwise, in the third phase, negative eigenvalues are changed to small, positive values (e.g., .000001). The diagonal matrix of adjusted eigenvalues is then multiplied by the associated matrix of eigenvectors (also produced using the Jacobi Eigenvalue Algorithm). That product is, in turn, multiplied by the inverse of the matrix of eigenvectors to arrive at a new matrix that is similar, but not equal to, \mathbf{I} . Lastly, the diagonals are reset to 1.0 as they may have changed during the transformation. This third section of the algorithm is repeated until all eigenvectors of the adjusted matrix are positive. At this point, the user input matrix has been adjusted such that it is a consistent correlation matrix.

Though the **User-Input Correlation Matrix** is now a consistent correlation matrix, the transformation of \mathbf{I} is not complete and the **Adjusted Correlation Matrix (\mathbf{L})** has not been determined. As will be shown later, depending on the parent distributions being correlated, there may be a maximum achievable correlation between any two of the variables. Determination of \mathbf{L} will be covered later in the section: *Correcting the User-Input Correlation Matrix (Part II)*.

Correlating Input Random Variables

In order to understand how the **Applied Correlation Matrix (\mathbf{A})** is to be optimized such that the **Output Correlation Matrix (\mathbf{O})** is identical to the **Adjusted Correlation Matrix (\mathbf{L})**, the method for correlating the parent distributions must first be discussed. It is a well-known fact that normal random variables can be correlated by multiplying a vector of uncorrelated normal



random draws by the Cholesky decomposition¹ of the desired correlation matrix. The Lurie-Goldberg Method takes this one step further using normal random variates to generate correlated uniform random variates. These uniform random variates are then transformed via the inverse-CDF technique to generate draws from the desired parent distributions. In this method, although the correlations between the normal random draws are known, as these draws are transformed into other distributions, the correlations change. Hence, the core problem emerges: how can the **Optimal Applied Correlation Matrix (A')** be uncovered such that **O** matches **L**? Answering this question is key to implementing the Lurie-Goldberg Method. The authors have developed an algorithm that addresses this very question, without necessitating any runs of the simulation. Additionally, they have begun the process of optimizing this algorithm, finding heuristics that allow it to run with a minimal number of calculations.

Implementation and Application of the CCA

The CCA's chief advantage is that it is non-recurring and its implementation requires no simulation. Furthermore, because the algorithm only requires looking at pairs of parent distributions, once the applied matrix has been found for a set of parent distributions, the algorithm must only be run when distributions are added or changed, and even then, only for the new/altered distributions. The algorithm also uses Pearson's correlation while COTS risk tools substitute Spearman's rank correlation.

The applications of the CCA reach beyond the Cost and Risk analysis community; this algorithm is useful anywhere there is a need to induce Pearson's correlation between input variables. For example, this algorithm can be applied to auto correlating, stock market projections in the financial arena and to traditional modeling and simulation situations when correlation is needed. The algorithm was designed with a focus on portability. Because the algorithm is coded with Visual Basic, it can be easily integrated in existing tools and models.

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Open Source Code References:

- The Foxes Team, Italy—<http://digilander.libero.it/foxes>
- Axel Vogt, Germany—http://www.axelvogt.de/axalom/bivariateNormal_Series.zip

¹ The Cholesky Decomposition Matrix of any matrix M is L such that $M = LL^T$



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