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EFFICIENT STAND-ALONE GENERALIZED INVERSE ALGORITHMS AND

SOFTWARE FOR ENGINEERING/SCIENCES APPLICATIONS: RESEARCH

AND EDUCATION

by

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A Dissertation Submitted to the Faculty of Old Dominion University in Partial Fulfillment of the Requirements for the Degree of

DOCTOR OF PHILOSOPHY

CIVIL AND ENVIRONMENTAL ENGINEERING

OLD DOMINION UNIVERSITY August 2012

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ABSTRACT

EFFICIENT STAND-ALONE GENERALIZED INVERSE ALGORITHMS AND SOFTWARE FOR ENGINEERING/SCIENCES APPLICATIONS: RESEARCH AND EDUCATION

Subhash Chandra Bose S V Kadiam Old Dominion University, 2012 Director: Dr. Duc T Nguyen

Efficient numerical procedures for finding the generalized (or pseudo) inverse of a general (square/rectangle, symmetrical/unsymmetrical, non-singular/singular, real/complex numbers) matrix and solving systems of Simultaneous Linear Equations (SLE) are formulated and explained. The developed procedures and its associated computer software (under MATLAB computer environment) have been based on "special Cholesky factorization schemes" (for a singular matrix), the generalized inverse of the matrix product, and were further enhanced by the Domain Decomposition (DD) formulation.

Test matrices from different fields of applications have been chosen, tested and compared with other existing algorithms. The results of the numerical tests have indicated that the developed procedures are far more efficient than existing algorithms.

Furthermore, an educational version of the generalized inverse algorithms and software for solving SLE has also been developed to run any FORTRAN and/or 'C' programs over the web. This developed technology and software is freely available and can run on any device with internet connectivity and browser capability. Copyright, 2012, by Subhash Chandra Bose S V Kadiam, All Rights Reserved

This dissertation is dedicated to my father.

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Most importantly, I would like to thank my family, who have been supportive throughout my life and academic career.

NOMENCLATURE

SVD	Singular Value Decomposition
SLE	Simultaneous Linear Equations
SPD	Symmetric Positive Definite
GMRES	Generalized Minimal Residual
CG	Conjugate Gradient Algorithm
DD	Domain Decomposition Formulations
LDL ^T	LDL Transpose Algorithm
nsu	Number of Sub-domains
pinv()	MATLAB Command to Compute Generalized
	Inverse
inv()	MATLAB Command to Compute Regular Inverse
A ⁺	Generalized Inverse of A
CG	Conjugate Gradient Algorithm
Bi – CG	Bi-Conjugate Gradient Algorithm
GMRES	Generalized Minimal Residual Algorithm
PCG	Preconditioned Conjugate Gradient Algorithm
ODU – ginverse	ODU Generalized Inverse Solver
geninv	Generalized Inverse Algorithm Discussed in [13]
MATLAB — pinv	MATLAB Generalized Inverse Solver
ODU — ginverse iterative	CG Iterative Method in ODU Generalized Inverse
	Solver

ODU - DD	ODU Generalized Inverse Domain Decomposition
	Solver
Original System ginverse MV	ODU Generalized Inverse Solver
Original System geninv	Generalized Inverse Algorithm discussed in [13]

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1. INTRODUCTION

In scientific computing, most computational time is spent on solving systems of Simultaneous Linear Equations (SLE) which can be represented in matrix notations as

$$Ax = b \tag{1.1}$$

where $A \in \mathbb{R}^{n \times n}$ is a singular/non-singular matrix, and b is a given vector in \mathbb{R}^n . For practical engineering/science applications, matrix A can be either sparse (for most cases), or dense (for some cases). Solving large scale system of SLE has been (and continues to be) a major challenging problem for many real-world engineering and science applications.

The generalized (or pseudo) inverse of a matrix is an extension of the ordinary/regular square (non-singular) matrix inverse, which can be applied to any matrix (such as singular, rectangular, etc.). The generalized inverse has numerous important engineering and science applications. Over the past decades, generalized inverses of matrices and their applications have been investigated by many researchers [1-8].

1.1 Literature Survey

Various methods have been proposed for finding the generalized inverse and its associated SLE. Xuzhou Chen et. al. [9] has proposed a method based on a finite recursive algorithm. The approach was based on the symmetric rank-one update. The algorithm proposed by Xuzhou Chen, however, was inefficient (in terms of computational time) and requires lot of computer memory. It has been shown that this algorithm [9] can be only effective for the computation of generalized inverse/MoorePenrose inverse of rectangular matrices (with rows<<cols or cols<<rows) and is inefficient for square matrices.

The most commonly implemented method in programming languages to compute generalized inverse (and its associated SLE) was based on Singular Value Decomposition (SVD) [3, 10-11]. This method is numerically very stable, however, it is computationally expensive for practical applications. MATLAB [12] uses SVD to compute the pseudo/generalized inverse by invoking the built-in function *pinv()*. It should be noted here that in finding the solution for SLE (with square/singular, or rectangular coefficient matrices), MATLAB and most (if not all) other researchers have computed the generalized inverse explicitly. Then, the solution can be found by a simple matrix times vector operation.

Since the standard Eigen-value problems (of $A \times A^{H}$, and $A^{H} \times A$) need to be solved in SVD, this method is computationally expensive. Despite of this fact, solving Eq. (1.1) by using SVD is still more efficient than Xuzhou Chen's proposed finite recursive algorithm [9].

In [6], an efficient algorithm for finding the generalized inverse of a (full rank) rectangular (or square) matrix has been proposed. However, this algorithm has not been able to handle the cases where the matrix has rank deficiency (such as a matrix which has some dependent rows and/or columns)

Pierre Courrieu [13] has proposed an algorithm to explicitly compute the generalized inverse using full-rank Cholesky factorization on the coefficient matrix. His algorithm was based on a theorem to compute the generalized inverse of a product of two

matrices. Pierre Courrieu's algorithm has proven to be more efficient than finite recursive method [9] and SVD [10-11].

1.2 SVD and the Generalized Inverse

A general (square or rectangular) matrix $A \in \mathbb{R}^{n \times n}$ can be decomposed as

$$A = U\Sigma V^H \tag{1.2}$$

where

 $\Sigma = a \ diagonal \ matrix \ (does \ NOT \ have \ to \ be \ a \ square \ matrix)$

$$=\begin{cases} \Sigma_{ij} = 0, for \ i \neq j\\ \Sigma_{ij} \ge 0, for \ i = j \end{cases}$$
(1.3)

[U] and [V] = unitary matrices

and
$$\begin{cases} U^{H} = U^{T}(for \ real \ matrices) \\ U^{H} = U^{-1} \end{cases}$$
(1.4)

Let A be a singular matrix of size $m \times n$ and let k be the rank of the matrix. Based on Eq. (1.2), one has

$$A = U\Sigma V^{H};$$
where $\Sigma = \begin{bmatrix} \sigma_{1} & & \\ & \sigma_{2} & \\ & & \sigma_{k} \end{bmatrix}$
with $\sigma_{1} \ge \sigma_{2} \ge \cdots \ge \sigma_{k} > 0;$

and $\sigma_i = \sqrt{Eigen - Values of A^T A (or AA^T)}$

<u>Note</u>: Eigen-values of $A \times A^{H}$ and Eigen-values of $A^{H} \times A$ are the same. However, the Eigen-vectors of $A \times A^{H}$ and Eigen-vectors of $A^{H} \times A$ are "NOT" the same.

Then, the generalized inverse A^+ of A is the $n \times m$ matrix and is given as

$$A^+ = V \Sigma^+ U^{\rm H} \tag{1.6}$$

(1.5)

where

$$\Sigma^{+} = \begin{bmatrix} E \\ 0 \end{bmatrix} \begin{bmatrix} 0 \\ 0 \end{bmatrix} \text{ and } E \text{ is the } k \times k \text{ diagonal matrix, with}$$
$$E_{ii} = \Sigma_{i}^{-1} \text{ for } 1 \le i \le k$$

More details about computing the SVD from a given matrix [A] can be found in the Appendix A.

1.3 Objective

The main objective of this dissertation is to develop an efficient (in terms of computational time and computer memory requirement) generalized inverse formulation to solve SLE with full or deficient rank of the coefficient matrix. The coefficient matrix can be singular/non-singular, symmetric/unsymmetric, square/rectangular, and with real/complex numbers. The proposed generalized inverse procedures can also be integrated in to the Domain Decomposition (DD) formulation for solving general, large scale SLE commonly encountered in engineering/sciences applications. Due to popular MATLAB software, which is widely accepted by researchers and educators worldwide, the developed code from this work is written in MATLAB language and has the following capabilities/features:

a) A stand-alone generalized inverse software to solve SLE

b) A stand-alone DD generalized inverse software to solve SLE

c) Utilizing sparse storage scheme (whenever possible) for storing data and solving SLEd) Developing user friendly interfaces to test new problems (including the numerical data downloaded from popular web sites [14-15]).

e) Additional wall-time reduction for DD generalized inverse solver can be realized/achieved by performing "parallel matrix times matrix" operations under MATLAB-MPI computer environment.

f) Developing an "educational version" (written in FORTRAN language) of the software, which uses the generalized inverse for solving general SLE on the internet.

In Chapter 2, some major algorithms for solving SLE by direct and iterative methods are reviewed. These methods are mainly designed for solving non-singular SLE. Simple/basic domain decomposition (DD) algorithms, using mixed direct-iterative solvers, are discussed in Chapter 3. Major works in this dissertation are presented in Chapter 4, where efficient "generalized (or pseudo-) inverse" algorithms are thoroughly explained with and without incorporating the DD formulation. The numerical performance of the proposed algorithms are conducted in Chapter 5, through extensive set of coefficient matrices (including rectangular, square, symmetrical, non-symmetrical, singular, non-singular matrices) obtained from well established/popular websites [14-15]. Detailed procedures for executing any FORTRAN code (such as the "educational version" of the developed generalized inverse code for solving SLE) on the internet are explained and demonstrated in Chapter 6. Basic/simple parallel MATLAB-MPI functions (including parallel matrix times matrix operations) are summarized in Chapter 7. Finally, conclusions and future research works are summarized in Chapter 8.

2. DIRECT AND ITERATIVE METHODS FOR SYSTEM OF NON-SINGULAR SLE

Many real life, practical problems in scientific computing require efficient solution of Simultaneous Linear Equations (SLE), which can be conveniently expressed in the matrix notation as

$$Ax = b \tag{2.1}$$

In general, solutions for Eq. (2.1) can be classified into 2 categories: Direct and Iterative methods. In the subsequent sections, basic ideas behind these two types of solution approaches will be briefly summarized and discussed.

2.1 Direct Methods for Solving SLE

Depending on the nature of the coefficient matrix A, shown in Eq. (2.1), different direct methods/algorithms are available, such as:

a) Cholesky algorithm [if matrix A is Symmetric Positive Definite (SPD)]

b) LDL^{T} algorithm [if matrix A is Symmetric, could be either positive or negative definite]

c) LU decomposition algorithm [if matrix A is unsymmetric]

2.1.1 Cholesky Method

If the coefficient matrix A is Symmetric Positive Definite (SPD), then the following three step Cholesky algorithm can be used to obtain the solution for Eq. (2.1) Step1: Matrix Factorization Phase The coefficient matrix A is decomposed into

$$[A] = [U]^T U \tag{2.2}$$

where U is an $n \times n$ upper triangular matrix. For a general $n \times n$ SPD [A], the diagonal and off-diagonal terms of the factorized matrix U can be computed from the following formulas [3, 10-11]

$$u_{ii} = \left(A_{ii} - \sum_{k=1}^{i-1} (u_{ki})^2\right)^{\frac{1}{2}}$$

and

$$u_{ij} = \frac{A_{ij} - \sum_{k=1}^{i-1} u_{ki} u_{kj}}{u_{ii}}$$

<u>Note:</u> Since the "square root" operation is required for computing the diagonal terms of U, positive definite is a requirement for matrix A to assure the number under the square root is positive.

Step2: Forward solution phase

 Substituting Eq. (2.2) in Eq. (2.1)

$$[U]^{T}[U]\{x\} = \{b\}$$
 (2.3)

 Let's define

 $[U]\{x\} = \{y\}$
 (2.4)

 Eq. (2.3) becomes
 (2.5)

The intermediate unknown $\{y\}$ can be easily solved from Eq. (2.5) and hence the name "forward solution".

From Eq. (2.4), the unknown vector $\{x\}$ can be effectively solved and hence the name "backward solution".

The matrix factorization phase (step 1) is the most-time consuming part of solving SLE in the Cholesky algorithm. However, if the right-hand-side (RHS) vector $\{b\}$, shown in Eq. (2.1), becomes a matrix (with multiple columns), then the combined Forward and Backward solution time may become significant (as compared to the matrix factorization phase). As a general rule of thumb, computational time/effort for one matrix factorization is roughly equivalent to 20-25 times the efforts for one Forward and Backward solution phases.

2.1.2 LDL^T Method

In many engineering and science applications, the coefficient matrix in Eq. (2.1) is symmetric, however it may not be positive definite. The coefficient matrix could be negative definite. In this case, LDL^{T} algorithms can be used for solving Eq. (2.1), which also requires the following 3 computational steps

Step1: Matrix Factorization phase

$$[A] = [L][D][L]^T (2.6)$$

In Eq. (2.6), matrices [L] and [D] represent the lower triangular (with 1 on its diagonal) and diagonal matrices, respectively.

Step2: Forward Solution and Diagonal Scaling phase

Substituting Eq. (2.6) in Eq. (2.1), one gets

 $[L][D][L]^{T}\{x\} = \{b\}$ (2.7)

Let's define

$$[L]^{T}\{x\} = \{y\}$$
(2.8)

$$[D]\{y\} = \{z\} \Leftrightarrow \begin{bmatrix} D_{11} & 0 & 0\\ 0 & D_{22} & 0\\ 0 & 0 & D_{33} \end{bmatrix} \begin{pmatrix} y_1\\ y_2\\ y_3 \end{pmatrix} = \begin{pmatrix} z_1\\ z_2\\ z_3 \end{pmatrix}$$
(2.9)

or,
$$y_i = \frac{z_i}{D_{ii}}$$
; for $i = 1, 2, 3, ... N$ (2.10)

Then Eq. (2.7) becomes:

$$[L]\{z\} = \{b\} \Leftrightarrow \begin{bmatrix} 1 & 0 & 0 \\ L_{21} & 1 & 0 \\ L_{31} & L_{32} & 1 \end{bmatrix} \begin{pmatrix} z_1 \\ z_2 \\ z_3 \end{pmatrix} = \begin{pmatrix} b_1 \\ b_2 \\ b_3 \end{pmatrix}$$
(2.11)

or,
$$z_i = b_i - \sum_{k=1}^{i-1} L_{ik} z_k$$
; for $i = 1, 2, 3, ..., N$ (2.12)

Step3: Backward Solution phase

In this step, Eq. (2.8) can be effectively solved for the original unknown vector $\{x\}$

$$[L]^{T}\{x\} = \{y\} \Leftrightarrow \begin{bmatrix} 1 & L_{21} & L_{31} \\ 0 & 1 & L_{32} \\ 0 & 0 & 1 \end{bmatrix} \begin{pmatrix} x_{1} \\ x_{2} \\ x_{3} \end{pmatrix} = \begin{pmatrix} y_{1} \\ y_{2} \\ y_{3} \end{pmatrix}$$
(2.13)

or,
$$x_i = y_i - \sum_{k=i+1}^N for \ i = N, N - 1, \dots 1$$

2.1.3 LU Decomposition Method

LU decomposition can be used to solve Eq. (2.1), when the coefficient matrix A is unsymmetric.

Step1: Factorization phase

The coefficient matrix in Eq. (2.1) can be factorized as a product of two matrices

$$A = L.U \tag{2.14}$$

where L is lower triangular (with values 1 on its diagonal) and U is upper triangular.

For the case of 4×4 matrix A. Eq. (2.14) would look like

$$\begin{bmatrix} a_{11} & a_{12} & a_{13} & a_{14} \\ a_{21} & a_{22} & a_{23} & a_{24} \\ a_{31} & a_{32} & a_{33} & a_{34} \\ a_{41} & a_{42} & a_{43} & a_{44} \end{bmatrix} = \begin{bmatrix} 1 & 0 & 0 & 0 \\ a_{21} & 1 & 0 & 0 \\ a_{31} & a_{32} & 1 & 0 \\ a_{41} & a_{42} & a_{43} & 1 \end{bmatrix} \begin{bmatrix} \beta_{11} & \beta_{12} & \beta_{13} & \beta_{14} \\ 0 & \beta_{22} & \beta_{23} & \beta_{24} \\ 0 & 0 & \beta_{33} & \beta_{34} \\ 0 & 0 & 0 & \beta_{44} \end{bmatrix}$$
(2.15)

Various terms inside matrices [L] and [U] can be computed by equating both sides of Eq.

(2.15).

Step2: Forward solution phase

Substituting Eq. (2.14) into Eq. (2.1), one obtains

$$[L.U]\mathbf{x} = \mathbf{b} \tag{2.16}$$

Let us define

$$[U]\{x\} = y \tag{2.17}$$

Substituting Eq. (2.17) in Eq. (2.16), one gets

$$[L]\{y\} = b \tag{2.18}$$

In this "forward solution" phase, Eq. (2.18) can be easily solved for the "intermediate" unknown vector $\{y\}$.

Step3: Backward solution phase

Solving the unknown vector $\{x\}$ in Eq. (2.17) is called "backward solution" phase

 $[U]{x} = y$ (2.17, repeated)

While direct methods have offered advantages in terms of its robustness, accuracy, and reliability, etc., for large/sparse SLE (especially for 3-D problems), these direct methods may become excessively expensive. Furthermore, direct methods also have the following limitations:

a) The amount of required computer memory can be high.

b) The operation counts can be high, especially when many non-zero fill-in terms occurred during the factorization phase, even though reordering algorithms have commonly used (to minimize the non-zero fill-in terms) prior to numerical factorization phase, and

c) These methods have low degree of parallelism (or not easy to parallelize).

The above-mentioned drawbacks have motivated researchers to investigate iterative methods as possible alternative choices.

2.2 Iterative Methods for Solving SLE

Iterative methods can be superior to direct methods in all the above mentioned three aspects. Some of the popular iterative methods are Conjugate Gradient (CG), Bi-Conjugate Gradient (Bi-CG) with or without stabilizers [1, 3, 10], Generalized Minimal Residual (GMRES) [1, 3, 10], etc. These methods (CG for symmetrical, while Bi-CG and GMRES for unsymmetrical systems of SLE) can lead to low memory requirement and make effectively use of parallelism. Most (if not all) existing iterative algorithms require "matrix times vector" and "dot product of 2 vectors" operations. For these reasons, iterative methods are much more easier to parallelize (for improving computational efficiency) as compared to direct methods. These advantages make iterative linear system solvers as attractive alternatives to direct methods, particularly for large (3-D) problems. Despite of these desirable features, iterative methods may also have difficulties for fast convergence (or even have divergence) to a specified (small) error tolerance etc..., unless these iterative methods were used in conjunction with "efficient preconditioned" algorithms [1-3, 10-11] !

2.2.1 Conjugate Gradient (CG) Algorithm with Preconditioner

For systems of Symmetrical Positive Definite (SPD) SLE, the Preconditioned Conjugate Gradient (PCG) algorithms can be considered as the method of choice. PCG algorithms can be summarized in the following step-by-step numerical procedures [3, 10-

11], for solving Eq. (2.1)

Eq. (2.1) can be re-casted as:

$$PAP^T P^{-T} \vec{x} = P \vec{b} \tag{2.19}$$

Eq. (2.19) can be expressed in the following general form [1, 3, 10-11]:

$$[A^*]\vec{y} = \vec{b^*} \tag{2.20}$$

where matrix

 $[A^*] = [P] \times [A] \times [P^T] =$ symmetrical matrix, and the right-hand-side vector $\{b^*\}$ is defined as

$$\{b^*\} = [P] \times \{b\}$$
(2.21)

$$\vec{y} = P^{-T} \vec{x} \tag{2.22}$$

and [P]= preconditioned matrix

The step-by-step PCG algorithm is summarized in Fig. 2.1

Given the initial guessed vector $\vec{x}^{(0)}$ Compute (or input) the preconditioned matrix [P]Compute $\vec{r}^{(0)} = \{Pb\} - |PAP^{T}|\vec{x}^{(0)}|$ Set $\vec{a}^{(0)} = 0; \rho_{-1} = 1$ **Do** i = 1, 2, ... $\rho_{i-1} = \left\{ r^{(i-1)} \right\}^{T} \left\{ r^{(i-1)} \right\}$ $\beta_{i-1} = \frac{\rho_{i-1}}{\rho_{i-2}}$ $a^{(i)} = r^{(i-1)} + \beta_{i-1} a^{(i-1)}$ $q^{(i)} = \left| PAP^{T} \right| d^{(i)}$ $\alpha_i = \frac{\rho_{i-1}}{\left\{d^{(i)}\right\}^r \left\{q^{(i)}\right\}}$ $x^{(i)} = x^{(i-1)} + \alpha_i a^{(i)}$ $r^{(i)} = r^{(i-1)} - \alpha_i q^{(i)}$ Converge?? End do

If converged, then set

 $\vec{x} = \left[P \right]^r \vec{x}$

Figure 2.1 Preconditioned Conjugate Gradient Algorithm

2.2.2 GMRES Algorithm [1, 3, 10, 16]

For systems of "unsymmetrical" SLE, popular algorithms such as GMRES, Bi-Conjugate Gradient (Bi-CG), Bi-Conjugate Gradient with Stabilizers (Bi-CG Stab) [1, 3, 10, 16] are recommended. For readers' convenience, a version of GMRES algorithms can be summarized in the following step-by-step numerical procedures [1, 3, 11], for solving $[A^*]\vec{y} = \vec{b^*}$

$$r_{0} = b - Ax_{0}$$

$$v_{1} = \frac{r_{0}}{|r_{0}|_{2}}$$
start
for $j = 1 : m$

$$w = Av_{j}$$
for $i = 1 : j$

$$H(i, j) = (w, v_{i})$$

$$w = w - H(i, j)v_{i}$$
end
$$v_{j+1} = \frac{w}{|w|_{2}}$$
if $||v_{j+1}||_{2} < tolerance$ Break
end
$$V_{j+1} = [v_{1}, v_{2}, v_{3}, \dots, v_{j+1}]$$

$$H = V_{j}^{T} A V_{j}$$

$$y = \arg \min ||V_{j+1}r_{0} - Hy|_{2}$$

$$x = x_{0} + V_{j}y$$
if $||Ax - b|_{2} < tolerance$ Stop
else $x_{0} = x$ and goto start

Figure 2.2 GMRES Algorithm

3. DOMAIN DECOMPOSITION SOLVER

Domain decomposition [1-2, 4, 11, 16] algorithm is a powerful method for solving large scale system of equations arising from discretization of partial differential equations (PDE) in finite element procedures. The computational domain is decomposed into smaller sub-domains each of which is easier to solve.

Domain decomposition (DD) is an application of the divide-and-conquer problem-solving strategy, which consists of expressing a large problem as a set of smaller sub-problems defined on sub-domains and provides a way to determine the solution of the original problem in terms of solutions to sub-problems.

The goal of DD is to divide the original problem into sub-problems that can be solved independently. A critical issue in DD is to assure that the sub-problems preserve the solution to the original problem.

Let us assume the system of linear algebraic equations

$$Kr = f \tag{3.1}$$

where the matrix of system is $K \in \mathbb{R}^{n \times n}$ and vectors $r \in \mathbb{R}^n$, $f \in \mathbb{R}^n$. It is possible to split Eq. (3.1) into blocks (or sub-matrices)

$$\begin{bmatrix} K_{11} & K_{12} \\ K_{21} & K_{22} \end{bmatrix} \begin{bmatrix} r^{(1)} \\ r^{(2)} \end{bmatrix} = \begin{bmatrix} f^{(1)} \\ f^{(2)} \end{bmatrix}$$
(3.2)

Eq. (3.2) can be written as

$$K_{11}r^{(1)} + K_{12}r^{(2)} = f^{(1)}$$
(3.3)

$$K_{21}r^{(1)} + K_{22}r^{(2)} = f^{(2)}$$
(3.4)

From Eq. (3.3), the vector $r^{(1)}$ can be expressed in the form

$$r^{(1)} = K_{11}^{-1} \left(f^{(1)} - K_{12} r^{(2)} \right) \tag{3.5}$$

Substituting Eq. (3.5) in Eq. (3.4)

$$(K_{22} - K_{21}K_{11}^{-1}K_{12})r^{(2)} = f^{(2)} - K_{21}K_{11}^{-1}f^{(1)}$$
(3.6)

From Eq. (3.6) we can observe that the number of unknowns have been reduced as this matrix equation is only related to the unknown vector $r^{(2)}$.

The matrix $(K_{22} - K_{21}K_{11}^{-1}K_{12})$ is often referred as the Schur's complement in the mathematical community.

The basic ideas in DD solver is to solve for the unknown vector $\{r\}$, shown in Eq. (3.1), in the following 2 major steps:

Step 1: The "boundary unknown" vector $\{r^{(2)}\}$ is solved from Eq. (3.6). Since the triple matrix products appeared in Eq. (3.6) is usually dense, and computationally expensive (because K_{12} is a matrix, and <u>not</u> a vector), iterative solver (which is based on matrix times vector operations) is usually recommended in this step.

Step 2: The "interior unknown" vector $\{r^{(1)}\}$ is solved from Eq. (3.5). Since the coefficient matrix $[K_{11}]$ is usually sparse (and $f^{(1)}$ is a vector, not a matrix), direct solver (such as Cholesky, or LDL^{T} , or LU algorithms) is strongly recommended.

4. GENERALIZED INVERSE ALGORITHMS FOR SINGULAR/NON-SINGULAR, SQUARE/RECTANGULAR SYSTEM OF SLE

In Chapter 2, various direct and iterative methods for solving square/non-singular system of SLE have been summarized. The (direct and iterative) methods described in Chapter 2 can be significantly enhanced/improved by the Domain Decomposition (DD) formulation [1-2, 11, 16], described in Chapter 3. Domain decomposition formulation has been widely adopted, since it can take full advantage of "parallel processing" capability offered by most (if not all) today super-computers, and even to desktop/laptop computers, which have multiple processors. In this Chapter 4, however, the main focus is shifted into algorithms (numerical procedures) that can solve much more general classes of SLE, shown in Eq. (2.1), for which the coefficient matrix [A] can be square/rectangular, non-singular/singular, well-posed/illed condition. These desirable algorithms have been based on the so called Generalized (or Pseudo, or Moore-Penrose) inverse of a general matrix [5-9, 11-13, 17-18]. Furthermore, DD formulation will also be used in this chapter to further improve the numerical performance of the generalized inverse.

4.1 Basic Conditions for the Generalized Inverse

The Moore-Penrose inverse (or generalized inverse or pseudo inverse) of a $m \times n$ matrix K (not necessarily a square matrix) is the unique $n \times m$ matrix K⁺ which satisfies the following four conditions:

- 1. General condition: $KK^+K = K$,
- 2. Reflexive condition: $K^+KK^+ = K^+$,
- 3. Normalized condition: $(KK^+)' = KK^+$,

4. Reverse normalized condition: $(K^+K)' = K^+K$

4.2 Potential Engineering/Science Generalized Inverse Applications

There are some applications that result in (or lead to) an inconsistent system of SLE. A solution may not exist in this case. However, we can consider to fit a vector \vec{x} to a given inconsistent system. That is, we can define a least-squares error problem for finding \vec{x} that minimizes the absolute error $||A\vec{x} - \vec{b}||_2$ which is equivalent to minimizing the summation of the square of the error $||A\vec{x} - \vec{b}||_2^2$. As should be obvious, such a solution vector \vec{x} may not be unique, and it can be obtained/computed by $\vec{x} = A^+\vec{b}$, where A^+ is generalized (or pseudo-) inverse of A.

It has been well documented in the literature that many real-life engineering/science/statistical applications can be efficiently solved by efficient computation of the generalized inverse in conjunction with SLE. In the following section, it can be shown that the popular "least square problem" can be formulated such that generalized inverse algorithms can be incorporated for obtaining the desired solution.

4.3 Least Squares Curve Fitting Problem

Let us assume we are given a set of data points $(x_1, y_1), (x_2, y_2), ..., (x_n, y_n)$, and we wish to find parameters $c_1, c_2, ..., c_k$ to be multiplied by the basis functions $\emptyset_1, \emptyset_2, ..., \emptyset_k$ such that the scalar function S, defined as

$$S = \sum_{i=1}^{n} |\varepsilon_i|^2 \tag{4.1}$$

is minimized, where the errors ε_i can be computed by

$$\varepsilon_{1} = c_{1}\phi_{1}(x_{1}) + c_{2}\phi_{2}(x_{1}) + \dots + c_{k}\phi_{k}(x_{1}) - y_{1}$$

$$\varepsilon_{2} = c_{1}\phi_{1}(x_{2}) + c_{2}\phi_{2}(x_{2}) + \dots + c_{k}\phi_{k}(x_{2}) - y_{2}$$

$$\vdots \qquad \vdots \qquad \vdots \qquad \vdots$$

$$\varepsilon_{n} = c_{1}\phi_{1}(x_{n}) + c_{2}\phi_{2}(x_{n}) + \dots + c_{k}\phi_{k}(x_{n}) - y_{n}$$
(4.2)

In Eq. (4.1), the scalar parameter S represents the "summation of the square of the errors". Ideally, we would like to set Eq. (4.1) to zero, i.e., each of the ε_i to be zero. Otherwise, we would like to minimize the scalar function S. If c and y are the vectors

$$\begin{bmatrix} c_1 \\ c_2 \\ \vdots \\ c_k \end{bmatrix} and \begin{bmatrix} y_1 \\ y_2 \\ \vdots \\ y_n \end{bmatrix},$$
(4.3)

respectively, and A is the $n \times k$ matrix

$$A = \begin{bmatrix} \phi_1(x_1) & \phi_2(x_1) & \cdots & \phi_k(x_1) \\ \phi_1(x_2) & \phi_2(x_2) & \cdots & \phi_k(x_2) \\ \vdots & \vdots & \vdots & \vdots \\ \phi_1(x_n) & \phi_2(x_n) & \cdots & \phi_k(x_n) \end{bmatrix}$$
(4.4)

Then, the Right-Hand-Side (or RHS) of Eq. (4.2) can also be written in the form

$$\begin{pmatrix} \phi_{1}(x_{1}) & \phi_{2}(x_{1}) & \cdots & \phi_{k}(x_{1}) \\ \phi_{1}(x_{2}) & \phi_{2}(x_{2}) & \cdots & \phi_{k}(x_{2}) \\ \vdots & \vdots & \vdots & \vdots \\ \phi_{1}(x_{n}) & \phi_{2}(x_{n}) & \cdots & \phi_{k}(x_{n}) \end{pmatrix}_{n \times k} \begin{pmatrix} c_{1} \\ c_{2} \\ \vdots \\ c_{k} \end{pmatrix}_{k \times 1} = \begin{pmatrix} y_{1} \\ y_{2} \\ \vdots \\ y_{n} \end{pmatrix}_{n \times 1}$$
(4.5)

The requirement S = 0 is equivalent to

$$Ac = y \tag{4.6}$$

From Eq. (4.5), if n > k, we have an over-determined system of linear equations, since the number of equations is larger than that of the unknowns. It is generally not possible to find a solution to this system, but we may find $c_1, c_2, ..., c_k$ such that Ac is close to y (in the least squares sense). If there is a solution c^+ of the least squares problem, then we write

$$c^+ = A^+ y \tag{4.7}$$

The matrix A^+ is called the pseudo-inverse (or generalized inverse) of A. We know that when n = k and A is invertible (i.e., A has an inverse), then

$$A^{+} = A^{-1} \tag{4.8}$$

4.3.1 The Normal Equations [3, 10, 19]

In order to minimize S, we must minimize

$$S_{\min} \equiv \sum_{i=1}^{n} \left[c_1 \phi_1(x_i) + c_2 \phi_2(x_i) + \dots + c_k \phi_k(x_i) - y_i \right]^2$$
(4.9)

If we take for j = 1, 2, ..., k the partial derivatives of S with respect to c_j and set them equal to zero, we will get the following normal equations

$$2\sum_{i=1}^{n} \left[c_1 \phi_1(x_i) + c_2 \phi_2(x_i) + \dots + c_k \phi_k(x_i) - y_i \right] \phi_j(x_i) = 0$$
(4.10)

or, equivalently,

$$\sum_{i=1}^{n} \left[c_{1}\phi_{1}(x_{i}) + c_{2}\phi_{2}(x_{i}) + \dots + c_{k}\phi_{k}(x_{i}) \right] \phi_{j}(x_{i}) = \sum_{i=1}^{n} \phi_{j}(x_{i}) y_{i}$$
(4.11)

The above normal equations can also be expressed (in matrix notations) as

$$A^T A c = A^T y \tag{4.12}$$

where the matrix A has the entries $a_{ij} = \phi_j(x_i)$.

As an example, let k = 4 and $\phi_j(x) = x^{j-1}$, where j = 1, 2, ..., k = 4

Then,

$$\phi_1(x) = 1$$

$$\phi_2(x) = x$$

$$\phi_3(x) = x^2$$

$$\phi_4(x) = x^3$$

and

$$A = \begin{pmatrix} 1 & x_1 & x_1^2 & x_1^3 \\ 1 & x_2 & x_2^2 & x_2^3 \\ \vdots & \vdots & \vdots & \vdots \\ 1 & x_n & x_n^2 & x_n^3 \end{pmatrix}$$
(4.13)

$$A^{T} = \begin{pmatrix} 1 & 1 & \cdots & 1 \\ x_{1} & x_{2} & \cdots & x_{n} \\ x_{1}^{2} & x_{2}^{2} & \cdots & x_{n}^{2} \\ x_{1}^{3} & x_{2}^{3} & \cdots & x_{n}^{3} \end{pmatrix}$$
(4.14)

so that

$$A^{T}A = \begin{pmatrix} n & \sum x_{i} & \sum x_{i}^{2} & \sum x_{i}^{3} \\ \sum x_{i} & \sum x_{i}^{2} & \sum x_{i}^{3} & \sum x_{i}^{4} \\ \sum x_{i}^{2} & \sum x_{i}^{3} & \sum x_{i}^{4} & \sum x_{i}^{5} \\ \sum x_{i}^{3} & \sum x_{i}^{4} & \sum x_{i}^{5} & \sum x_{i}^{6} \end{pmatrix}$$
(4.15)

where in the summations, *i* runs from 1 to n and

$$A^{T} y = \begin{pmatrix} \sum y_{i} \\ \sum x_{i} y_{i} \\ \sum x_{i}^{2} y_{i} \\ \sum x_{i}^{3} y_{i} \end{pmatrix}$$
(4.16)

If the matrix $A^{T}A$ in Eq. (4.15) is invertible, then there is a solution to the least squares problem.

We then have

$$c = \left(A^T A\right)^{-1} A^T y, \qquad (4.17)$$

so that the pseudo-inverse A^+ (of the given matrix A, see Eqs. 4.6-4.7) is given by

$$A^{+} = (A^{T} A)^{-1} A^{T}$$
(4.18)

In subsequent sections, Eq. (4.18) can be generalized into the form shown in Eq. (4.27).

4.4 Special Cholesky Algorithms for Factorizing a Singular Matrix

The stiffness matrix of the "floating" sub-domain is singular due to the fact that there are not enough (support) constraints to prevent its rigid body motion. To facilitate the discussions, let's assume the "floating" sub-domain's stiffness matrix is given as:

$$\begin{bmatrix} K_{float} \end{bmatrix} = \begin{bmatrix} 1 & 2 & -3 & 2 & -2 & -3 & -2 \\ 2 & 4 & -6 & 4 & -4 & -6 & -4 \\ -3 & -6 & 9 & -6 & 6 & 9 & 6 \\ 2 & 4 & -6 & 5 & -1 & -5 & -7 \\ -2 & -4 & 6 & -1 & 13 & 9 & -5 \\ -3 & -6 & 9 & -5 & 9 & 13 & 9 \\ -2 & -4 & 6 & -7 & -5 & 9 & 27 \end{bmatrix}$$
(4.19)

 $row2 = 2 * row1 \tag{4.20}$

$$row3 = -3 * row1 \tag{4.21}$$

$$row5 = -8 * row1 + 3 * row4$$
 (4.22)

In Eq. (4.19), it can be observed that rows number 2, 3 and 5 are dependent rows (and columns). Thus, the above matrix is singular, and the regular/standard Cholesky factorization algorithms will not work. To facilitate the development of efficient "generalized inverse" algorithms (and its applications) in subsequent chapters, "special

Cholesky" factorization algorithm is needed. The "special Cholesky factorization" algorithm is essentially identical to the regular/standard one, with the following two modifications (or differences):

- (a) During the numerical factorization phase, if the dependent row(s) is/are detected, then these dependent row(s) is/are skipped!
- (b) Factorization of the current i^{th} row, in general, will require the previously already factorized rows k = 1, 2, ..., i 1. However, if the previous k^{th} row was amongst the dependent row(s), then the contribution from this k^{th} row to this current i^{th} row will be ignored.
- (c) Having obtained the "special Cholesky" factorized square matrix, those factorized/dependent rows will be deleted to obtain the so-called "truncated Cholesky factorized rectangular matrix U^* ".

Using MATLAB, the Eigen-values of matrix in Eq. (4.19) can be computed as:

 $\vec{\lambda} = \{0.0000, 0.0000, 0.0000, 0.2372, 4.9375, 24.9641, 41.8612\}^T$

Since there are 3 zero Eigen-values, it implies there are 3 rigid body modes (or 3 dependent rows/columns) in Eq. (4.19)

If row-by-row Cholesky factorization scheme is applied to Eq. (4.19), we will encounter that the factorized $u_{22} = 0 = u_{33} = u_{55}$, which indicated that row number 2, 3 and 5 are dependent rows. Thus, if we set all factorized terms of rows number 2, 3 and 5 are zero, and "ignoring" these three rows in the factorization of subsequent rows, one obtains the following Cholesky factorized matrix [U] of a given matrix [K_{float}], shown in Eq. (4.19):

$$[U] = \begin{bmatrix} 1 & 2 & -3 & 2 & -2 & -3 & -2 \\ \cdot & 0 & 0 & 0 & 0 & 0 & 0 \\ \cdot & \cdot & 0 & 0 & 0 & 0 & 0 \\ \cdot & \cdot & \cdot & 1 & 3 & 1 & -3 \\ \cdot & \cdot & \cdot & 0 & 0 & 0 \\ \cdot & \cdot & \cdot & \cdot & 1.7321 & 3.464 \\ \cdot & \cdot & \cdot & \cdot & \cdot & 1.4145 \end{bmatrix}$$

Based on the above "special Cholesky factorization" algorithm, and using the 7×7 singular matrix data as shown in Eq. (4.19), the "truncated Cholesky factorized matrix U^* " (corresponding to the independent rows # 1, 4, 6, and 7) of the product $[K_{float}]^T * [K_{float}]$ can be obtained/computed as: For row #1 of the truncated 4×7 matrix U^* : 5.9160 11.8321 - 17.7482 11.6631 - 12.3392 - 21.4669 - 19.1004 For row #4 of the truncated 4×7 matrix U^* : 0.0000 0.0000 0.0000 4.4689 13.4068 0.9781 - 21.7565 For row #6 of the truncated 4×7 matrix U^* : 0.0000 0.0000 0.0000 0.0000 4.4960 10.0650 For row #7 of the truncated 4×7 matrix U^* :

 $0.0000 \quad 0.0000 \quad 0.0000 \quad 0.0000 \quad 0.0000 \quad 0.0000 \quad 0.7209$

4.5 Special LDL^T Algorithms for Factorizing a Singular Matrix

If the square/symmetrical/singular coefficient matrix is NOT positive definite, then the symmetrical, floating stiffness matrix (shown in Eq. 4.19) can be factorized by the familiar sparse LDL^{T} algorithms [11], with the following minor modifications:

(a) Whenever a dependent i^{th} row is encountered (such as the factorized $u_{ii} = 0$), then the following things need be done: (a.1) Recording the dependent row number(s). For the data given by Eq. (4.19), the dependent rows are rows number 2, 3 and 5.

(a.2) Set all the nonzero terms of the factorized i^{th} row (of L^{T}) to be zero

(a.3) Set
$$\frac{1}{u_{ii}} \equiv D_{ii} = 0$$

(b) Whenever an independent i^{th} row is encountered, the factorized i^{th} row will have contributions from all appropriated previously factorized rows. However, contributions from the previously factorized (dependent) rows will be ignored.

(c) Finally, the truncated/rectangular LDL^{T} factorized matrix U^{*} can be obtained by deleting those dependent row(s) from the "special LDL^{T} factorized" square matrix.

As an example, the (LDL^{T}) factorized matrix (for the matrix data $[K_{float}]$ shown in Eq. (4.19)) can be computed as

	г 1	2	-3	2	-2	-3	-2-
	.	0	0	0	0	0	0
			0	0	0	0	0
[U] =			•	1	3	1	-3
	.				0	0	0
	.					0.333	2
	L,	•				-3 0 1 0 0.333	0.5

and the truncated factorized LDL^{T} matrix U^{*} of the product $[K_{float}]^{T} * [K_{float}]$ can be obtained by deleting rows #2, 3, and 5 of the above matrix U.

A complete educational version of "special LDL^{T} " software code (written in FORTRAN language) is listed/given in the Appendix B.

4.6 Efficient Generalized Inverse Algorithms

Moore-Penrose inverse can be computed using Singular Value Decomposition (SVD) [10-13], Least Squares Method, QR factorizations, Finite Recursive Algorithm [9], etc. In this work, our numerical algorithms have been based on:

(a) The "special Cholesky factorization" (for symmetrical/singular coefficient matrix) see section 4.4 and the Appendix C, and

(b) The generalized inverse of a product of 2 matrices [5, 13]

and can be described in the following paragraphs.

Consider Eq. (2.1) $[G]\vec{x} = \vec{b}$, with a square coefficient $n \times n$ matrix, and let the rank be less than the size of the matrix (if r is the rank of the matrix, then $r \leq n$). Let the size of the known right-hand-side vector \vec{b} be $n \times 1$. Consider a symmetric positive $n \times n$ matrix G'G, with rank $r \leq n$ (here, the matrix [G] plays the same role as matrix [A] in Eq. (2.1)), then based on the theorem presented in [5, 13, 17-18], there exists a unique [M] such that:

$$G'G = M'M \tag{4.24}$$

In Eq. (4.24), matrices [G'] and [G] have the dimensions $n \times m$ and $m \times n$, respectively.

M is the upper triangular (special) Cholesky factorized matrix and contains exactly n-r zero rows. Removing the zero rows from *M*, one obtains a $r \times n$ (upper, rectangular) matrix *L'*.

$$A \equiv M'M = LL' \tag{4.25}$$

In this work, the upper triangular (special) Cholesky factorized matrix [M] can be obtained by the regular/standard Cholesky factorization, with the following modifications:

a) When the diagonal term of the current i^{th} row is very close to zero, then factorization of this dependent row is skipped.

b) When the current i^{th} row is factorized, all previous rows k = 1, 2, ..., i - 1 were used except those dependent row(s).

Consider the generalized inverse of a matrix product AB [5, 13, 17-18]

$$(AB)^{+} = B'(A'ABB')^{+}A'$$
(4.26)

From Eq. (4.26), if B = I then

$$A^{+} = (A'A)^{+}A' \tag{4.27}$$

Eq. (4.27) can be considered as a general version of the earlier Eq. (4.18).

If B = A' and A is a $n \times r$ matrix of rank r, then one obtains from Eq. (4.26)

$$(AA')^{+} = (A')' \Big(A'AA'(A')' \Big)^{+} A'$$
(4.28)

Let us consider regular inverse in Eq. (4.28) in place of generalized inverse

$$(AA')^{+} = A(A'AA'A)^{-1}A'$$

= $A(A'A)^{-1}(A'A)^{-1}A'$ (4.29)

Using Eq. (4.27),

$$G^{+} = (G'G)^{+}G'$$
(4.30)

From Eqs. (4.24 - 4.25) and Eq. (4.29) one obtains,

$$(G'G)^{*} = (LL')^{*} = L(L'L)^{-1}(L'L)^{-1}L'$$
(4.31)

Thus, Eq. (4.30) becomes

$$G^{+} = (G'G)^{+}G' = L(L'L)^{-1}(L'L)^{-1}L'G'$$
(4.32)

While MATLAB solution can be obtained by $\vec{x} = pinv(G) \times \vec{b}$, implying the generalized inverse G^+ [see Eq. 4.32] to be formed explicitly, our main idea is to solve SLE where \vec{b} is a known right-hand-side vector, as described in the next section.

To facilitate the discussion of Generalized Inverse, and its usage for solving general systems of SLE, the following (small-scale) numerical examples are used:

Example 4.1

The coefficient matrix [G] is a rectangular (tall type) matrix, and the RHS vector $\{b\}$ is a linear combinations of columns of [G].

In this example, we wish to solve for $\{x\}$ from the SLE $[G] * \{x\} = \{b\}$, where the numerical values of the coefficient matrix [G], and the RHS vector $\{b\}$ are given as [refer to Eq. (4.32)]:

٢2	0	0	1	ן 1	(7.)	r 6 1
-1	0	0	-2	-1	$\begin{pmatrix} \mathbf{z}_1 \\ \mathbf{z}_2 \end{pmatrix}$	-5
3	0	0	0	0	$\int_{r_0}^{r_2}$	6
0	1	2	-3	1	$\begin{pmatrix} \lambda_3 \\ \gamma_1 \end{pmatrix} =$	1
0	-2	5	1	0	$\begin{pmatrix} \lambda_4 \\ \chi \end{pmatrix}$	4
l 1	-2	3	1 -2 0 -3 1 4	_1J	$ \begin{cases} x_1 \\ x_2 \\ x_3 \\ x_4 \\ x_5 \end{cases} =$	L 6 J

Using MATLAB built-in function, the rank of G can be computed as

rank(G) = 5

and
$$G^T * G = \begin{bmatrix} -15 & -2 & 3 & 8 & 2 \\ -2 & 9 & -14 & -13 & 3 \\ 3 & -14 & 38 & 11 & -1 \\ 8 & -13 & 11 & 31 & -4 \\ 2 & 3 & -1 & -4 & 4 \end{bmatrix}$$

Special Cholesky factor of $G^T * G$

	[3.8729]	-0.5163	0.7745	2.0655	ן 0.5163	
	0.0000	2.9552	-4.6020	-4.0380	1.1053	
=	0.0000	0.0000	4.0275	-2.2800	0.9154	
	0.0000	0.0000	0.0000	2.2866	0.64909	
	0.0000	0.0000	0.0000	0.0000	1.1189]	

of independent rows = 5

Independent rows = 12345

Dependent row = 6

Product of $L^T * L$

	20.4000	-12.8609	-1.1172	5.0584	ן0.5778
	-12.8609	47.4396	-8.3159	-8.5160	1.2368
I	-1.1172	-8.3159	22.2581	-4.6195	1.0243
	5.0584	-8.5160	-4.6195	5.6500	0.7263
	L 0.5778	1.2368	1.0243	0.7263	1.2520 ^J

Regular Cholesky factorization of $L^T * L$

	4.5166	-2.8474	-0.2473	1.1199	ן0.1279
	0.0000	6.2714	-1.4383	-0.8494	0.2553
=	0.0000	0.0000	4.4864	-1.2402	0.3172
	0.0000	0.0000	0.0000	1.4615	0.8164
	L0.0000	0.0000	0.0000	0.0000	0.6350J

Generalized inverse of [G]

	r0.07957		0.2930	0.0117	-0.0206	ן 0.0265
	-0.4289	-0.3494	-0.1162	0.9364	-0.8888	0.8570
=	-0.1342	-0.0723	-0.0308	0.3504	-0.1133	0.2885
	-0.0645	-0.3494 -0.0723 -0.2377	-0.1401	0.1385	-0.2425	0.3118
1	L 0.4337	-0.2761	-0.2072	-0.2320	0.4060	-0.5220

and the solution vector \vec{x} is obtained as

$\int x_1$)	(2.0000)
x_2		0.9999
$\{x_3$	$\rangle = \langle$	0.9999
$ x_4 $		0.9999
x_5)	(1.0000)

Example 4.2

The coefficient matrix [G] is a square/singular matrix, and the RHS vector $\{b\}$ is a

random vector.

٢2	-1	0	0	1	ן1–	(x_1)		r 0 1
4	-2	0	0	1 2	-2	<i>x</i> ₂		0
0	0	1	2	0	2	$\int x_3$		7
0	0	-2	5	-1	1) X ₄	(4
1	3	0		1	3	<i>x</i> 5		12
l_1^-	1	0	2	5	$ \begin{bmatrix} -1 \\ -2 \\ 2 \\ 1 \\ 3 \\ 7 \end{bmatrix} $	(x_6)		L23J

Using MATLAB built-in function, the rank of G can be computed as

rank(G) = 5

	r 22	-6	0	3	16	ך 0
	-6	15	0	5	3	21
$C^T \star C =$	0	0	5	-8	2	0
u +u	3	5	-8	34	6	26
	16	3	2	6	32	32
$G^T * G =$	LO	21	0	26	32	ل68

Special Cholesky factor of $G^T * G$

	4.6904	-1.2792	0.0000	0.6396	0.3411	ן0.0000
	0.0000	3.6556	0.0000	1.5915	2.0143	5.7445
=	0.0000	0.0000	2.2360	-3.5777	0.8944	0.0000
	0.0000	0.0000	0.0000	4.2729	0.8921	3.9451
	L0.0000	0.0000	0.0000	0.0000	3.8353	4.4086

of independent rows = 5

Independent rows = 12345

Dependent row = 6

Product of $L^T * L$

	35.6 8 18]	3.2129	0.7627	5.7764	ן13.0832
	3.2129	52.9542	-3.8924	31.2607	33.0513
=	0.7627	-3.8924	18.6000	-14.4892	3.4304
	5.7764	31.2607	-14.4892	34.6177	20.8144
	L _{13.0832}	33.0513	3.4304	20.8144	34.1462 []]

Regular Cholesky factorization of $L^T * L$

	5.9734]	0.5378	0.1276	0.9670	ן2.1902
	0.0000	7.2570	-0.5458	4.2359	4.3920
=	0.0000	0.0000	4.2761	-2.8765	1.2974
	0.0000	0.0000	0.0000	2.7321	1.3996
	L0.0000	0.0000	0.0000	0.0000	2.5331

Generalized inverse of [G]

	ſ 0.0749	0.1498	0.0823	-0.0161	0.1882	–0.0742ן
	-0.0107	-0.0215	-0.1076	-5.146×10^{-3}	0.3118	-0.0879
_	0.0182	0.0364	0.5157	-0.1939	0.0305	-0.0980
	0.0151	0.0302	0.1518	0.1278	-7.108×10^{-4}	-0.0271
	0.0150	0.0301	-0.1820	-3.678×10^{-3}	-0.0793	0.1366
	L-0.0242	-0.0485	0.0902	-0.0308	-0.0145	0.0761 J

and the solution vector \vec{x} is obtained as

$$\begin{pmatrix} x_1 \\ x_2 \\ x_3 \\ x_4 \\ x_5 \\ x_6 \end{pmatrix} = \begin{cases} 1.0638 \\ 0.9459 \\ 0.9465 \\ 0.9423 \\ 0.9029 \\ 2.0845 \end{cases}$$

Example 4.3

The coefficient matrix [G] is a square/non-singular matrix, and the RHS vector

 $\{b\}$ is a linear combination of columns of [G].

<u>٢</u> 2	-1	0	0	1	ר_1	(x1)		[1]	I
4	-2	0	0	-2	-2	<i>x</i> ₂		-2	
0	0	1	2	0	2	$\int x_3$	L_	5	
0	0	-2	5	-1	1	x_4	(=)	3	
1	3	0	1	1	3	<i>x</i> 5		9	
l_1	3	0	1	1 -2 0 -1 1 6	1 J	(x ₆)		L ₁₂	

Using MATLAB built-in function, the rank of G can be computed as

rank(G) = 6

$$G^{T} * G = \begin{bmatrix} 22 & -4 & 0 & 2 & 1 & -6 \\ -4 & 23 & 0 & 6 & 24 & 17 \\ 0 & 0 & 5 & -8 & 2 & 0 \\ 2 & 6 & -8 & 31 & 2 & 13 \\ 1 & 24 & 2 & 2 & 43 & 11 \\ -6 & 17 & 0 & 13 & 11 & 20 \end{bmatrix}$$

Special Cholesky factor of $G^T * G$

	[4.6904	-0.8528	0.0000	0.4204	0.2132	–1.2792 ₇
	0.0000	4.7193	0.0000	1.3483	5.1239	3.3709
_	0.0000	0.0000	2.2360	-3.5777	0.8944	0.0000
Π	0.0000	0.0000	0.0000	4.0249	-0.4472	2.2360
	0.0000	0.0000	0.0000	0.0000	3.9623	-1.2618
	L0.0000	0.0000	0.0000	0.0000	0.0000	0.6384 J

of independent rows = 6

Independent rows = 123456

Dependent row = none

Product of $L^T * L$

	24.5909	-6.6695	-1.3348	-1.2395	2.4589	-0.8167
=	-6.6695	61.7090	-0.2412	10.6735	16.0488	2.1522
	-1.3348	-0.2412	18.6000	-14.8000	3.5440	0.0000
	-1.2395	10.6735	-14.8000	21.4000	-4.5936	1.4276
	2.4589	16.0488	3.5440	-4.5936	17.2923	-0.8056
	L-0.8167	2.1522	0.0000	1.4276	-0.8056	0.4076 []]

Regular Cholesky factorization of $L^T * L$

	F4.9589	-1.3449	-0.2691	-0.2499	0.4958	–0.1646 ן
	0.0000	7.7390	-0.0779	1.3356	2.1597	0.2494
_	0.0000	0.0000	4.3036	-3.4303	0.8936	-5.78×10^{-3}
-	0.0000	0.0000	0.0000	2.7903	-1.5370	0.3703
	0.0000	0.0000	0.0000	0.0000	3.0365	-0.2266
	L0.0000	0.0000	0.0000	0.0000	0.0000	0.36012 J

Generalized inverse of [G]

	Г 0.2341	0.0972	-0.0317	-0.0158	0.1825	–0.0396 ₁
	-1.2817	0.5694	-0.0634	-0.0317	-0.1349	0.4206
	-1.2222	0.6111	0.5555	-0.2222	-0.4444	-0.4444
	-0.6388	0.3194	0.2222	0.1111	-0.2777	0.2777
	0.5000	-0.2500	0.0000	0.0000	0.0000	0.0000
	L 1.2500	-0.6250	0.0000	0.0000	0.5000	-0.5000J

and the solution vector \vec{x} is obtained as

$$\begin{pmatrix} x_1 \\ x_2 \\ x_3 \\ x_4 \\ x_5 \\ x_6 \end{pmatrix} = \begin{pmatrix} 1.0000 \\ 0.9999 \\ 0.9999 \\ 0.9999 \\ 1.0000 \\ 1.0000 \end{pmatrix}$$

Example 4.4

The coefficient matrix [G] is a square/singular matrix, and the RHS vector $\{b\}$ is a

random vector.

<u>۲</u> 2	-1	0	0	0	1	ר1–	(x_1)		ר1]	
4	2	0	0		2	-2	x_2		6	
0		1	2	-5	-3	-1 -2 6 1 0 3 6	x_3		5	
0	0	-2	5	6	2	1	$\{x_4\}$	} =	3	
0	0	3	5	1	7	0	x_5		9	
1	2	-3	6	-2	1	3	x_6		9	
L2	4	-6	12	-4	2	6 J	(x_7)	1	191	l

Using MATLAB built-in function, the rank of G can be computed as

rank(G) = 6

$$G^{T} * G = \begin{bmatrix} 25 & 16 & -15 & 30 & -10 & 15 & 5 \\ 16 & 25 & -30 & 60 & -20 & 13 & 27 \\ -15 & -30 & 59 & -83 & 16 & -1 & -41 \\ 30 & 60 & -83 & 234 & -35 & 69 & 107 \\ -10 & -20 & 16 & -35 & 82 & 24 & -54 \\ -15 & 13 & -1 & 69 & 24 & 72 & -6 \\ 5 & 27 & -41 & 107 & -54 & -6 & 87 \end{bmatrix}$$

Special Cholesky factor of $G^T * G$

	<mark>5.0000</mark> آ	3.2000	-3.0000	6.0000	-2.0000	3.0000	ן 1.0000
	0.0000	3.8418	-5.3099	10.6198	-3.5399	0.8849	-1.0933
_	0.0000	0.0000	4.6695	-1.8438	-1.8838	2.7195	-1.0933
	0.0000	0.0000	0.0000	9.0454	1.2293	5.1535	3.6698
	0.0000	0.0000	0.0000	0.0000	7.7722	4.1069	-4.7144
	L0.0000	0.0000	0.0000	0.0000	0.0000	3.3756	-3.2763

of independent rows = 6

Independent rows = 123456

Dependent row = 7

Product of $L^T * L$

1	94.2400	107.8723	-14.2385	70.9443	-7.9381	ן 6.8505
	107.8723	207.4266	-42.0738	119.0041	-53.0841	-17.3091
_	-14.2385	-42.0738	37.3448	-8.9911	1.6820	12.7625
-	70.9443	119.0041	-8.9911	123.3579	13.4190	5.3727
	-7.9381	-53.0841	1.6820	13.4190	99.5613	29.3094
	6.8505 6 .8505	-17.3091	12.7625	5.3727	29.3094	22.1291 J

Regular Cholesky factorization of $L^T * L$

	[9.7077	11.1120	-1.4667	7.3080	-0.8177	ן 0.7056
	0.0000	9.1624	-2.8131	4.1252	-4.8019	-2.7449
	0.0000	0.0000	5.2229	2.55271	-2.4940	1.1632
_	0.0000	0.0000	0.0000	0.81297	6.6882	1.2578
	0.0000	0.0000	0.0000	0.0000	4.9813	2.2468
	L0.0000	0.0000	0.0000	0.0000	0.0000	2.4723 J

Generalized inverse of [G]

	0.2050	0.1624	0.1051	0.0748	-0.0440	-0.0119	–0.0239 ₁
	-0.4999	0.2500	0.0000	-1.0547	0.0000	0.0000	0.0000
	-0.1361	0.1144	0.1337	0.0393	0.0869	-0.0364	-0.0729
=	0.0579	-0.0468	2.017×10^{-4}	0.0254	0.0272	0.0151	0.0303
	-0.0938	0.0780	0.0410	0.1627	-0.0238	-0.0249	-0.0499
	0.0303	-0.0267	-0.0633	-0.0582	0.0895	8.3858×10^{-3}	0.0167
	L-0.0596	0.0482	0.1468	0.0914	1.3269×10^{-3}	-0.0156	-0.0312 ^J

and the solution vector \vec{x} is obtained from Eq. (4.32) as

1	(x_1)		(1.2092)
	<i>x</i> ₂		1.0000
	<i>x</i> ₃		1.1341
ł	x ₄	} = {	0.5089
	x 5		0.1789
	x 6		0.4105
	(x_7)		0.8290

4.7 Mixed Direct-Iterative Generalized Inverse Algorithms for Solving SLE

Instead of explicitly computing the generalized inverse from Eq. (4.32), which involves a lot of "matrix times matrix" operations, one can/should repeatedly use "matrix times vector" operations for improving its computational efficiency. Furthermore, it is noted that the "regular (not generalized) inverse" of the matrix product [L' * L] should be replaced by the more efficient SLE, either by Direct or by Iterative algorithms. More details can be explained in the following sub-sections.

4.7.1 Direct methods in Generalized Inverse for solving SLE

Using the matrix-product operations in Eq. (4.32), one can compute the unknown solution vector \vec{x} according to the following sequence of steps:

From Eq. (2.1), with $[A] \equiv [G]$, and from Eq. (4.32),

$$\vec{x} = L(L'L)^{-1}(L'L)^{-1}L'G'\vec{b}$$

Let

 $Minv = (L'L)^{-1}$

Then one computes the following sequence of matrix times vector:

$$tempol = G' \times \vec{b}$$

 $tempo3 = L' \times tempo1$

 $tempo2 = Minv \times tempo3$

 $tempol = Minv \times tempo2$

 $tempo2 = L \times tempo1$

and the unknown solution vector \vec{x} is stored in the temporary vector tempo2.

4.7.2 Iterative Method in Generalized Inverse for solving SLE

From Eq. (4.32),

$$\vec{x} = L(L'L)^{-1}(L'L)^{-1}L'G'\vec{b}$$

Let

M1 = (L'L)

Then one computes the following sequence of matrix times vector:

$$tempol = G' \times \overline{b}$$

$$tempo3 = L' \times tempo1$$

Using MATLAB built-in (Conjugate Gradient) function, one computes the following vectors:

$$tempo2 = cg(M1, tempo3)$$

tempol = cg(M1, tempo2)

$$tempo2 = L \times tempo1$$

and the unknown solution vector \vec{x} is stored in the temporary vector tempo2.

where cg is the Conjugate Gradient (iterative) Algorithm for solving SLE and the unknown solution vector \vec{x} is stored in the temporary vector tempo2.

Important Notes:

(a) Inside the "generalized inverse" algorithms, one needs to find the "regular inverse" of the coefficient matrix (= L' * L, in this case). This "regular matrix inversion" should be equivalently solved by SLE, which can be done either by direct, or iterative solvers.

(b) For certain large-scale (especially for 3-D) problems, iterative solver (with appropriated pre-conditioned strategies) can be a more preferable method of choice (as compared to the direct method), due to the following desirable features:

- No fill-in terms occurred
- Much easier to parallelize

4.8 Domain Decomposition Generalized Inverse Solver

The efficient generalized inverse algorithms discussed in sections 4.6 and 4.7 can be further improved by utilizing the Domain Decomposition (or DD) formulation. In Chapter 3, some key equations resulted from DD formulation have already been derived, based on the assumption that the coefficient of a square matrix is non-singular. In this section, let us consider the system of linear algebraic equations

$$Kr = f \tag{4.33}$$

where the matrix of system is $K \in \mathbb{R}^{n \times n}$ (K can be either non-singular, or singular) and vectors $r \in \mathbb{R}^n$, $f \in \mathbb{R}^n$.

Let the original domain be decomposed into m sub domains. The coefficient matrix can be represented in a special form as shown below

$$\begin{bmatrix} K_{1}^{[ii]} & O & K_{1}^{[ib]} \\ K_{2}^{[ii]} & K_{2}^{[ib]} \\ O & K_{3}^{[ii]} & K_{3}^{[ib]} \\ \vdots & \ddots & \vdots \\ K_{m}^{[bi]} & K_{m}^{[ii]} & K_{m}^{[ib]} \\ K_{1}^{[bi]} & K_{2}^{[bi]} & K_{3}^{[bi]} & \cdots & K_{m}^{[bi]} \\ \end{bmatrix} \begin{bmatrix} r_{[i]}^{(1)} \\ r_{[i]}^{(3)} \\ \vdots \\ r_{[i]}^{(m)} \\ r_{[b]} \end{bmatrix} = \begin{bmatrix} f_{[i]}^{(1)} \\ f_{[i]}^{(2)} \\ f_{[i]}^{(3)} \\ \vdots \\ f_{[i]}^{(m)} \\ f_{[b]} \end{bmatrix}$$
(4.34)

From Eq. (4.34) the following notations are used

 $r_{[i]}^{[j]}$, where $j \in \{1, 2, ..., m\}$ - vectors of displacements of interior nodes on the j^{th} sub domain,

 $r_{[b]}$ - vector of boundary displacements,

$$f_{[j]}^{[j]}$$
, where $j \in \{1, 2, ..., m\}$ - vectors of interior loads acting on j^{th} sub domain

 $f_{[b]}$ - vector of boundary loads on the boundaries.

 $K^{[bb]}$ represents the summation of boundary stiffness matrices from all sub-domains.

For "structural engineering" applications, and using DD formulation, Eq. (4.33) can be naturally expressed in the form as shown in Eq. (4.34). For general "field" problems (or mathematical problems), one usually starts with Eq. (4.33). For these problems, however, Nested Disection or METiS re-ordering algorithms [20-22] can be used to transform Eq. (4.33) into the form shown in Eq. (4.34).

From Eq. (4.34), matrix $K^{[bb]}$ contains contributions from "all" sub-domains and can be expressed as

$$K^{[bb]} = \sum_{j=1}^{m} K_{j}^{[bb]}$$
(4.35)

For symmetrical cases, $K_j^{[ib]}$ and $K_j^{[bi]}$ are the transpose of each other. Vectors $r_{[i]}^{[j]}$ can be expressed in the form

$$\eta_{i]}^{[i]} = \left(K_{j}^{[ii]}\right)^{-1} \left(f_{i}^{[j]} - K_{j}^{[ib]} r_{b}^{[j]}\right)$$
(4.36)

where $j \in \{1, 2, ..., m\}$, and $r_b^{[j]}$ is a sub-set of the vector $r_{[b]}$.

Also,

$$\left(K^{[bb]} - \sum_{j=1}^{m} K^{[bi]}_{j} \left(K^{[ii]}_{j}\right)^{-1} K^{[ib]}_{j}\right) r_{[b]} = f_{b} - \sum_{j=1}^{m} K^{[bi]}_{j} \left(K^{[ii]}_{j}\right)^{-1} f^{[j]}_{i}$$
(4.37)

Eq. (4.37) is called the reduced system or resulting system.

If the matrix K, shown in Eq. (4.33), is singular, we can apply the concept of generalized inverse to invert the (possible) singular coefficient matrix in Eq. (4.37) to find the unknown boundary displacement vector $r_{[b]}$ first, and then use them in computing the remaining unknown interior displacement vector $r_{[i]}^{[j]}$, as shown in Eq. (4.36).

For a typical j^{th} sub-domain $j \in \{1, 2, ..., m\}$, one obtains

$$\boldsymbol{\eta}_{[l]}^{(j)} = \left[K_{j}^{(l)} \right]^{-1} \cdot \left(f_{l}^{(j)} - K_{j}^{(lB)} \boldsymbol{r}_{[b]}^{[j]} \right)$$
(4.38)

5. ENGINEERING/SCIENCES NUMERICAL APPLICATIONS

Based on the discussions presented in the previous chapters (chapters 1-4), a fairly extensive set of numerical examples are used in this chapter for validating the accuracy and evaluating the (computational time) performance of the proposed algorithms discussed in Chapter 4. Test examples to cover the cases where the known/given coefficient matrix in the SLE can be a square/rectangular, singular/non-singular, symmetrical/non-symmetrical matrix, and the known/given right-hand-side (rhs) vector can be random vector, or it can be a linear combinations of columns of the coefficient matrix are all investigated in this chapter.

5.1 Description of the Test Examples

Test matrices are collected from Tim Davis Sparse Matrix Collection [14], University of Florida. Rank deficient (singular) matrices derived from various engineering and science applications such as Linear Programming, Combinatorial Problem, Directed Graph, Fluid Dynamics, Linear Programming, Chemical Process Simulations, Cell traffic matrices, etc. are included in Tables 5.1-5.4

SI. No	Name	Size	Rank	nnz	Group	Description
1	lock_700	700x700	165	22,175	HB	Finite Element Problem
2	dwt_1005	1005x1005	995	8,621	HB	Structural Problem
3	bcspwr06	1454x1454	1446	5,300	HB	Power network problem
4	bcsstm13	2003x2003	1241	21,181	HB	Symmetric Mass Matrix, Fluid Flow Generalized Eigen Values
5	lock2232	2232x2232	368	80,352	HB	Finite Element Problem
6	cegb2802	2802x2802	289	277,362	HB	Finite Element Problem

Table 5.1: Symmetric Singular Test Matrices for ODU Generalized Inverse Solver

Sl. No	Name	Size	Rank	nnz	Group	Description
1	tomo_900	900x900	893	35,598	Regtools	2-D tomography test problem
2	GD00_c	638x638	300	1,041	Pajek network	Directed Multigraph
3	GD96_a	1096x1096	827	1,677	Pajek network	Directed Multigraph
4	ex_6	1651x1651	1650	49,062	FIDAP	CFD
5	CS_phd	1882x1882	705	1,740	Pajek network	Directed Graph
6	tomo_2500	2500x2500	2496	166,782	Regtools	2-D tomography test problem

Table 5.2: Un-symmetrical Singular Test Matrices for ODU Generalized Inverse Solver

Sl. No	Name	Size	Rank	nnz	Group	Description
1	D_6	970x435	339	6,491	JGD_SL6	Differentials of
						Voronoi complex of perfect forms
2	mk9-b2	1260x378	343	3,780	JGD_Homology	Combinatorial problem
3	n3c6-b3	1365x455	364	5,460	JGD_Homology	Combinatorial problem
4	Franzl	2240x768	755	5,120	JGD_Franz	Combinatorial problem
5	mk10-b2	3150x630	586	9,450	JGD_Homology	Simplical complexes
6	n4c6-b3	5970x1330	1140	23,880	JGD_Homology	Simplical complexes from Homology from Volkmar Welker

Table 5.3: Rectangular Singular Test Matrices (Tall type: rows>>cols) for ODU Generalized Inverse Solver

Sl. No	Name	Size	Rank	nnz	Group	Description
1	lp_standgub	361x1383	360	3,338	LPnetlib	Linear Programming Problem
2	lp_ship04l	402x2166	360	6,380	LPnetlib	Linear Programming Problem
3	lp_ship08s	778x2467	712	7,194	LPnetlib	Linear Programming Problem
4	Trec12	551x2726	550	151,219	JGD_Kocay	combinatorial problem
5	lp_ship08l	778x4363	712	12,882	LPnetlib	Linear Programming Problem
6	lp_d6cube	415x6184	404	37,704	LPnetlib	Linear Programming Problem

Table 5.4: Rectangular Singular Test Matrices (Fat type: rows<<cols) for ODU Generalized Inverse Solver

5.2 Numerical Performance of ODU Generalized Inverse Solver

Based on the detailed algorithms explained in Chapter 4, and using the rankdeficient matrices as coefficient matrices described in section 5.1, the numerical performance of our proposed procedures [for solving SLE, shown in Eq. (4.33)] are evaluated in this section. The known RHS vector $\{b\}$ can be random vector, or can be chosen such a way that the unknown solution vector $\{x\} = \{1, 1, ..., 1\}$.

We also compared the performance of our algorithm with the efficient algorithm described in [13] and also with MATLAB built-in function *pinv*() [12] for computing the generalized inverse explicitly. We use MATLAB version 7.6.0.324 (R2008a) on Intel Core 2 CPU, 2.13GHZ, 2GB RAM, Windows XP Professional SP3 for numerical comparisons. To be consistent and fair, sparse test matrices obtained from tables 5.1-5.4 are converted to full matrices (in this section).

Tables 5.5 through 5.16 records the times (in seconds) taken by our proposed algorithm, the algorithm mentioned in [13] and MATLAB built-in function [12] pinv(). For our convenience, we represent our algorithm with ODU - ginverse(), algorithm in [13] with geninv and MATLAB built-in function with MATLAB - pinv(). In addition, we have also presented the error norm for all the test matrices.

5.2.1 Direct Method in Generalized Inverse to Solve SLE with RHS Vector as Linear Combination of Columns of the Coefficient Matrix

In this sub-section, the explicitly inverse of the matrix product [L' * L], shown in Eq. (4.32), is implemented by MATLAB built-in function inv(L' * L), and the results are shown in Tables 5.5-5.12.

<i>Sl</i> .	Name	Size	Rank	ODU – ginverse	geninv	MATLAB - pinv()
No				Error Norm	Error Norm	Error Norm
1	lock_700	700x700	165	0.1514 1.033×10 ⁻⁸	0.3446 1.1399×10 ⁻⁶	1.2967 2.215×10 ¹¹
2	dwt_1005	1005x1005	995	2.6634 7.1302×10 ⁻⁹	4.2889 6.764×10 ⁻⁶	14.0320 4.5736×10 ⁻¹²
3	bcspwr06	1454x1454	1446	8.5029 1.477×10 ⁻⁸	13.3176 1.829×10 ⁻⁵	40.3646 2.7131×10 ⁻¹²
4	bcsstm13	2003x2003	1241	11.5997 6.7629×10 ⁻⁹	19.1901 <i>1.826×10⁻⁸</i>	36.3413 5.6493×10 ⁻¹³
5	lock2232	2232x2232	368	5.5518 7.9519×10 ⁻⁹	10.8755 2.5797×10 ⁻⁷	40.7582 1.0761 × 10 ⁻¹¹
6	cegb2802	2802x2802	289	8.9 571 <i>9.7558×10⁻⁹</i>	18.6816 3.7220×10 ⁻⁷	69.9847 1.7532×10 ⁻¹¹

Table 5.5: Computational Times (in seconds) for Symmetric Rank-Deficient Test Matrices with RHS Vector as Linear Combination of Columns of Coefficient Matrix

Sl. No	Name	Size	Rank	ODU – ginverse	geninv	MATLAB - pinv()
				Error Norm	Error Norm	Error Norm
1	tomo_900	900x900	893	1.8545	3.0092	10.1396
				7.667×10 ⁻⁹	1.055×10 ⁻⁶	2.789×10 ¹¹
2	GD00_c	638x638	300	0.1805	0.3961	1.8696
				7.909×10 ⁻¹²	2.602×10 ⁻¹¹	5.704×10 ⁻¹³
3	GD96_a	1096x1096	8 27	2.5991	4.1606	7.3928
				3.179×10 ⁻¹³	1.776×10 ⁻¹³	3.148×10 ⁻¹³
4	ex 6	1651x1651	1650	12.78109	19.9238	44.6059
	_			2.547×10 ⁵	0.00657	4.6022×10 ¹²
5	CS_phd	1882x1882	705	8.5161	12.8672	41.8010
				7.724×10 ⁻¹⁴	9.295×10 ¹⁴	2.707×10 ⁻¹³
6	tomo_2500	2500x2500	2496	44.7203	69.0133	221.6490
	_			2.031×10 ⁷	0.0002893	2.8190×10 ⁻¹⁰

Table 5.6: Computational Times (in seconds) for Non-Symmetric Rank-Deficient Test Matrices with RHS Vector as Linear Combination of Columns of Coefficient Matrix

<i>Sl.</i>	Name	Size	Rank	ODU – ginverse	geninv	MATLAB – pinv()
No				Error Norm	Error Norm	Error Norm
1	D_6	970x435	339	0.1347 1.216× <i>10⁻¹¹</i>	0.2809 1.333 × 10 ⁻¹¹	1.3240 7.403×10 ⁻¹³
2	mk9-b2	1260x378	343	0.1162 5.950×10 ⁻¹⁴	$\begin{array}{c} 0.2478 \\ 1.018 \times 10^{-13} \end{array}$	0.6098 1.681×10 ⁻¹³
3	Franz1	2240x768	755	1.3077 1.457×10 ⁻¹³	2.3649 1.290×10 ⁻¹³	6.0490 2.806×10 ⁻¹³
4	mk10- b2	3150x630	586	0.8094 1.599×10 ⁻¹³	$ \begin{array}{c} 1.5776 \\ 2.057 \times 10^{-13} \end{array} $	3.2363 2.573×10 ⁻¹³

Table 5.7: Computational Times (in seconds) for Rectangular Rank-Deficient Test Matrices (Tall type: Rows>>Cols) with RHS Vector as Linear Combination of Columns of Coefficient Matrix

Sl.	Name	Size	Rank	ODU – ginverse	geninv	MATLAB – pinv()
No				Error Norm	Error Norm	Error Norm
1	lp_standgub	361x1383	360	0.1242 6.321 × <i>10</i> ⁻⁷	0.4238 6.387×10 ⁻⁷	1.0215 3.579×10 ⁻¹¹
2	lp_ship041	402x2166	360	0.1760 1.421×10 ⁻¹¹	0.6712 1.056×10 ⁻¹¹	1.3390 1.851×10 ¹²
3	lp_ship08s	778x2467	712	1.2747 1.178×10 ⁻¹¹	3.4073 1.052×10 ⁻¹¹	5.3489 1.583×10 ⁻¹²
4	lp_ship081	778x4363	712	1.5622 2.955×10 ⁻¹¹	5.2861 1.517×10 ¹¹	8.5278 4.243 × 10 ⁻¹²

Table 5.8: Computational Times (in seconds) for Rectangular Rank-Deficient Test Matrices (Fat type: Rows<Cols) with RHS Vector as Linear Combination of Columns of Coefficient Matrix

Sl.	Name	Size	Rank	ODU – ginverse	geninv	MATLAB - pinv()
No				Error Norm	Error Norm	Error Norm
1	lock_700	700x700	165	0.1483 1495.46	0.3423 1495.46	1.2948 1495.46
2	dwt_1005	1005x1005	995	2.6740 295	4.3156 295	14.0342 295
3	bcspwr06	1454x1454	1446	8.5747 102.551	13.3752 102.551	41.1197 102.551
4	bcsstm13	2003x2003	1241	11.7270 29034.7	19.3159 29034.7	36.3659 29034.7
5	lock2232	2232x2232	368	5.5934 106.489	10.9423 106.489	40.7723 106.489
6	cegb2802	2802x2802	289	9.1066 28002	18.7714 28002	69.7151 28002

5.2.2 Direct Method in Generalized Inverse to Solve SLE with Randomly Generated RHS Vector $\{1, 2, ..., n\}^{T}$.

 Table 5.9: Computational Times (in seconds) for Symmetric Rank-Deficient Test

 Matrices with Randomly Generated RHS Vector

SI.	Name	Size	Rank	ODU – ginverse	geninv	MATLAB - pinv()
No				Error Norm	Error Norm	Error Norm
1	tomo 900	900x900	893	1.8725	3.0287	10.0959
	-			669.871	669 .871	669.871
2	GD00_c	638x638	300	0.2012	0.4092	1.8748
	_			6204.66	6204.66	6204.66
3	GD96_a	1096x1096	827	2.6079	4.1399	7.3863
		4		4821.7	4821.7	4821.7
4	ex_6	1651x1651	1650	12.7676	20.0046	44.5628
	_			1682.04	1682.04	1682.04
5	CS_phd	1882x1882	705	8.5641	12.9633	41.7883
				40498.9	40498.9	40498.9
6	tomo_2500	2500x2500	2496	45.0147	69.2692	221.4820
				1561.25	1561.25	1561.25

 Table 5.10: Computational Times (in seconds) for Non-Symmetric Rank-Deficient Test

 Matrices with Randomly Generated RHS Vector

SI.	Name	Size	Rank	ODU – ginverse	geninv	MATLAB - pinv()
No				Error Norm	Error Norm	Error Norm
1	D_6	970x435	339	0.1991	0.3480	1.3203
	_			14795	14795	14795
2	mk9-b2	1260x378	343	0.1615	0.2935	0.6090
				6376.16	6376.16	6376.16
3	Franzl	2240x768	755	1.2964	2.3602	6.0413
				37127.5	37127.5	37127.5
4	mk10-b2	3150x630	586	0.8063	1.5845	3.2379
				26222.6	26222.6	26222.6

Table 5.11: Computational Times (in seconds) for Rectangular Rank-Deficient Test Matrices (Tall type: Rows>>Cols) with Randomly Generated RHS Vector

Sl.	Name	Size	Rank	ODU – ginverse	geninv	MATLAB – pinv()
No				Error Norm	Error Norm	Error Norm
1	lp_standgub	361x1383	360	0.1586	0.4580	1.0321
				2	2	2
2	lp_ship041	402x2166	360	0.2250	0.7126	1.3383
				1508.7	1508.7	1508.7
3	lp_ship08s	778x2467	712	1.2945	3.3961	5.3447
				833468	833468	833468
5	lp_ship081	778x4363	712	1.5722	5.2922	8.5757
	1 00.000			3724.81	3724.81	3724.81

Table 5.12: Computational Times (in seconds) for Rectangular Rank-Deficient Test Matrices (Fat type: Rows<<Cols) with Randomly Generated RHS Vector

5.2.3 Iterative Methods in Generalized Inverse to Solve SLE with Randomly Generated RHS vector

In this sub-section, the explicit inverse of the matrix product [L' * L], shown in

Eq. (4.32), is implemented by MATLAB built-in (Conjugate Gradient iterative solver),

and the results are shown in Tables 5.13-5.16.

Iterative solver used: Conjugate Gradient Algorithm

Error tolerance used: 10⁻⁷

Sl. No	Name	Size	Rank	ODU — ginverse iterative	ODU – ginverse direct	geninv()	MATLAB pinv(
				Error Norm	Error Norm	Error Norm	Error Norm
1	lock1074	1074x1074	155	0.5649	0.4351	1.0278	4.1070
				47.5347	47.5347	47.5347	47.5347
2	lock2232	2232x2232	368	5.9314	5.5499	10.8729	40.7251
				63.1189	63.1189	63.1189	63.1189
3	cegb2802	2802x2802	289	9.3092	9.0585	18.6983	69.7249
	-			<i>77.0117</i>	77.0117	77.0117	77.0117

Table 5.13: Computational Times (in seconds) for Symmetric Rank-Deficient Test Matrices (Using Iterative Solver inside Generalized Inverse) with Randomly Generated RHS Vector

Name	Size	Rank	ODU – ginverse	ODU – ainverse	geninv()	MATLAB pinv()
			iterative	direct	Error Norm	Error Norm
	ļ					
GD00 c	638x638	300	0.2835	0.2125	0.4097	1.8861
_			46.1318	46.1318	46.1318	46.1318
GD96_a	1096x1096	827	2.7595	2.6291	4.1385	7.3832
			28.0286	28.0286	28.0286	28.0286
CS_phd	1882x1882	705	8.5563	8.5541	12.9385	41.6813
			96.906	96.906	96.906	96.906
	GD00_c GD96_a	GD00_c 638x638 GD96_a 1096x1096	GD00_c 638x638 300 GD96_a 1096x1096 827	Nume Size Nume - ginverse iterative GD00_c 638x638 300 0.2835 GD96_a 1096x1096 827 2.7595 28.0286 CS_phd 1882x1882 705 8.5563	Nume Size Nume - ginverse iterative - ginverse direct GD00_c 638x638 300 0.2835 0.2125 46.1318 46.1318 46.1318 GD96_a 1096x1096 827 2.7595 2.6291 28.0286 28.0286 28.0286 CS_phd 1882x1882 705 8.5563 8.5541	Nume Size Nume Size Nume - ginverse iterative - ginverse direct - ginverse direct - ginverse direct - ginverse direct Error Norm GD00_c 638x638 300 0.2835 0.2125 0.4097 46.1318 46.1318 46.1318 46.1318 46.1318 GD96_a 1096x1096 827 2.7595 2.6291 4.1385 28.0286 28.0286 28.0286 28.0286 28.0286 CS_phd 1882x1882 705 8.5563 8.5541 12.9385

Table 5.14: Computational Times (in seconds) for Non-Symmetric Rank-Deficient Test Matrices (Using Iterative Solver inside Generalized Inverse) with Randomly Generated RHS Vector

Sl. No	Name	Size	Rank	ODU – ginverse	ODU – ginverse direct	geninv()	MATLAB pinv()
			iterative Error Norm	Error Norm	Error Norm	Error Norm	
1	mk9-b2	1260x378	343	0.1423 43.7972	0.1615 <i>43.7972</i>	0.2928 <i>43.7972</i>	0.6096 43.7972
2	Franzl	2240x768	755	1.2200 <i>84.9266</i>	1.3051 <i>84.9266</i>	2.3673 <i>84.9266</i>	6.0349 <i>84.9266</i>
3	n4c6-b3	5970x1330	1140	7.26591 128.086	7.8267 128.086	14.127 8 128.086	25.7394 128.086

Table 5.15: Computational Times (in seconds) for Rectangular (Tall) Rank-DeficientTest Matrices (Using Iterative Solver inside Generalized Inverse) with RandomlyGenerated RHS Vector

SI. No	Name	Size	Rank	ODU — ginverse	ODU — ginverse	geninv()	MATLAB pinv
NU				iterative Error Norm	direct Error Norm	Error Norm	Error Norm
1	lp_ship08s	778x2467	712	1.59103 21.3141	1.2930 21.3141	3.3817 21.3141	5.3262 21.3141
2	lp_ship081	778x4363	712	1.7496 24.7184	1.6065 24.7184	5.2864 24.7184	8.4837 24.7184
3	lp_d6cube	415x6184	404	0.5738 10.7908	0.46281 10.7908	2.1500 10.7908	4.1469 10.7908

 Table 5.16: Computational Times (in seconds) for Rectangular (Fat) Rank-Deficient Test

 Matrices (Using Iterative Solver inside Generalized Inverse) with Randomly Generated

 RHS Vector

Furthermore, graphical comparisons (in terms of the computational times) of ODUginverse with other algorithms have been presented in Appendix E.

5.3 Numerical Performance of ODU Generalized Inverse DD Solver

As can be seen from the previous Section 5.2, our proposed implementation of the generalized inverse for solving SLE have shown "significant time reduction" as compared to MATLAB built-in function, and also compared to [13]. In this sub-section, however, we want to investigate the numerical performance of our proposed generalized inverse within the frame work of DD formulation (for solving system of $[G] * \{x\} = \{b\}$, where the coefficient matrix [G] has the dimension $m \times n$ and can be either rectangular, or square/singular).

5.3.1 Description of Test Problems for ODU Generalized Inverse DD Solver

Test matrices are collected from Tim Davis Sparse Matrix Collection, University of Florida [14]. Rank deficient (singular) matrices derived from various engineering and science applications such as Linear Programming, Combinatorial Problem, Directed Graph, Fluid Dynamics, Linear Programming, Chemical Process Simulations, Cell traffic Matrices, etc. are included in Tables 5.17-5.18

Sl. No	Name	Size	Rank	nnz	Group	Description/Kind
1	GD98_c	112x112	100	336	Pajek	Pajek network, undirected graph
2	dwt_209	209x209	208	1,743	HB	Structural Problem
3	dwt_307	<u>307x307</u>	288	2,523	HB	Structural Problem

Table 5.17: Symmetric Singular Test Matrices for ODU Generalized Inverse DD Solver

Sl. No	Name	Size	Rank	nnz	Group	Description/Kind
1	gent113	113x113	107	655	HB	Statistical/Mathematical Problem
2	gre_216b	216x216	215	812	HB	Directed Weight Graph
3	GD00_a	352x352	178	458	Pajek	Directed Graph

Table 5.18: Un-Symmetrical Singular Test Matrices for ODU Generalized Inverse DD Solver

5.3.2 Problem Formulation with Same Sub Matrix

Let us construct a system matrix (Eq. (4.33)) in the form shown by Eq. (4.34). For the sake of discussion, let us consider 3 sub domains. The system matrix takes the following form

$$\begin{pmatrix} K_{1}^{ii} & O & O & K_{1}^{ib} \\ O & K_{2}^{ii} & O & K_{2}^{ib} \\ O & O & K_{3}^{ii} & K_{3}^{ib} \\ K_{1}^{bi} & K_{2}^{bi} & K_{3}^{bi} & K_{0}^{bb} \end{pmatrix} \begin{pmatrix} r_{i}^{1} \\ r_{i}^{2} \\ r_{i}^{3} \\ r_{b} \end{pmatrix} = \begin{pmatrix} f_{i}^{1} \\ f_{i}^{2} \\ f_{i}^{3} \\ f_{b} \end{pmatrix}$$
(5.1)

A singular coefficient matrix is considered from the test matrix collection [14-15] as K_1^{ii} . For simplicity/convenience, we also assume

(a)
$$K_1'' = K_2'' = K_3'' = K^{bb}$$
 (5.2)

(b)
$$K_1^{ib} = K_2^{ib} = K_3^{ib} = K_1^{bi} = K_2^{bi} = K_3^{bi} = K_1^{ii}$$
 (5.3)

The known right hand side (RHS) vector is chosen as *factor*×*random column vector*. Where *factor* is a user defined variable. The user can also specify the number of sub domains and, accordingly, the system matrix and the right hand side vector will be automatically generated. This section presents a comparison of numerical results (in terms of timings) of the developed Domain Decomposition generalized inverse formulations with other existing algorithms. MATLAB sparse storage scheme has been adopted on the input test matrices.

Sl. No	nsu	Size	Rank	ODU – DD	Original system ginverse MV	Original system geninv
				Error Norm	Error Norm	Error Norm
1	2	336x336	300	0.06130	0.07158	0.11555
				8.21375	8.21375	8.21375
2 3	448x448	400	0.09911	0.15793	0.26940	
				12.2835	12.2835	12.2835
3	4	560x560	500	0.12202	0.30984	0.49831
				9.95631	9.95631	9.95631
4	5	672x672	600	0.15667	0.56595	0.88289
				15.2519	15.2519	15.2519
5	6	784x784	700	0.2000	0.95323	1.45982
				13.1714	13.1714	13.1714
6	7	896x896	800	0.20218	1.44812	2.19521
				14.9267	<i>14.9267</i>	14.9267
7	8	1008x1008	900	0.24897	2.16118	3.21151
		l		14.6983	14.6983	14.6983

Table 5.19: Computational Times (in seconds) for Rank-Deficient Test Matrices with same K^{ii} (GD98_c) Sub-Matrices using Domain Decomposition

Sl. No	nsu	Size	Rank	ODU – DD	Original system ginverse MV	Original system geninv
				Error Norm	Error Norm	Error Norm
1	2	627x627	624	0.28611	0.53233	0.859960
				2.9162	2.9162	2.9162
2	3	836x836	832	0.401712	1.37505	2.10186
				3.87701	3.87701	<i>3.87701</i>
3	4	1045x1045	1040	0.53614	2.98979	4.44309
				2.0488	2.0488	2.0488
4	5	1254x1254	1248	0.59997	5.08382	7.43006
				2.54757	2.54757	2.54757
5	6	1463x1463	1456	0.71609	8.2171	12.06235
				1.06024	1.06024	1.06024
6	7	1672x1672	1664	0.87530	12.46184	18.24484
				2.14487	2.14487	2.14487
7	8	1881x1881	1872	1.00417	18.2003	25.81453
				8.99101	8.99101	8.99101

Table 5.20: Computational Times (in seconds) for Rank-Deficient Test Matrices with same K^{ii} (dwt_209) Sub-Matrices using Domain Decomposition

Sl. No	nsu	Size	Rank	ODU – DD	Original system ginverse MV	Original system geninv
				Error Norm	Error Norm	Error Norm
1	2	921x921	864	0.67732	1.70853	2.59315
				14.6443	14.6443	14.6443
2	3	1228x1228	1152	1.0100	4.38649	6.43760
				12.991	12.991	12.991
3	4	1535x1535	1440	1.2845	8.6928	12.76263
				11.2669	11.2669	11.2669
4	5	1842x1842	1728	1.62886	15.2775	22.0046
				17.1461	17.1461	17.1461
5	6	2149x2149	2016	1.96057	26.66948	35.4461
				17.6758	17.6758	17.6758
6	7	2456x2456	2304	2.28013	36.77451	53.29277
				17.0915	17.0915	17.0915
7	8	2763x2763	2592	2.70171	52.6385	75.09115
				21.1027	21.1027	21.1027

 Table 5.21: Computational Times (in seconds) for Rank-Deficient Test Matrices with same Kⁱⁱ (dwt_307) Sub-Matrices using Domain Decomposition

SI. No	nsu	Size	Rank	ODU – DD	Original system ginverse MV	Original system geninv
				Error Norm	Error Norm	Error Norm
1	2	339x339	321	0.075614	0.093916	0.14541
				4.83871	4.83871	4.83871
2	3	452x452	428	0.010403	0.19469	0.305175
				7.01463	7.01463	7.01463
3	4	565x565	535	0.134215	0.35666	0.575419
				9.03996	9.03996	9.03996
4	5	678x678	642	0.166451	0.65779	1.03111
				10.8548	10.8548	10.8548
5	6	791x791	749	0.20053	1.31634	1.89611
				10.1139	10.1139	10.1139
6	7	904x904	856	0.224461	1.64580	2.46109
				10.1037	10.1037	10.1037
7	8	1017x1017	963	0.30215	2.433860	3.65811
				8.71228	8.71228	8.71228

 Table 5.22: Computational Times (in seconds) for Rank-Deficient Test Matrices with same K^{ii} (gent113) Sub-Matrices using Domain Decomposition

Sl. No	nsu	Size	Rank	ODU – DD	Original system ginverse MV	Original system geninv
				Error Norm	Error Norm	Error Norm
1	2	648x648	643	0.213104	0.48045	0.73973
				29.2826	29.2826	29.2826
2	3	864x864	857	0.32912	1.17688	1.77665
				<i>32.058</i>	32.058	32.058
3	4	1080x1080	1066	0.45151	2.500164	3.57136
				36.0931	36.0931	36.0931
4	5	1296x1296	1279	0.53808	4.33999	6.28897
				40.1448	40.1448	40.1448
5	6	1512x1512	1492	0.64925	7.19551	10.12725
				45.0203	45.0203	45.0203
6	7	1728x1728	1705	0.699806	10.70643	15.20850
				44.2327	44.2327	44.2327
7	8	1944x1944	1918	0.80099	16.44834	23.81176
				54.8367	54.8367	54.8367

Table 5.23: Computational Times (in seconds) for Rank-Deficient Test Matrices with same Kⁱⁱ (gre_216b) Sub-Matrices using Domain Decomposition

Sl. No	nsu	Size	Rank	ODU – DD	Original system ginverse MV	Original system geninv
				Error Norm	Error Norm	Error Norm
1	2	1056x1056	534	0.38620	1.05101	1.62067
				50.4652	50.4652	50.4652
2	3	1408x1408	712	0.56582	2.96705	4.05349
				56.6697	56.6697	56.6697
3	4	1760x1760	890	0.71932	5.83588	8.01557
				63.9613	63.9613	63.9613
4	5	2112x2112	1068	0.888563	10.48973	14.15703
				69.849	69.849	69.849
5	6	2464x2464	1246	1.16060	16.4323	22.84216
				79.2452	79.2452	79.2452
6	7	2816x2816	1424	1.48376	25.37213	34.37000
				82.7129	82.7129	<i>82.7129</i>
7	8	3168x3168	1602	1.69688	36.3547	49.16511
				85.8676	85.8676	85.8676

 Table 5.24: Computational Times (in seconds) for Rank-Deficient Test Matrices with same Kⁱⁱ (GD00_a) Sub-Matrices using Domain Decomposition

5.3.3 Problem Formulation with Different Sub Matrix

Let us construct a system matrix (Eq. (4.33)) in the form shown by Eq. (4.34). For the sake of discussion, let us consider 3 sub domains. The system matrix takes of the form

$$\begin{pmatrix} K_1^{ii} & O & O & K_1^{ib} \\ O & K_2^{ii} & O & K_2^{ib} \\ O & O & K_3^{ii} & K_3^{ib} \\ K_1^{bi} & K_2^{bi} & K_3^{bi} & K^{bb} \end{pmatrix} \begin{pmatrix} r_i^1 \\ r_i^2 \\ r_i^3 \\ r_b \end{pmatrix} = \begin{pmatrix} f_i^1 \\ f_i^2 \\ f_i^3 \\ f_b \end{pmatrix}$$

A singular coefficient matrix is considered from the test matrix collection [14-15] as K_1^{ii} . For simplicity, we also assume

- (a) $K_1'' \neq K_2'' \neq K_3''$
- (b) $K_1^{ii} = K_1^{ib} = K_1^{bi}$
- (c) $K_2^{ii} = K_2^{ib} = K_2^{bi}$
- (d) $K_3^{ii} = K_3^{ib} = K_3^{bi}$
- (e) $K_1^{ii} = K^{bb}$

The known right hand side vector is chosen as $f_i = 1: m, f_b = 1: m$ where $[m, n] = size(K_1^{ii})$. The user can also specify the number of sub domains and accordingly, the system matrix and the right hand side vector will be generated.

Sl. No	nsu	Size	Rank	ODU – DD Error Norm	Original system ginverse MV Error Norm	Original system geninv Error Norm
				12.8452	12.8452	12.8452
2	10	1771x1771	1725	0.65966	14.52799	21.00736
				15.7321	15.7321	15.7321
3	11	1932x1932	1876	0.71032	19.01395	26.94307
				18.9077	18.9077	18.9077

Numerical results for these test cases are presented in Tables 5.25-5.27

Table 5.25: Computational Times (in seconds) for Rank-Deficient Test Matrices with
different Kⁱⁱ (can_161) Sub-Matrices using Domain Decomposition

SI. No	nsu	Size	Rank	ODU DD Error Norm	Original system ginverse MV Error Norm	Original system geninv Error Norm
				2990.92	2990.92	2990.92
2	4	3500x3500	825	2.83254	22.08674	29.95532
				3343.95	3343.95	3343.95
3	5	4200x4200	990	3.88394	37.47156	49.44602
				3663.12	3663.12	3663.12

 Table 5.26: Computational Times (in seconds) for Rank-Deficient Test Matrices with different Kⁱⁱ (lock_700) Sub-Matrices using Domain Decomposition

SI. No	nsu	Size	Rank	ODU – DD	Original system ginverse MV	Original system geninv
				Error Norm	Error Norm	Error Norm
1	4	1445x1445	1438	1.23158	7.89617	11.5578
				2.73861	2.73861	2.73861
2	5	1734x1734	1723	1.33245	14.00956	20.0745
				4.1833	4.1833	4.1833
3	6	2023x2023	2007	1.55488	22.07078	31.8324
				5.91608	5.91608	5.91608

 Table 5.27: Computational Times (in seconds) for Rank-Deficient Test Matrices with different Kⁱⁱ (mesh3e1) Sub-Matrices using Domain Decomposition

5.3.4 Problem Formulation with Same Sub-Matrices and RHS as Linear Combinations of Columns

The known right hand side vector is chosen as $C_1 + C_2 + C_3 + \dots + C_n$, where n is the number of columns, and C_i (with i = 1, 2, ..., n) represents the i^{th} column of the coefficient matrix [G] in the big matrix.

Numerical results for the test cases are presented in Tables 5.28-5.30

Sl.	nsu	Size	Rank	ODU – DD	Original system ginverse MV	Original system geninv
No				$\sum x_i $	$\sum_{i=1}^{n} x_i $	$\sum_{i=1}^{n} x_i $
				Error Norm	Error Norm	Error Norm
1	3	2800x2800	660	2.52323	10.26388	14.11214
				2764	2764	2764
				1.9738×10 ⁻⁶	5.0396×10 ⁻⁷	0.00016343
2	4	3500x3500	825	3.31814	20.38195	27.63699
				3455	3455	3455
				2.2792×10 ⁻⁶	3.0938×10 ⁻⁷	0.0001999
3	5	4200x4200	990	4.04986	35.52977	48.13018
				4146	4146	4146
				2.5482×10 ⁻⁶	7.7514×10 ⁻⁷	0.0002406

Table 5.28: Computational Times (in seconds) for Rank-Deficient Test Matrices with same K^{ii} (lock_700) Sub-Matrices and RHS as Linear Combinations of Columns $C_1 + C_2 + C_3 + \dots + C_n$ using Domain Decomposition

Sl. No	nsu	Size	Rank	$\frac{ODU - DD}{\sum x_i }$	$\begin{array}{c} \textit{Original system} \\ \textit{ginverse MV} \\ \sum x_i \end{array}$	$\frac{Original system}{geninv} \\ \sum x_i $
				Error Norm	Error Norm	Error Norm
1	3	1228x1228	1152	1.05252 1228 3.5175×10 ⁻⁷	4.20519 1228 2.2430×10 ⁻⁸	6.20720 1228 2.4071×10 ⁻⁵
2	4	1535x1535	1440	1.37372 1535 <i>4.0617×10</i> -7	8.58372 1535 5.7408×10 ⁻⁸	12.5739 1535 3.1755×10 ⁻⁵
3	5	1842x1842	1728	1.70161 1842 <i>4.5413×10</i> -7	14.99398 1842 <i>4.0396×10⁸</i>	21.9261 1842 <i>4.3923×10⁻⁵</i>

Table 5.29: Computational Times (in seconds) for Rank-Deficient Test Matrices with same K^{ii} (dwt_307) Sub-Matrices and RHS as Linear Combinations of Columns $C_1 + C_2 + C_3 + \dots + C_n$ using Domain Decomposition

SI. No	nsu	Size	Rank	$\frac{ODU - DD}{\sum_{i=1}^{j} x_{i}}$	Original system ginverse MV $\sum_{ x_i } x_i $ Error Norm	Original system geninv $\sum_{ x_i } x_i $ Error Norm
1	5	2112x2112	1068	1.05579 1164 <i>9.7350×10⁻¹³</i>	9.72518 1164 9.69202 × 10 ⁻¹²	13.58581 1164 <i>3.7717×10¹¹</i>
2	6	2464x2464	1246	1.21400 1358 1.0599×10 ⁻¹²	15.7301 1358 1.0988×10 ⁻¹¹	21.73368 1358 5.7415×10 ¹¹
3	7	2816x2816	1424	1.42343 1552 1.3897×10 ⁻¹²	23.75027 1552 1.3625×10 ¹¹	32.9830 1552 7.8649×10 ¹¹

Table 5.30: Computational Times (in seconds) for Rank-Deficient Test Matrices with same K^{ii} (GD00_a) Sub-Matrices and RHS as Linear Combinations of Columns $C_1 + C_2 + C_3 + \dots + C_n$ using Domain Decomposition

5.3.5 Problem Formulation with Different Sub-Matrices and RHS as Linear

Combinations of Columns

The known right hand side vector is chosen as $C_1 + C_2 + C_3 + \dots + C_n$ where n is the number of columns and C_i (with i = 1, 2, ..., n) represents the i^{th} column of the coefficient matrix [G] in the big matrix.

Numerical results for the test cases are presented in Tables 5.31-5.32

Sl. No	nsu	Size	Rank	$\frac{ODU - DD}{\sum \mathbf{x}_i }$	$\frac{Original system}{ginverse MV}$	$\begin{array}{c} \textit{Original system} \\ \textit{geninv} \\ \sum_{i} x_i \end{array}$
				Error Norm	Error Norm	Error Norm
1	3	1228x1228	1152	1.04227	4.381476	6.44660
				1228	1228	1228
				3.9554×10 ⁻⁷	2.7835×10 ⁻⁸	2.4688×10 ⁵
2	4	1535x1535	1439	1.39427	8.91303	12.98718
				1535	1535	1535
				4.2179×10 ⁷	5.0078×10 ⁻⁸	3.237×10 ⁻⁵
3	5	1842x1842	1725	1.64756	15.51327	22.31969
				1842	1842	1842
				4.42625×10 ⁻⁷	5.60808×10 ⁻⁸	4.6407×10 ⁻⁵

Table 5.31: Computational Times (in seconds) for Rank-Deficient Test Matrices with different K^{ii} (dwt_307) Sub-Matrices and RHS as Linear Combinations of Columns $C_1 + C_2 + C_3 + \dots + C_n$ using Domain Decomposition

Sl. No	nsu	Size	Rank	$\frac{ODU - DD}{\sum \mathbf{x}_t }$	Original system ginverse MV $\sum x_t $	Original system $ \begin{array}{c} geninv\\ \sum x_i \end{array} $
				Error Norm	Error Norm	Error Norm
1	15	1792x1792	1590	0.47628 1791.76 6.7768×10 ⁻¹²	13.16311 1791.76 2.0569×10 ¹⁰	19.00465 1791.76 <i>1.4315×10⁻⁹</i>
2	16	1904x1904	1686	0.52987 1903.66 7.9443×10 ⁻¹²	15.83865 1903.66 2.9199×10 ¹⁰	22.6178 1903.66 7.944×10 ⁻¹²
3	17	2016x2016	1781	0.57382 2015.52 7.2594×10 ⁻¹²	18.88350 2015.52 <i>3.3548×10¹⁰</i>	26.91830 2015.52 2.16341×10 ⁹

Table 5.32: Computational Times (in seconds) for Rank-Deficient Test Matrices with different K^{ii} (GD98_c) Sub-Matrices and RHS as Linear Combinations of Columns $C_1 + C_2 + C_3 + \dots + C_n$ using Domain Decomposition

6. EDUCATIONAL GENERALIZED INVERSE SOFTWARE FOR INTERNET USERS

Most of the currently available commercialized software (and/or freely available public source codes) for "large-scale" Engineering/Science computation has been written in either FORTRAN, C, or C++ languages. Legacy (commercialized) Finite Element Analysis (FEA) software, such as MSC-NASTRAN, SAP-2000, etc. have been written in FORTRAN language. While these languages are efficient for "number crunching", these FORTRAN/C/C++ software are NOT suitable for internet (educational) users. On the other hand, it is too time consuming if one has to re-write these (large) source codes in JAVA, or FLASH, etc... for internet/educational purposes. Based on our earlier research works [23], a general procedure for executing "any" FORTRAN software on the internet is explained and summarized in this chapter. More specifically, this chapter will explain how to use the developed FORTRAN generalized software.

6.1 Description for Executing FORTRAN Software on the Internet

Since the 1960's, scientific programs have been developed in FORTRAN for the solution of various structural, environmental, mathematical, chemical, etc. problems.

With the growing popularity and possibilities of the internet, web-based learning is becoming more and more popular these days. The new trend focuses on developing more effective learning methods for large pre-existing scientific languages like FORTRAN, C etc. In this chapter, a web-based environment is utilized as a means to introduce numerical methods concepts in civil engineering and other related fields of engineering. Software development and implementation is presented, including detailed descriptions of the techniques employed to link software written in high level computer languages, such as C and FORTRAN, to a web-based, user friendly interface for both input and output.

Web-based instruction systems represent a developing branch of computer-aided instruction. This type of educational information emphasizes the use of the web for transfer of educational information, and may be considered as a replacement to traditional delivery methods of lectures and textbooks.

The motivation for developing this educational software is the challenges faced while developing a 3-D truss analysis module in FLASH Actionscript language [24]. Initially, a web-based module was developed by converting 3-D truss analysis FORTRAN code to FLASH Actionscript language. This module analyzes a 3-D truss with the user specified input data. This module needs to re-write the entire FORTRAN code to a different language (in this case FLASH Actionscript). This process is not only time consuming; it requires a good knowledge of the other programming language (FLASH Actionscript). Fig. 6.1 shows a sample of the developed web based 3D truss module and can be found at http://www.lions.odu.edu/~skadi002/3d/. This emerged as one of the many situations where developed FORTRAN computer programs are no longer available for the public use (easy use) and are still considered to be valuable source of both research and educational material. A better and convenient way to interact the FORTRAN programs directly from the server is in needed to serve the purpose.

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Figure 6.1: Online 3-D Truss Analysis.

6.2 Client Server Interface

Web application software is an application that uses a web browser as a client. Commonly used web browsers are Internet explorer, Mozilla Firefox, Google Chrome, Apple Safari, etc. A client is a system that accesses a remote service on another computer commonly known as a server.

Web applications commonly use a combination of server-side scripting languages like PHP, ASP, etc., and client-side languages like HTML, PHP, JAVA, etc. In the developed software, PHP [25] is used as both client side and server side technologies. In Fig. 6.2 client can be any one of desktop, laptop, Mac PC or PDA with browser capabilities and an internet connection. The UNIX server is a machine where the compiled FORTRAN programs are stored.

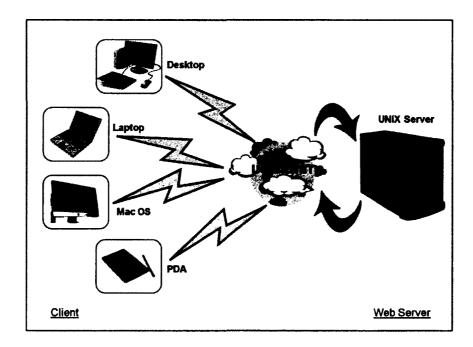


Figure 6.2: Client-Server Interface

6.3 Detailed Step-by-Step Procedures

Input and output interfaces are made user-friendly to attract users' attention, and guide them into use of the software in an efficient manner.

(a) The user references http://www.lions.odu.edu/~skadi002/work/ as URL which directs to a webpage containing number of educational applications. The user has a choice to select from various FORTRAN applications.

(b) For the sake of discussion, let us assume the user has interest in solving SLE (singular coefficient matrix) using generalized inverse. After clicking the appropriate link, the browser directs to a new webpage containing the description and usage of the application (see Fig. 6.3).

(c) The user clicks the "New" button to enter the input data in the textbox provided. It is a hypertext preprocessor (PHP) form that is interpreted by the browser to allow the user to enter the input data (see Fig 6.4).

(d) A sample/demo input data is provided in the input page so that the user can prepare in the specified fashion. Well documented instructions are provided for the users to prepare the input data to the application.

(e) Once the input data has been entered in the text box, the user hits the "Click here to submit input file" button which is located below the input textbox.

(f) At this point the input data has been sent to the FORTRAN program. The user clicks the "EXECUTE" button to execute/run the application.

(g) The output of the application is instantly seen on the same webpage (see Fig. 6.5)

6.4 Demonstrated Examples

In this section, various examples including the one explained in section 6.3 have been demonstrated. Below are the descriptions of the applications.

6.4.1 Solving system of Simultaneous Linear Equations using Generalized Inverse Algorithms

This application solves SLE of singular coefficient matrix using Generalized inverse algorithms. Fig. 6.3 shows a screenshot of the application homepage. The input data page with the sample of how input data page is prepared/entered is shown in Fig. 6.4. The output of the application is shown in Fig. 6.5.

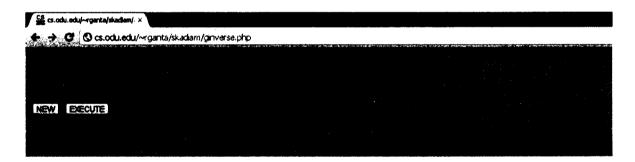


Figure 6.3 Sample of Generalized Inverse Home Page

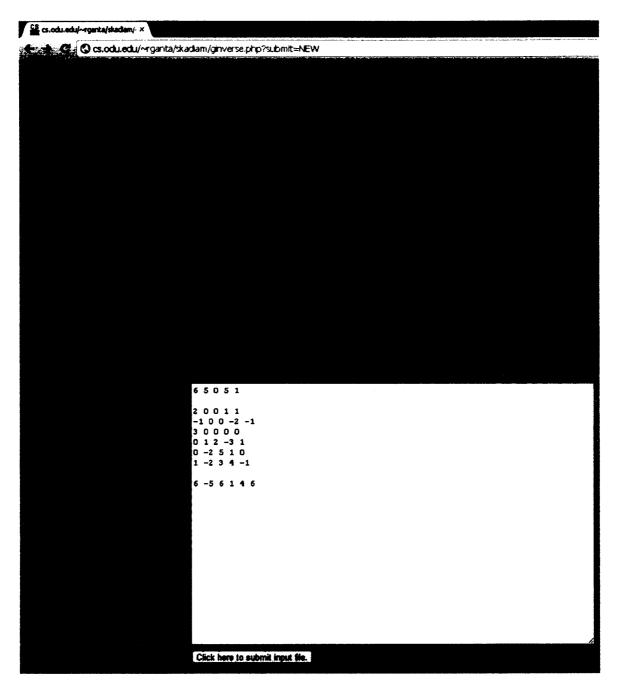


Figure 6.4 Generalized Inverse Input Page

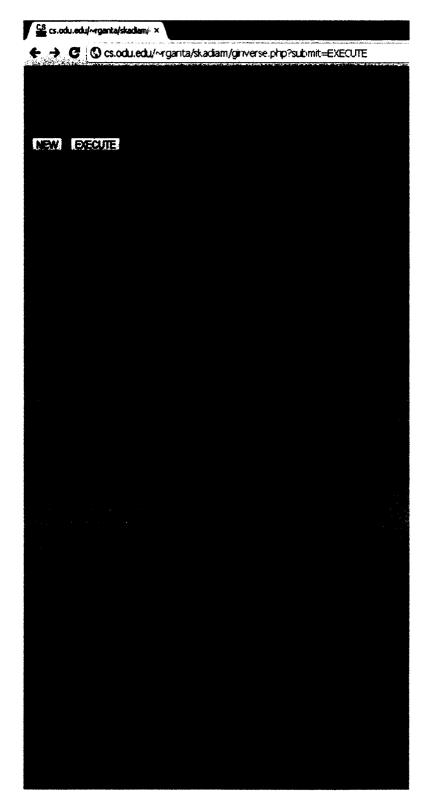


Figure 6.5 Generalized Inverse Output Page

6.4.2 Solving System of Simultaneous Linear Equations using LU Decomposition Algorithms

This application solves SLE of non-singular coefficient matrix using LU decomposition algorithms. Fig. 6.6 shows a screenshot of the application homepage. The input data page with the sample of how input data page is prepared is shown in Fig. 6.7. The output of the application is shown in Fig. 6.8.



Figure 6.6: Sample of LU Decomposition Home Page.

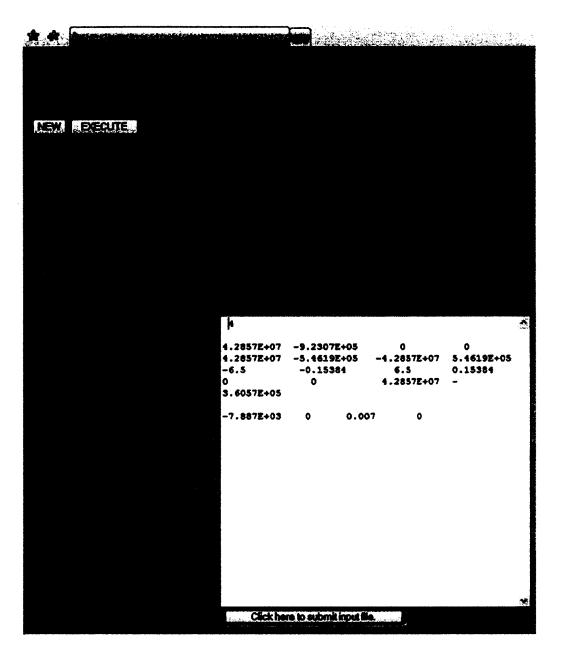


Figure 6.7: LU Decomposition Input Page

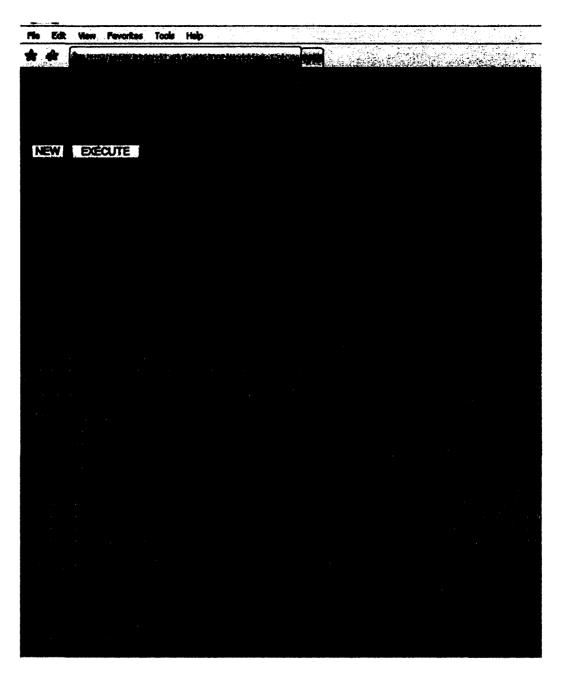


Figure 6.8: LU Decomposition Output Page

6.4.3 Solving System of Simultaneous Linear Equations using LU Decompositions Algorithms on Portable Devices

Nowadays, most of the portable devices such as smart phones, tablet PCs, personal digital assistants (PDA), e-book readers, etc., have internet browsing capabilities. The developed application can be accessed on the above said portable devices. One such example is shown in Fig. 6.9. In this example, the developed educational software is accessed on I-Phone.

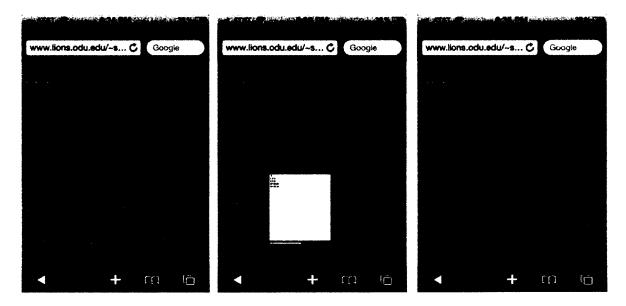


Figure 6.9: I-Phone Homepage, Input Data and Output Data.

7. MATLAB – MPI BUILT-IN FUNCTIONS FOR PARALLEL COMPUTING APPLICATIONS

7.1 Introduction

Matlab (MATrix LABoratory) is a tool to do numerical computations, solve engineering and sciences applications, display 2D and 3D graphical information, algorithm development, simulation, etc. It is a high-level scientific and engineering programming environment which provides many useful capabilities and has an extensive library of built-in functions.

Message Passing Interface (MPI) is widely used in large scale (intensive numerical) computations. This is especially true for "generalized inverse" computer implementation, where as matrix times matrix and/or Cholesky factorization operations were required. MPI is a library of functions/routines that can be used to create parallel programs in scientific languages such as FORTRAN, C, C++, etc.

Similar to traditional Message Passing Interface (MPI), MatlabMPI developed in Lincoln Laboratory, MIT, allows any Matlab programs to run in parallel. MatlabMPI implements the widely used MPI "look and feel" on the top of standard Matlab file input/output, resulting in Matlab implementation of MPI.

MatlabMPI can be downloaded into user's local ZORKA account from http://www.ll.mit.edu/mission/isr/matlabmpi/matlabmpi.html

7.2 MatlabMPI Functions

Some basic, most often used MatlabMPI built-in functions are briefly discussed below:

a. function MPI_Init

This function is called at the start of an MPI program. It also initializes MPI in Matlab environment.

Example:

MPI_Init;

b. function NP = MPI_Comm_size(comm)

This function returns the numbers of processors in the communicator. The "comm" in this function is an MPI communicator which is typically a copy of MPI_COMM_WORLD.

Example:

NP = MPI_Comm_size(comm)

c. function my_ID = MPI_Comm_rank(comm)

This function returns the rank (or processor ID #) of the current processor.

Example:

٠

my_ID = MPI_Comm_rank(comm)

d. function MPI_Send(dest,tag,comm,varargin)

This function sends variable to a destination. It sends message containing variables to a specified destination with a given tag. The argument "dest" contains processor ID #, "tag" can be an any integer and "varargin" represents variable argument inputs.

Example:

MPI_Send (dest,tag,comm,data1,data2,data3,..)

e. function varargout = MPI_Recv(source,tag,comm)

This function receives a message from a specified source processor with a given tag and returns the output variable(s).

Example:

[var1,var2,var3,...] = MPI_Recv(source,tag,comm)

f. function MPI_Abort()

This function will abort any currently running MatlabMPI sessions by looking for leftover Matlab jobs and killing them.

g. function MPI_Finalize()

This function is the last statement indicating the end of a MatlabMPI program.

h. function MPI_Run(m_file,n_proc,machines)

This function runs a Matlab file by name "m_file" on multiple processors. It also runs "n_proc" number of copies of m_file on machines. To run on multiple processors, the argument "machines" are to be designated with "machine1, machine2,...'.

Example:

MPI_Run('example1',2,{}); for the case a single node (with 2 processors) is used.

MPI_Run('example2',4,{zorka1, zorka2}); for the case multiple nodes (assuming zorka1 and zorka2 are both available) are used.

The above discussed functions are used within traditional Matlab source code to run in parallel environment. In addition to the MPI functions, Matlab uses other built-in functions to perform various operations/tasks. Some of the additional functions are discussed below:

i. function eval()

This function execute string with Matlab expression and is also used with MPI_Run. Example:

eval(MPI_Run('example1',2,{}))

j. function disp()

This function displays an array without printing the array name. It can also be used to display a string or a text inside the Matlab code.

Example:

disp(['Example1 from rank: ',num2str(my_rank)]);

If the rank is 0 (master processor), the output appears on the screen. And if the rank is more than 0, i.e 1,2,3..., the output prints to the corresponding file.

k. function MPI_Bcast(source, tag, comm, varargin)

This function broadcasts the variable(s) to all processors.

Example:

[var1, var2, ..] = MPI_Bcast(source,tag,comm,data1,data2,..)

7.3 Example 1: Display Rank of Processors

In this example, a simple MatlabMPI source code to print/display rank of different processors is shown below.

% MPI INITIALIZE

MPI_Init;

% MPI COMMUNICATOR

comm = MPI_COMM_WORLD;

% GET SIZE OR NUMBER OF PROCESSORS IN THE COMMUNICATOR

NP = MPI_Comm_size(comm)

% GET RANK (or ID #) OF CURRENT PROCESSOR

my_rank = MPI_Comm_rank(comm)

% DISPLAY RANK OF EACH PROCESSOR

disp(['Hello Message from rank: ', num2str(my_rank)]);

% FINALIZE Matlab MPI

MPI_Finalize;

% DISPLAY SUCCESS MESSAGE

disp('Success');

Let the file name for this MatlabMPI application be example1.m. In order to run in parallel environment, type the following statements in Matlab command prompt.

% ADDING PATH TO THE MatlabMPI SOURCE DIRECTORY TO INVOKE MPI FUNCTIONS.

addpath /local/MatlabMPI/src

% example1.m IS A MatlabMPI APPLICATION CODE WHICH NEEDS TO EXIST IN THE SAME WORKING DIRECTORY AS THE OTHER MPI FUNCTIONS. IN THIS CASE WE ARE USING 4 PROCESSORS.

eval(MPI_Run('example1',4,{}));

% ONCE THE DESIRED OUTPUT IS OBTAINED/PRINTED, THE FUNCTION MatMPI_Delete_all HAS TO BE INVOKED. THIS FUNCTION DELETES LEFTOVER MatlabMPI FILES FROM THE PREVIOUS RUN. THIS FUNCTION IS ALSO INVOKED BEFORE THE START OF A NEW MatlabMPI APPLICATION.

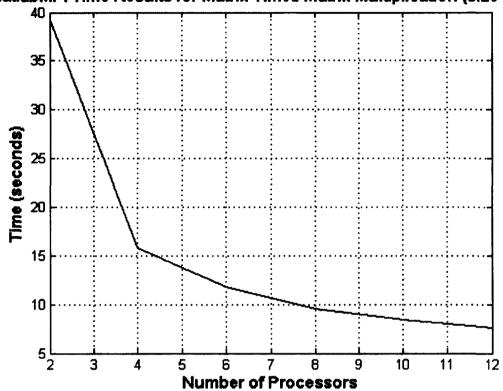
7.4 Example 2: Matrix-Matrix Multiplication

The MatlabMPI source code for matrix-matrix multiplication (dense format) can be found in Appendix D.

Below are the time results for matrix times matrix multiplication (size = 1000).

Sl.No	Number of Processors (NP)	Time (seconds)		
1	2	39.4087		
2	4	15.7884		
3	6	11.7958		
4	8	9.5475		
5	10	8.3805		
6	12	7.5807		

Table 7.1 Time Results (in seconds) for Matrix-Matrix Multiplication using MatlabMPI



MatlabMPI Time Results for Matrix Times Matrix Multiplication (size=1000)

Figure 7.1 Graphical Representation of Time Results (in seconds) for Matrix-Matrix Multiplication using MatlabMPI

8. CONCLUSIONS AND FUTURE WORKS

In this dissertation, various efficient algorithms for solving SLE with full rank or rank deficient have been reviewed, proposed and tested. These algorithms were based on efficient generalized inverse algorithms, which had also been incorporated into the DD formulation. Users are provided the options of incorporating either direct, or iterative solvers into the developed DD generalized inverse formulation. Extensive numerical results have been used to evaluate the performance (in terms of numerical accuracy, calculated error norm, CPU/wall-clock time) of the proposed procedures. The developed numerical procedures can be applied to solve "general" SLE (in the form $[G]{x} = {b}$, coefficient [G] where the matrix could square/rectangular, be symmetrical/unsymmetrical, non-singular/singular). Numerical results have shown that the proposed algorithms are highly efficient as compared to existing algorithms [6, 9, 13] (including the popular MATLAB built-in function pinv(G) * b) [12]. Further reduction in wall-time can also be realized/achieved by taking advantages of "parallel matrix times matrix operations" under MATLAB-MPI computer environment [26].

Furthermore, this dissertation has also contributed the "educational value" to the educational communities, by providing the tools/technologies to execute any existing FORTRAN code for internet users, without requiring them to download any (commercial) software on their desktop/laptop computers. The only requirement for the users to use/learn/execute our FORTRAN-web application is to have access to the internet, which is readily available not only in every home, but also in most public places (such as in the airports, hotels, universities, restaurants, etc.).

Extensions to this current work may include a variation of DD formulation proposed in [4], parallel implementation of the proposed DD generalized inverse solver, and incorporating METiS [20] reordering algorithm for automatically partitioned a given coefficient matrix into diagonal blocks, in such a way to minimize the total number of boundary (interface) nodes, etc.

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APPENDIX A

SINGULAR VALUE DECOMPOSITION (SVD) AND THE GENERALIZED INVERSE

Example A1

Given
$$A = \begin{bmatrix} 2 & 3 \\ 4 & 6 \end{bmatrix}$$
 (A.1)

Step1: Compute $AA^{H} = AA^{T} = \begin{bmatrix} 2 & 3 \\ 4 & 6 \end{bmatrix} \begin{bmatrix} 2 & 4 \\ 3 & 6 \end{bmatrix} = \begin{bmatrix} 13 & 26 \\ 26 & 52 \end{bmatrix}$ (A.2) Also: $AA^{H} = (U\Sigma V^{H})(V\Sigma^{H}U^{H}) = U\Sigma^{2}U^{H}$ Similarly, $A^{H}A = V\Sigma^{2}V^{H}$

Also compute
$$A^{H}A = A^{T}A = \begin{bmatrix} 2 & 4 \\ 3 & 6 \end{bmatrix} \begin{bmatrix} 2 & 3 \\ 4 & 6 \end{bmatrix} = \begin{bmatrix} 20 & 30 \\ 30 & 45 \end{bmatrix}$$
 (A.3)

Step2: Compute the standard Eigen-solution individually for $AA^{T}and A^{T}A$

Using MATLAB built-in function "eig", Eigen-values and the corresponding Eigenvectors for AA^{T} are given as $[u1, lamba1] = eig(AA^{T})$

$$u1 = \begin{bmatrix} -0.8944 & 0.4472\\ 0.4472 & 0.8944 \end{bmatrix} \equiv U \tag{A.4}$$

$$lambda1 = \begin{bmatrix} 0 & 0\\ 0 & 65 \end{bmatrix}$$
(A.5)

Also, for $A^T A$

 $[u2, lamba2] = eig(A^{T}A)$ $u2 = \begin{bmatrix} -0.8321 & 0.5547\\ 0.5547 & 0.8321 \end{bmatrix} \equiv V$ (A.6)

$$lambda2 = \begin{bmatrix} 0 & 0\\ 0 & 65 \end{bmatrix}$$
(A.7)

From the above equations, we can observe that Eigen-values in both the cases are the same, but their corresponding Eigen-vectors are different.

Also,
$$\sigma = \sqrt{lambda1 = lambda2}$$

Computing $\sigma = \sqrt{lambda1 = lambda2} = \begin{bmatrix} 0 & 0 \\ 0 & \sqrt{65} \end{bmatrix}$ (A.8)

From Eqs. (A.4, A.7 and A.8), the SVD of Eq. (A.1) can be obtained as

$$A = U\Sigma V = \begin{bmatrix} -0.8944 & 0.4472\\ 0.4472 & 0.8944 \end{bmatrix} \begin{bmatrix} 0 & 0\\ 0 & \sqrt{65} \end{bmatrix} \begin{bmatrix} -0.8321 & 0.5547\\ 0.5547 & 0.8321 \end{bmatrix}$$
(A.9)

The generalized inverse A^+ of Eq. (A.1) is computed as

$$A^+ = V \Sigma^+ U^H \tag{A.10}$$

where
$$\Sigma^+ = \begin{bmatrix} 0 & 0\\ 0 & \frac{1}{\sqrt{65}} \end{bmatrix}$$
 (A.11)

Hence,

$$A^{+} = \begin{bmatrix} -0.8321 & 0.5547 \\ 0.5547 & 0.8321 \end{bmatrix} \begin{bmatrix} 0 & 0 \\ 0 & \frac{1}{\sqrt{65}} \end{bmatrix} \begin{bmatrix} -0.8944 & 0.4472 \\ 0.4472 & 0.8944 \end{bmatrix}$$
(A.12)

$$A^{+} = \begin{bmatrix} 0.0308 & 0.0615\\ 0.0462 & 0.0923 \end{bmatrix}$$
(A.13)

The result obtained in Eq. (A.13) has been checked with MATLAB generalized inverse function pinv() and same result is obtained.

APPENDIX B

AN EDUCATIONAL FORTRAN SOURCE CODE OF "SPECIAL LDL"" ALGORITHM FOR FACTORIZATION OF SINGULAR/SQUARE/SYMMETRICAL COEFFICIENT MATRIX

```
С
       Implicit real*8 (a-h, o-z)
С
C=
С
       Remarks :
С
              Identifying which are dependent rows of a "floating" substructure
       (a)
С
              Factorizing (by LDL transpose) of a floating substructure stiffness
       (b)
С
              Whenever a dependent row is encountered during LDL factored
С
              process, then we just :
С
              [1] set all factorized values of the dependent row to be ZEROES
С
              [2] ignore the dependent row(s) in all future faztorized rows
С
              [K "float"] = [K11]
                                    [K12]
       (c)
С
                            [K21] [K22]
С
              where [K11] = full rank (=non-singular)
С
       (d)
              The LDL transpose of [K11] can be obtained by taking the results
С
              of part (b) and deleting the dependent rows/columns
С
       Author(s) : Prof. Duc T. Nguyen
С
       Version : 04-30-2004 (EDUCATIONAL purpose, LDL/FULL matrix is
С
assumed)
       Stored at : cd ~/cee/*odu*clas*/generalized inverse by ldl.f
С
С
C=
С
       dimension u(99,99), idepenrows(99), tempo1(99)
С
       iexample=1
                             ! can be 1, or 2, or 3
С
       if (iexample . eq. 1)
                             n=3
С
if (iexample . eq. 2) n=12
С
       if (iexample . eq. 3)
                             n=7
С
       do 1 i=1.n
       do 2 i = 1, n
       u(i,j)=0
2
       continue
       continue
 1
```

if (iexample . eq. 1) then С u(1,1)=2.! non-singular case u(1,1)=1.! С singular case u(1,2) = -1. u(2,2)= 2. u(2,3) = -1.u(3,3)=1.С elseif (iexample . eq. 2) then С u(1,1)= 1.88*10**5 $u(1,2) = -4.91 \times 10^{**4}$ u(1,3) = -1.389*10**5 $u(1,7) = -4.91 \times 10^{*}4$ $u(1,8) = 4.91 \times 10^{**4}$ С u(2,2)=1.88*10**5 u(2,6) = -1.389 * 10 * * 5 $u(2,7) = 4.91 \times 10^{10}$ $u(2,8) = -4.91 \times 10^{**4}$ ¢ u(3,3)=1.88*10**5 $u(3,4) = 4.91 \times 10^{**4}$ $u(3,5) = -4.91 \times 10^{*}4$ $u(3,6) = -4.91 \times 10^{*}4$ С $u(4,4) = 1.88 \times 10^{10}$ $u(4,5) = -4.91 \times 10^{10}$ $u(4,6) = -4.91 \times 10^{*}4$ u(4,8) = -1.389 * 10 * * 5С u(5,5)=2.371*10**5u(5,7) = -1.389 * 10 * * 5u(5,11) = -4.91 + 10 + 4 $u(5,12) = 4.91 \times 10^{*}4$ С $u(6,6) = 3.76 \times 10^{10}$ u(6,10) = -1.389*10**5 $u(6,11) = 4.91 \times 10^{**4}$ $u(6,12) = -4.91 \times 10^{**4}$ С u(7,7)=2.371*10**5 $u(7,9) = -4.91 \times 10^{10}$ u(7,10)= -4.91*10**4

С

85

C u(8,8)= 3.76*10**5 $u(8,9) = -4.91 \times 10^{**4}$ u(8,10)= -4.91*10**4 u(8,12) = -1.389*10**5С u(9,9)= 1.88*10**5 u(9,10) = 4.91 * 10 * * 4u(9,11)= -1.389*10**5 С u(10,10) = 1.88*10**5С u(11,11)= 1.88*10**5 $u(11,12) = -4.91 \times 10^{*}4$ С $u(12,12) = 1.88 \times 10^{**5}$ С elseif (iexample . eq. 3) then С u(1,1)=1. u(1,2)=2.u(1,3) = -3.u(1,4)=2.u(1,5)= -2. u(1,6) = -3.u(1,7) = -2.С u(2,2)= 4. u(2,3) = -6.u(2,4)=4.u(2,5) = -4.u(2,6) = -6.u(2,7) = -4.С u(3,3)=9.u(3,4) = -6.u(3,5)=6.u(3,6)= 9. u(3,7)=6.С u(4,4)=5.u(4,5) = -1.u(4,6)= -5. u(4,7) = -7.С u(5,5)=13.

u(5,6)=9.u(5,7) = -5.С u(6,6)=13.u(6,7)=9.u(7,7)=27. Endif С do 4 i=1,ndo 5 j=1,nu(j,i)=u(i,j)5 continue continue 4 С call generalized_inverse_ldl (n, u, idependrows, ndependrows) С write(6,*) '# dependent rows = ', ndependrows if (ndependrows .ge. 1) then write(6,*) ' dependent rows = ',(idependrows(i),i=1, ndependrows) endif write(6,*) 'LDL factorized u(-, -) =', ((u(i,j), j=i,n), i=1,n) С extracting & writing the LDL factorized of full rank of [K11] С by deleting the dependent row(s) /column(s) of [u] С do 52 i=1,n iskiprow=0 do 53 j=1,ndependrows if (idependrows(j) .eq. i) iskiprow=1 53 continue if (iskiprow.eq. 1) go to 52 icount=0 do 54 j=i,n iskipcol=0 do 55 k=1,ndependrows if (idepenrows(k) .eq. 0) iskipcol=1 55 continue if (iskipcol .eq. 0) then icount=icount+1 tempol(icount)=u(i,j)endif 54 continue write(6,*) 'LDL of [K11] = ',(tempo1(k),k=1,icount) 52 continue С stop end

```
С
С
      subroutine generalized_inverse_ldl (n, u, idependrows, ndependrows)
      Implicit real*8 (a-h, o-z)
      dimension u(99,*), idepenrows(*)
С
C
С
      Remarks:
С
      (a)
             Identifying which are dependent rows of a "floating" substructure
С
      (b)
             Factorizing (by LDL transpose) of a floating substructure stiffness
С
              Whenever a dependent row is encountered during LDL factored
              process, then we just :
С
              [1] set all factorized values of the dependent row to be ZEROES
C
              [2] ignore the dependent row(s) in all future faztorized rows
С
      (c)
              [K "float"] = [K11]
                                  [K12]
С
                          [K21] [K22]
С
             where [K11] = full rank (=non-singular)
С
             The LDL transpose of [K11] can be obtained by taking the results
      (d)
С
              of part (b) and deleting the dependent rows/columns
      Author(s) : Prof. Duc T. Nguyen
С
      Version: 04-30-2004
С
      Stored at : cd ~/cee/*odu*clas*/generalized inverse by ldl.f
С
С
C⁼
С
 eps=0.000000001
      do 11 i=2.n
        do 22 k=1,i-1
        if (dabs( u(k,k) ) .lt. eps)
                                                ! check for "previous"
                                  go to 22
                                            ! dependent row(s)
С
        xmult=u(k,i)/u(k,k)
          do 33 j=i,n
          u(i,j)=u(i,j)-xmult^{*}u(k,j)
33
          continue
        u(k,i)=xmult
22
        continue
С
C=
С
      to zero out entire dependent row
С
        if (dabs( u(i,i) ).lt. eps)
                                 then
        write(6,*) 'dependent row # i, u(i,i) = ', i, u(i,i)
        ndependrows= ndependrows+1
        idependrows(ndependrows)= i
```

do 42 i=i.n42 u(i,j)=0.do 44 k=1,i-144 u(k,i)=0.endif С C≃ С continue 11 С return end С

L D L_t factorized of the "full rank" sub-matrix [K11] of Example 3

```
dependent row # i, u(i,i) =
                        2
                            0.0E+0
dependent row # i, u(i,i) =
                        3
                            0.0E+0
dependent row \# i, u(i,i) =
                        5
                            0.0E+0
\# dependent rows = 3
dependent rows =
                        5
                 2
                    3
LDL of [K11] = 1.0 \ 2.0 \ -3.0 \ -2.0
LDL of [K11] = 1.0 \ 1.0 \ -3.0
LDL of [K11] = 3.0 2.0
LDL of [K11] = 2.0
┿╊┼╋┼╊╉╄╋╋╪┼╋╋┼┼╋╋┿┿╋┿┼╋┽┼╋┿┿┿┿┿╋┿┼
L D L t factorized of the "full rank" sub-matrix [K11] of Example 2
```

```
dependent row # i, u(i,i) =
                         10
                             -8.731149137020111E-11
dependent row \# i, u(i,i) =
                         11
                             -5.820766091346741E-11
dependent row \# i, u(i,i) =
                         12 -2.9103830456733703E-11
# dependent rows =
                   3
dependent rows = 10 11
                           12
LDL of [K11] = 188000.0 -0.26117002127659574 -0.7388297872340426 0.0E+0
0.0E+0 0.0E+0 -0.26117002127659574 0.26117002127659574 0.0E+0
LDL of [K11] = 175176.54255319148 -0.2070856178827499 0.0E+0 0.0E+0
-0.7929143821172502 0.2070856178827499 -0.2070856178827499 0.0E+0
LDL \text{ of } [K11] = 77864.19232391396 \ 0.6305851063829787 \ -0.6305851063829787
-1.0 -0.36941489361702123 0.36941489361702123 0.0E+0
LDL of [K11] = 157038.27127659574 -0.11550223476828982 0.0E+0
0.11550223476828982 -1.0 0.0E+0
```

```
LDL of [K11] = 204043.26040931543 -0.24063524519998494 -

0.7593647548000151 0.0E+0 0.0E+0

LDL of [K11] = 176184.80946068066 -0.211623292466662 -

2.0648651936724454E-17 0.0E+0

LDL of [K11] = 78494.47532361932 0.6255217300016221 -0.6255217300016221

LDL of [K11] = 157286.88305692037 -0.11690029517761087

LDL of [K11] = 155137.45100017014
```

```
L D L t factorized of the "full rank" sub-matrix [K11] of Example 1 (singular case)
```

```
dependent row # i, u(i,i) = 3 0.0E+0
# dependent rows = 1
dependent rows = 3
```

LDL of [K11] = 1.0 -1.0LDL of [K11] = 1.0

L D L_t factorized of the "full rank" sub-matrix [K11] of Example 1 (non-singular case) # dependent rows = 0

APPENDIX C

A COMPLETE LISTING OF AN EDUCATIONAL FORTRAN SOURCE CODE OF "CHOLESKY GENERALIZED INVERSE" ALGORITHMS FOR SLE

implicit real*8(a-h,o-z) real tar(2) integer r, ractual

```
\mathbf{c}=
c.....Remarks:
    (a) Identifying which are dependent rows of a "floating" substructure
С
    (b) Factorizing (by cholesky) of floating substructure stiffness
С
       Whenever a dependent row is encountered during cholesky factored process,
С
       then we just:
С
       [1] set all factorized values of the dependent row to be ZEROES
С
       [2] ignore the dependent row(s) in all future factorized rows
С
    (c) [K "float"] = [K11] [K12]
С
                [K21] [K22]
С
       where [K11] = full rank ( = non-singular )
С
              = {K "float"] with deleting the dependent rows/columns
С
    (d) The u transpose * U of [K11] can be obtained by taking the results
С
       of part (b) and deleting the dependent rows/columns
С
c.....Author(s): Prof. Duc T. Nguyen
c.....Version: 02-11-2012 (EDUCATIONAL purpose,LDL/FULL matrix is assumed)
c.....Stored at: cd ~/cee/*odu*clas*/generalized inverse by cholesky.f
С
c.....Notes: Prof. Duc Nguyen's "generalized cholesky" code has been correctly
         verified for AT LEAST 7-8 different examples
c.....
         [see generalize*cholesky*.dat; and see out1-keep]
c.....
C=
    dimension u(1999,1999), idependrows(1999), tempo1(1999),
           am inv(1999,1999), tempo2(1999)
   $
    dimension ut(1999,1999), am(1999,1999), amt(1999,1999),
           g(1999,1999), gt(1999,1999)
   $
    dimension itempo1(1999), rhs(1999)
С
    call pierrotime(t1)
c
    write(6, *)'
    write(6,*) '==
    write(6,*) 'today date: 04-23-2012; Prof. Duc T. Nguyen'
    write(6,*) '=
                                             1
    write(6, *)'
С
    maxnr1 = 1999
```

```
maxnc1 = 1999
    maxnc2 = 1999
С
c.....input (or randomly generate rectangular/square matrix [G] of dimension mxn
c.....where we assume/prefer m > n
С
    read(5,*) m, n, iautodata, irankn, iaxeqb
    write(6,*) 'user input: m,n,iautodata,irankn,iaxeqb = '
    write(6,*) m,n,iautodata,irankn,iaxeqb
    write(6, *)'
С
c.....read user's input matrix data
С
    if (iautodata .eq. 0) then
    do 32 i=1, m
     read(5,*)(g(i,j), j=1,n)
32
     continue
С
c.....user's input rhs vector {rhs} nx1
С
     read(5,*) (rhs(i), i=1,m)
С
c.....randomly generated input matrix data
С
    elseif (iautodata .eq. 1) then
     ndependcols = n - irankn
     icount = 0
     idum=0
     do 61 i=1.n
     if (j.LE. irankn) then
     irandom col = irand(1, n)
     write(6,*) 'irandom col = ', irandom col
С
     do 60 i=1, m
     g(i,j) = drand(idum) * 10000.d0
60 continue
     elseif (j.GT. irankn) then
     do 66 i=1, m
     g(i,j) = 0.d0
66 continue
     endif
61 continue
С
c.....generated rhs vector {rhs} nx1, such that solution vector = \{1, 1, ..., 1\}
С
    do 67 i=1, n
    tempol(i) = 1.d0
67 continue
```

```
call mtimesv(g, tempo1, rhs, maxnr1, maxnc1, m, n)
   endif
С
   write(6,*) 'user input, or randomly generated matrix G mxn = '
    do 63 i=1, m
    write(6,*) (g(i,j), j=1,n)
63 continue
                   T
   write(6, *)'
С
   write(6,*) 'user input: right-hand-side (rhs) vector mx1 = '
    do 92 i=1, m
    write(6,*) 'rhs(-) = ',rhs(i)
92 continue
                   1
   write(6, *)'
С
С
   call transpose(g, gt, maxnr1, maxnc1, m, n)
С
C-----
   if (m.ge. n) then
   call mtimesm(gt, g, am, maxnr1, maxnc1, maxnc2, n, m, n) ! compute G' * G
   neq = n
С
   write(6,*) 'print [am] = G transpose * G = '
   do 72 i=1.n
    write(6,*) (am(i,j),j=1,n)
72 continue
                     1
   write(6, *)'
С
   elseif (m.lt. n) then
   call mtimesm(g, gt, am, maxnr1, maxnc1, maxnc2, m, n, m) ! compute G * G'
   neq = m
С
   write(6,*) 'print [am] = G * G_transpose = '
   do 73 i=1,m
    write(6,*) (am(i,j),j=1,m)
73 continue
                          +
   write(6, *)'
С
   endif
С
       c--
С
   call generalized inverse cholesky(neq,am,idependrows,
  $ ndependrows,r, maxnr1, independrows, itempo1)
С
```

```
write(6,*) 'special cholesky factor of Gt*G, or G*Gt '
    do 22 i=1, r
    write(6,*) (am(i,j), j=1, neq)
22 continue
С
    write(6,*) '# independent rows = ',independrows
    write(6, *)'
    if (independrows .ge. 1) then
    write(6,*) 'independent rows = ',(itempo1(i),i=1,independrows)
    endif
С
    call transpose(am, amt, maxnr1, maxnc1, r, neg)
    call mtimesm(am, amt, u, maxnr1, maxnc1, maxnc2, r, neq, r) ! compute L'* L
    nactual = r
    call generalized inverse cholesky(nactual, u, idependrows,
   $ ndependrows, ractual, maxnr1, independrows, itempo1)
   write(6,*) 'regular cholesky factorization of M*Mt '
   write(6,*) 'M = factorized of Gt^*G, or G^*Gt with deleted rows'
    write(6,*) '# dependent rows = ndependrows = ',ndependrows
    write(6, *)'
    if (iaxeqb.eq. 0) then ! find generalized inverse explicitly
С
c.....find the actual inverse of [u] !!
С
    do 43 irow = 1, nactual
    do 42 i=1, nactual
    tempo1(i) = 0.d0
42 continue
    tempo1(irow) = 1.d0
    call fbe cholesky(nactual, u, tempo1, maxnr1)
С
    do 44 i=1, nactual
    am inv(i,irow) = tempol(i)
     write(6,*) 'i=row#, irow=col#, am inv(-,-) = ',i,irow,tempo1(i)
C
    continue
44
С
43 continue
С
c.....applying the French's generalized inverse formula
С
    if (m .ge. n) then
    call mtimesm(amt, am inv, u, maxnr1, maxnc1, maxnc2, neq,
   $ nactual, nactual)
                                              ! compute L * [am inv]
```

call mtimesm(u, am inv, ut, maxnr1, maxnc1, maxnc2, neq, \$ nactual, nactual) ! compute L * [am inv] * [am inv] call mtimesm(ut, am, u, maxnr1, maxnc1, maxnc2, neq, ! compute L * [am inv] * [am inv] * L' \$ nactual, neq) call mtimesm(u, gt, ut, maxnr1, maxnc1, maxnc2, neq. ! compute L * [am inv] * [am inv] * L' * G' \$ neq. m) write(6,*) 'generalized inverse of [G] = 'do 52 i=1, neg write(6,*) (ut(i,j), j=1,m) С 52 continue С elseif (m.lt. n) then call mtimesm(gt, amt, u, maxnr1, maxnc1, maxnc2, n, m, ! compute G' * L \$ nactual) call mtimesm(u, am inv, ut, maxnr1, maxnc1, maxnc2, n, ! compute G' * L * [am inv] \$ nactual, nactual) call mtimesm(ut, am inv, u, maxnr1, maxnc1, maxnc2, n, ! compute G' * L * [am inv] * [am inv] \$ nactual, nactual) call mtimesm(u, am, ut, maxnr1, maxnc1, maxnc2, n, \$ nactual, neq) ! compute G' * L * [am inv] * [am inv] * \mathbf{L}^{\prime} write(6,*) 'generalized inverse of [G] = 'do 54 i=1, nwrite(6,*) (ut(i,j), j=1,m) 54 continue . write(6, *)' endif С c.....solution of [G] $\{x\} = \{rhs\}$ mxn nx1 mx1 С С c.....thus, $\{x\} = [G+] * \{rhs\}$ nx1 nxm mx1 С С call mtimesv(ut, rhs, tempo1, maxnr1, maxnc1, n, m) elseif (iaxeqb.eq. 1) then ! AVOID computing generalized inverse explicitly if (m.ge. n) then c.....compute [G'] * {rhs}; with results stored in {tempo1} call mtimesv(gt, rhs, tempo1, maxnr1, maxnc1, n, m) write(6,*) 'Gt * rhs = tempol = ',(tempol(i),i=1,n) С c.....compute [L'] * {tempo1}; with results stored in {tempo2} call mtimesv(am, tempo1, tempo2, maxnr1, maxnc1, nactual, n) write(6,*) 'Gt * rhs = tempo2 = ',(tempo2(i),i=1,n) C c.....now, doing forward & backward solutions, stored the results in {tempo2}

```
call fbe cholesky(nactual, u, tempo2, maxnr1)
```

```
c.....now, doing forward & backward solutions AGAIN, stored the results in {tempo2}
    call fbe cholesky(nactual, u, tempo2, maxnr1)
c.....finally, compute [L] * {tempo2} = same as compute [G+] * {rhs} !!
    call mtimesv(amt, tempo2, tempo1, maxnr1, maxnc1, n, nactual)
     elseif (m.lt. n) then
c.....compute [L'] * {rhs}; with results stored in {tempo1}
    call mtimesv(am, rhs, tempo1, maxnr1, maxnc1, nactual, n)
c.....now, doing forward & backward solutions, stored the results in {tempo1}
    call fbe cholesky(nactual, u, tempo1, maxnr1)
c.....now, doing forward & backward solutions AGAIN, stored the results in {tempo1}
    call fbe cholesky(nactual, u, tempo1, maxnr1)
c.....compute [L] * {tempo1}; with results stored in {tempo2}
    call mtimesv(amt, tempo1, tempo2, maxnr1, maxnc1, n, nactual)
c.....finally, compute [G'] * \{tempo2\} = same as compute [G+] * \{rhs\} !!
    call mtimesv(gt, tempo2, tempo1, maxnr1, maxnc1, n, m)
     endif
    endif
С
c.....output unknown solution vector \{x\} with 3 numbers:
c.....smallest (dabs\{x\}), biggest (dabs\{x\}), sum (dabs\{x\})
С
    abs smallest = 10**8
    abs biggest = 0.d0
    sum abs = 0.d0
    write(6,*) 'solution vector \{x\} = pinv(G) * \{rhs\} is ...'
    do 102 i=1, n
     aa = dabs(tempol(i))
     bb = tempo1(i)
     write(6,*) 'i, x(i) = ',i, bb
     if (aa .LT. abs smallest) abs smallest = aa
     if (aa .GT. abs biggest) abs biggest = aa
     sum abs = sum abs + aa
102 continue
                                 ,
    write(6, *)'
    write(6,*) 'abs smallest, abs biggest, sum abs = '
    write(6,*) abs_smallest, abs_biggest, sum_abs
    write(6, *)'
С
c.....output absolute and relative error norm
С
    call error norm(g,tempo1,rhs,maxnr1,maxnc1,m,n,abserr,relerr,
   $ tempo2)
```

```
write(6,*) 'abserr, relerr = ',abserr, relerr
```

```
C
    call pierrotime(t2)
    time a to z = t2 - t1
    write(6,*) 'time a to z = ', time a to z
    write(6, *)'
С
С
    stop
    end
subroutine generalized inverse cholesky(n,u,idependrows,
  $ ndependrows, r, maxnr1, independrows, independentrows)
    implicit real*8(a-h,o-z)
   integer r, ractual
    dimension u(maxnr1,*),idependrows(*),independentrows(*)
C=
c.....Remarks:
    (a) Identifying which are dependent rows of a "floating" substructure
С
    (b) Factorizing (by cholesky) of floating substructure stiffness
С
       Whenever a dependent row is encountered during cholesky process,
С
       then we just:
С
       [1] set all factorized values of the dependent row to be ZEROES
С
С
       [2] ignore the dependent row(s) in all future factorized rows
    (c) [K"float"] = [K11] [K12]
С
               [K21] [K22]
С
С
       where [K11] = full rank ( = non-singular )
              = {K "float"] with deleting the dependent rows/columns
С
    (d) The U transpose * U of [K11] can be obtained by taking the results
С
       of part (b) and deleting the dependent rows/columns
С
c.....Author(s): Prof. Duc T. Nguyen
c.....Version: 02-11-2012
c.....Stored at: cd ~/cee/*odu*clas*/generalized_inverse_by_ldl.f
c=
С
С
    write(6,*) 'check point #01'
    eps=0.000000001
```

с

r = 0

do 11 ir = 1, n

ndependrows = 0independrows = 0

98

```
write(6,*)' ir, r = ', ir, r
С
     do 12 icol=ir. n
      sum = u(ir, icol)
      do 13 iprevrow=1, r-1
       sum = sum - u(iprevrow,ir) * u(iprevrow,icol)
 13
        continue
     write(6,*) 'check point #02'
С
     if (ir .eq. icol) then
c..... cholesky factorized diagonal terms of row # ir
     write(6,*) 'check point #03'
С
      if (sum .gt. eps) then
       u(r,ir)=dsqrt(sum)
       write(6,*) ' u(r,ir) = ',u(r,ir)
С
       independrows = independrows + 1
       independentrows(independrows) = ir
     write(6,*) 'check point #04'
С
      else
       ndependrows = ndependrows + 1
       idependrows(ndependrows) = ir
С
       write(6,*) 'sum, u(r,ir) = \text{diag term are } \dots ',sum,u(r,ir)
С
       r = r - 1
     write(6,*) ' ir, r = ', ir, r
С
     write(6,*) 'check point #05'
С
       go to 11
      endif
С
     else
c..... cholesky factorized off-diagonal terms of row # ir
     write(6,*) 'check point #06'
С
      u(r,icol) = sum/u(r,ir)
     endif
С
      continue
12
С
11
      continue
С
c.....all lower triangular of cholesky factorized [U] are set to zero !
```

c

r = r + 1

```
do 22 ir=1, r
  independr = independentrows(ir)
   do 23 icol=1, independr-1
23 u(ir,icol)=0.d0
22 continue
С
   write(6,*) 'ndependrows = ', ndependrows
С
   write(6,*) 'dependent rows = ',(idependrows(i), i=1,ndependrows)
С
С
  return
  end
subroutine transpose(a, at, maxnr1, maxnc1, nr1, nc1)
  implicit real*8(a-h,o-z)
  dimension a(maxnr1,*), at(maxnc1,*)
С
  do 1 i=1, nr1
   do 2 = 1, ncl
   at(j,i) = a(i,j)
2 continue
1
  continue
  return
  end
subroutine mtimesm(a, b, c, maxnr1, maxnc1, maxnc2, nr1,nc1,nc2)
  implicit real*8(a-h,o-z)
  dimension a(maxnr1,*), b(maxnc1,*), c(maxnr1,*)
С
  do 1 j=1, nc2
   do 2 i=1, nr1
   c(i,j) = 0.d0
   do 3 k=1, ncl
   c(i,j) = c(i,j) + a(i,k) * b(k,j)
   continue
3
2
  continue
  continue
1
  return
  end
subroutine fbe cholesky(n, u, rhs, maxnr1)
  implicit real*8(a-h, o-z)
С
  dimension u(maxnr1,*), rhs(*)
С
```

```
С
c.....forward cholesky solution
С
   do 1 i = 1, n
   sum = rhs(i)
   do 2 i=1, j-1
   sum = sum - u(i,j) * rhs(i)
2
  continue
   rhs(j) = sum/u(j,j)
1
   continue
С
c.....backward cholesky solution
С
   do 4 j=n, 1, -1
   sum = rhs(j)
   do 5 i=j+1, n
   sum = sum - u(j,i) * rhs(i)
5 continue
   rhs(j) = sum/u(j,j)
4
   continue
С
   return
   end
subroutine mtimesv(g, tempo1, rhs, maxnr1, maxnc1, m, n)
   implicit real*8(a-h,o-z)
С
   dimension g(maxnr1,*), tempo1(*), rhs(*)
Ç
   do 1 i=1, m
   sum = 0.d0
   do 2 j=1, n
   sum = sum + g(i,j) * tempo1(j)
2
   continue
   rhs(i) = sum
   continue
1
С
   return
   end
subroutine error norm(a,x,b,maxnr1,maxnc1,nr1,nc1,abserr,relerr,
  $ tempo1)
   implicit real*8(a-h,o-z)
   dimension a(maxnr1,*), x(*), b(*), tempo1(*)
С
```

```
call mtimesv(a, x, tempo1, maxnr1, maxnc1, nr1, nc1)
С
    abserr = 0.d0
    relerr = 0.d0
С
   do 1 i=1, nr1
   x(i) = tempol(i) - b(i)
   abserr = abserr + x(i)^{**2}
   relerr = relerr + b(i)^{**2}
1
   continue
    abserr = dabs(abserr)
    relerr = dabs(relerr)
    relerr = abserr/relerr
С
   return
   end
subroutine pierrotime (time)
   real tar(2)
   real*8 time
С
    С
   purpose :
    This routine returns the user + system execution time
С
    The argument tar returns user time in the first element and
С
    system time in the second element. The function value is the
С
    sum of user and system time. This value approximates the
С
    program's elapsed time on a quiet system.
С
С
    Uncomment for your corresponding platform
С
С
    Note: On the SGI the resolution of etime is 1/HZ
С
С
   Output
С
    time: user+sytem executime time
С
С
   *****
   SUN-Solaris
С
   time=etime(tar)
   HP - HPUX
С
С
   time=etime (tar)
                        !f90
   time=etime (tar)
                        !f77
С
С
   COMPAQ - alpha
```

101

c time=etime(tar)

c CRAY

- c time=tsecnd()
- c IBM
- c time=0.01*mclock()
- c SGI origin
- c time=etime(tar)

C.1: A Complete Input File of an Educational FORTRAN Source Code of "Generalized Inverse" Algorithms

For SLE

7 7 0 4 1

1.d0	2.d0	-3.d0	2.d0	-2.d0	-3.d0	-2.d0
2.d0	4.d0	-6.d0	4.d0	-4.d0	-6.d0	-4.d0
-3.d0	-6.d0	9.d0	-6.d0	6.d0	9.d0	6.d0
2.d0	4.d0	-6.d0	5.d0	-1.d0	-5.d0	-7.d0
-2.d0	-4.d0	6.d0	-1.d0	13.d0	9.d0	-5.d0
-3.d0	-6.d0	9.d0	-5.d0	9.d0	13.d0	9.d0
-2.d0	-4.d0	6.d0	-7.d0	-5.d0	9.d0	27.d0
-2.d0	-4.d0	6.d0	-7.d0	-5.d0	9.d0	27.d0

Note: MATLAB solution = (0, 0, 0, 0, 0, 0, 1) = satisfy SLE

C.2: A Complete Output File of an Educational FORTRAN Source Code of "Generalized Inverse" Algorithms

For SLE

today date: 04-23-2012; Prof. Duc T. Nguyen

user input: m,n,iautodata,irankn,iaxeqb = 77041

user input, or randomly generated matrix G mxn = 1.0 2.0 -3.0 2.0 -2.0 -3.0 -2.0 2.0 4.0 -6.0 4.0 -4.0 -6.0 -4.0 -3.0 -6.0 9.0 -6.0 6.0 9.0 6.0 2.0 4.0 -6.0 5.0 -1.0 -5.0 -7.0 -2.0 -4.0 6.0 -1.0 13.0 9.0 -5.0 -3.0 -6.0 9.0 -5.0 9.0 13.0 9.0 -2.0 -4.0 6.0 -7.0 -5.0 9.0 27.0

user input: right-hand-side (rhs) vector mx1 =rhs(-) = -2.0 rhs(-) = -4.0 rhs(-) = 6.0 rhs(-) = -7.0 rhs(-) = -5.0 rhs(-) = 9.0 rhs(-) = 27.0

```
print [am] = G_transpose * G =
35.0 70.0 -105.0 69.0 -73.0 -127.0 -113.0
70.0 140.0 -210.0 138.0 -146.0 -254.0 -226.0
-105.0 -210.0 315.0 -207.0 219.0 381.0 339.0
69.0 138.0 -207.0 156.0 -84.0 -246.0 -320.0
-73.0 -146.0 219.0 -84.0 332.0 278.0 -56.0
-127.0 -254.0 381.0 -246.0 278.0 482.0 434.0
-113.0 -226.0 339.0 -320.0 -56.0 434.0 940.0
```

```
special cholesky factor of Gt*G, or G*Gt

5.916079783099616 11.832159566199232 -17.74823934929885 11.663128715253528

-12.339252119036342 -21.46691807010432 -19.100486156864473

0.0E+0 0.0E+0 0.0E+0 4.468940430507951 13.406821291523839

0.9781800942313469

-21.756515429211084

0.0E+0 0.0E+0 0.0E+0 0.0E+0 4.496064087029694 10.065074253426936

0.0E+0 0.0E+0 0.0E+0 0.0E+0 0.0E+0 0.7209335773362233

# independent rows = 4
```

```
independent rows = 1 4 6 7
regular cholesky factorization of M*Mt
M = factorized of Gt*G, or G*Gt with deleted rows
# dependent rows = ndependrows = 0
```

```
solution vector \{x\} = pinv(G) * \{rhs\} is ...
i, x(i) = 1 \ 1.6874571068360796E-13
i, x(i) = 2 \ 3.3749142136721593E-13
i, x(i) = 3 \ -5.06237132050824E-13
```

i, x(i) = 4 -2.8787848147707946E-13 i, x(i) = 5 -2.2136011299001E-12 i, x(i) = 6 2.915937886998499E-12 i, x(i) = 7 0.9999999999987078

abs_smallest, abs_biggest, sum_abs = 1.6874571068360796E-13 0.999999999999987078 1.0000000000051376

abserr, relerr = 8.793202261721907E-25 9.354470491193518E-28 time_a_to_z = 9.539998136460781E-4

APPENDIX D

MatlabMPI SOURCE CODE FOR MATRIX-MATRIX MULTIPLICATION (MATRIX IN DENSE FORMAT)

%Initialize MPI MPI_Init;

%Create communicator comm = MPI_COMM_WORLD;

%Get Size and Rank comm_size = MPI_Comm_size(comm); %numtasks my_rank = MPI_Comm_rank(comm); %taskid

nn = 1000; % Matrix size

```
%tStart = tic;
```

% Master Processor task if (my_rank==0)

```
% a = 10*rand(nn);
% b = 10*rand(nn);
```

```
for i = 1: nn
    for j = 1:nn
        a(i,j) = (i-1)+(j-1);
    end
end
a;
for i = 1: nn
    for j = 1:nn
        b(i,j) = (i+1)*(j+1);
    end
end
b;
Z = zeros(nn);
tStart = tic;
```

domains = comm_size-1; % numworkers

```
%divide matrix "b" to parts (domains)
len = floor(length(b)/domains);
```

```
for i = 1:domains-1
     MPI_Send(i,1,comm,a(:,:),b(:,((i-1)*len)+1:i*len)) %send parts of matrices to slaves
     sent part = sprintf('\%g', i)
     i;
    disp('******');
  end
  MPI Send(domains, 1,comm,a(:,:),b(:,((domains-1)*len)+1:length(b))); %last part to
slave
  disp('last part sent');
  disp('******');
  for i = 1:domains
     Z = MPI Recv(i, 100, comm);
    Ζ;
     size(Z);
     recv = sprintf('%g', i);
  end
end %end master
if my rank > 0 %slave
  [matrix a matrix b] = MPI Recv(0,1,comm);
  %Computation
  [ra ca] = size(matrix_a);
  [rb cb] = size(matrix_b);
  for k = 1: cb
     for i = 1:ra
       c(i,k) = 0;
       for j = 1:ca
         c(i,k) = c(i,k) + matrix_a(i,j) * matrix_b(j,k);
       end
     end
  end
```

Z = c;

MPI_Send(0,100,comm,Z); exit;

end %slave

MPI_Finalize; tElapsed = toc(tStart) disp('Success');

APPENDIX E

GRAPHICAL COMPARISONS (IN TERMS OF COMPUTATIONAL TIMES) OF ODU-GINVERSE WITH OTHER ALGORITHMS

In this appendix, we graphically compare the computational times of ODUginverse with other existing algorithms. The description of the test problems can be found in section 5.

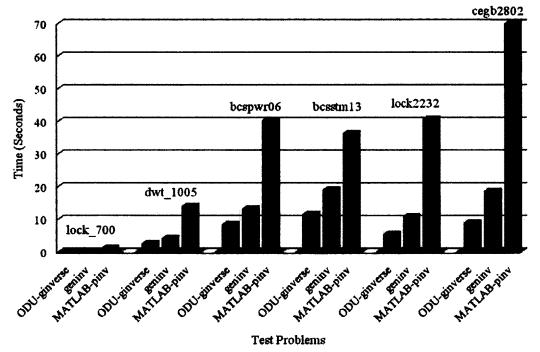


Figure E.1 Computational Times (in seconds) for Symmetric Rank-Deficient Test Matrices with RHS Vector as Linear Combination of Columns of Coefficient Matrix

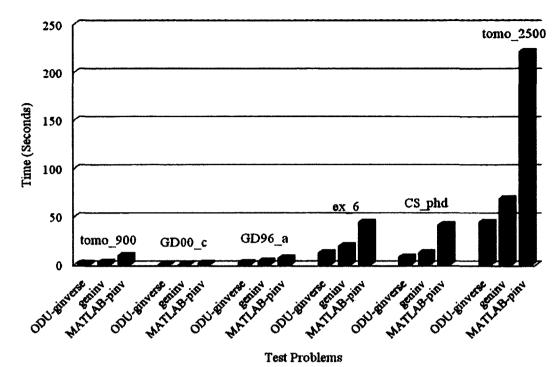


Figure E.2 Computational Times (in seconds) for Non-Symmetric Rank-Deficient Test Matrices with RHS Vector as Linear Combination of Columns of Coefficient Matrix

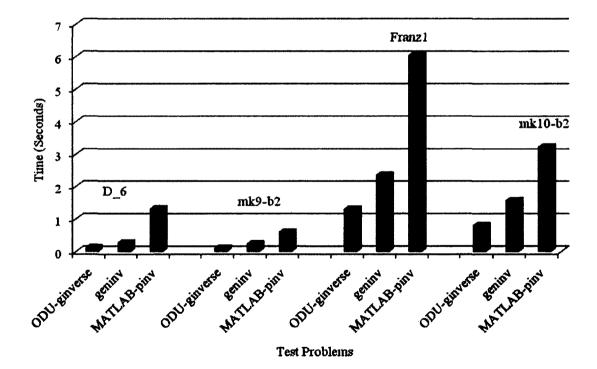


Figure E.3 Computational Times (in seconds) for Rectangular Rank-Deficient Test Matrices (tall type) with RHS Vector as Linear Combination of Columns of Coefficient Matrix

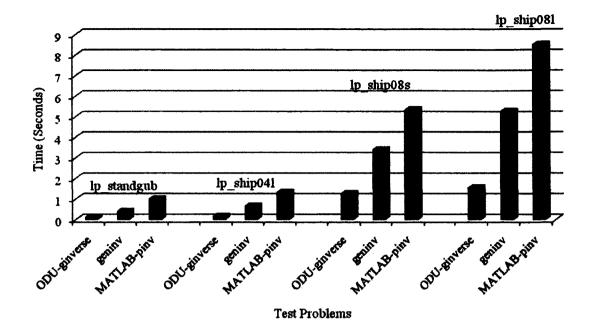


Figure E.4 Computational Times (in seconds) for Rectangular Rank-Deficient Test Matrices (fat type) with RHS Vector as Linear Combination of Columns of Coefficient Matrix

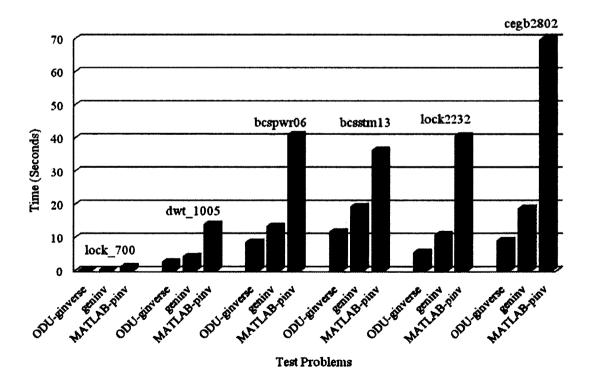


Figure E.5 Computational Times (in seconds) for Symmetric Rank-Deficient Test Matrices with Randomly generated RHS Vector

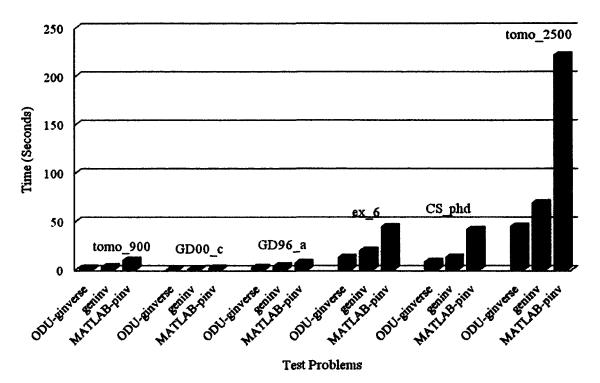


Figure E.6 Computational Times (in seconds) for Non-Symmetric Rank-Deficient Test Matrices with Randomly generated RHS Vector

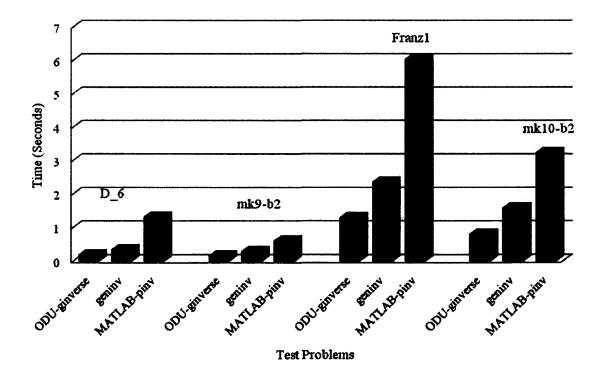


Figure E.7 Computational Times (in seconds) for Rectangular Rank-Deficient Test Matrices (tall type) with Randomly generated RHS Vector

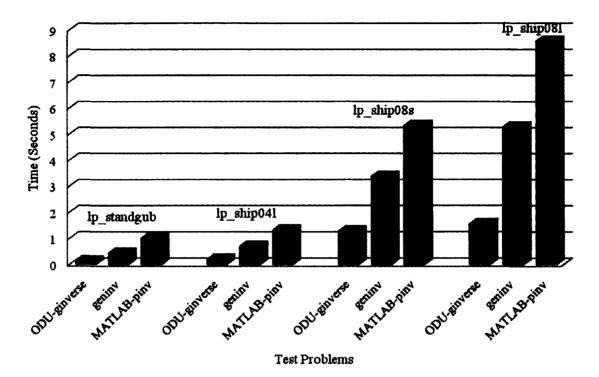


Figure E.8 Computational Times (in seconds) for Rectangular Rank-Deficient Test Matrices (fat type) with Randomly generated RHS Vector

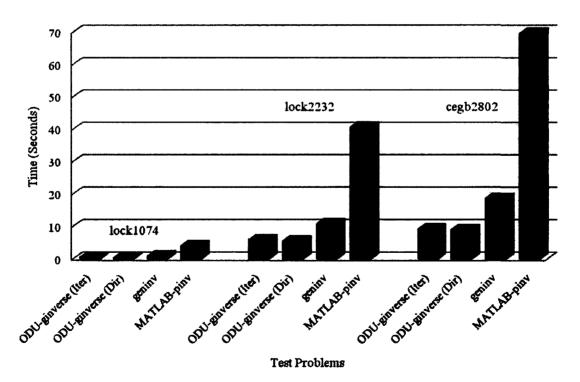


Figure E.9 Computational Times (in seconds) for Symmetric Rank-Deficient Test Matrices with Randomly generated RHS Vector (Iterative Solver inside Generalized Inverse)

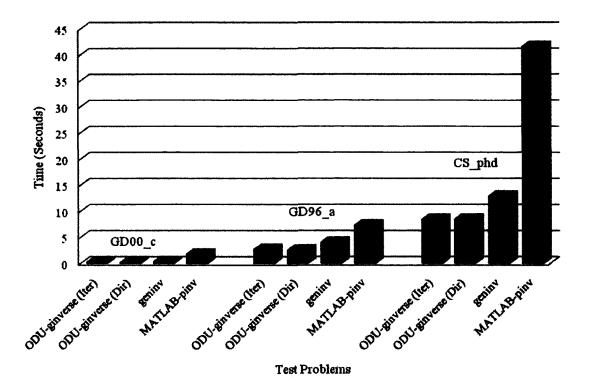


Figure E.10 Computational Times (in seconds) for Non-Symmetric Rank-Deficient Test Matrices with Randomly generated RHS Vector (Iterative Solver inside Generalized Inverse)

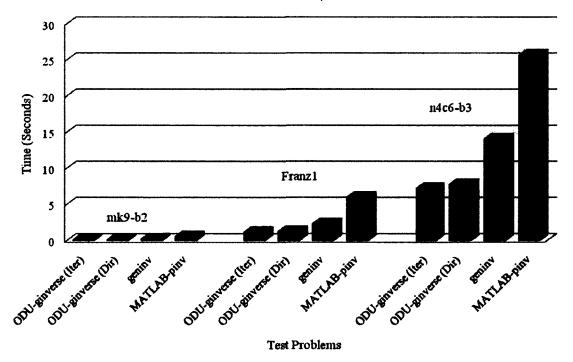


Figure E.11 Computational Times (in seconds) for Rectangular Rank-Deficient Test Matrices (tall) with Randomly generated RHS Vector (Iterative Solver inside Generalized Inverse)

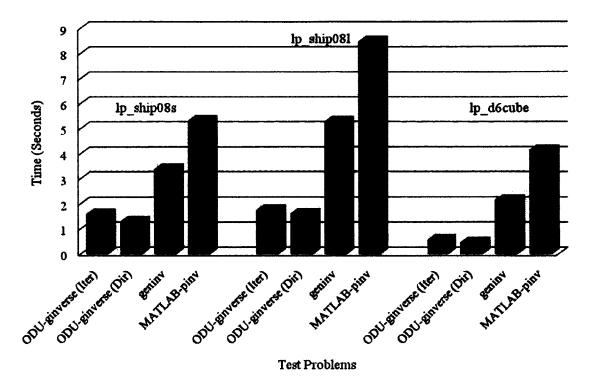


Figure E.12 Computational Times (in seconds) for Rectangular Rank-Deficient Test Matrices (fat) with Randomly generated RHS Vector (Iterative Solver inside Generalized Inverse)

VITAE

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