# A Comparative Evaluation of Matlab, Octave, FreeMat, Scilab, and R on Tara 

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#### Abstract

Matlab is the most popular commercial package for numerical computations in mathematics, statistics, the sciences, engineering, and other fields. Octave, FreeMat and Scilab are free numerical computational packages that have many of the same features as Matlab. $R$ is a free Statistical package. Although $R$ does not belong to the same line of products as Matlab, it is similar to Matlab in its computational capabilities. These packages are available to download on the Linux, Windows, and Mac OS X operating systems. We investigate whether they are viable alternatives to Matlab for uses in research and teaching. We compare the results on the cluster tara in the UMBC High Performance Computing Facility with 86 nodes, each with two quadcore Intel Nehalem processors and 24 GB of memory. The tests focused on usability lead us to conclude that the package Octave is the most compatible with Matlab, since it uses the same syntax and has the native capability of running m-files. Both FreeMat and Scilab were hampered by somewhat different syntax or function names and some missing functions. The tests focused on efficiency show that Matlab and Octave are fundamentally able to solve problems of the same size and with equivalent efficiency in absolute times, except in one test dealing with a very large problem. FreeMat and also Scilab exhibit significant limitations on the problem size and the efficiency of the problems they can solve in our tests. The syntax of $R$ is significantly different from that of Matlab, Octave, FreeMat, and Scilab. R too exhibited certain limitations on the size of problems it could solve for and its performance was similar to that of FreeMat and Scilab. In summary, we conclude that Octave is the best viable alternative to Matlab because it was not only fully compatible (in terms of syntax) with Matlab in our tests, but it also performed very well.


## 1 Introduction

### 1.1 Overview

There are several numerical computational packages that serve as educational tools and are also available for commercial use. Matlab is the most widely used such package. The focus of this study is to introduce the three additional numerical computational packages Octave, FreeMat, and Scilab, and the statistical package R, and provide information on which package is most compatible to Matlab users. Section 1.3 provides more detailed descriptions of these packages. To evaluate GNU Octave, FreeMat, and Scilab, a comparative approach is used
based on a Matlab user's perspective. To achieve this task, we perform some basic and some complex studies on Matlab, GNU Octave, FreeMat, Scilab, and R. The basic studies include basic operations solving systems of linear equations, computing the eigenvalues and eigenvectors of a matrix, and two-dimensional plotting. The complex studies include direct and iterative solutions of a large sparse system of linear equations resulting from finite difference discretization of an elliptic test problem. This report extends the report [3] by adding the package R to the comparison. That report was based on $[8,10]$, which considered the usability of software packages in a home computer setting, and extends them substantially by more substantive efficiency comparisons.

In Section 2, we perform the basic operations test using Matlab, GNU Octave, FreeMat, Scilab, and R. This includes testing the backslash operator, computing eigenvalues and eigenvectors, and plotting in two dimensions in all the packages except $R$, which follows a significantly different syntax and functions. The backslash operator works identically for all of the packages (except R) to produce a solution to the linear system given. The command eig has the same functionality in Octave and FreeMat as in Matlab for computing eigenvalues and eigenvectors, whereas Scilab uses the equivalent command spec to compute them. Plotting is another important feature we analyze by an m-file containing the twodimensional plot function along with some common annotations commands. Once again, Octave and FreeMat use the exact commands for plotting and similar for annotating as Matlab whereas Scilab requires a few changes. For instance in Scilab, the pi command is defined using \%pi and the command grid on from Matlab is replaced with xgrid or use the translator to create conversion. To overcome these conversions, we find that we can use the Matlab to Scilab translator, which takes care of these command differences for the most part. The translator is unable to convert the xlim command from Matlab to Scilab. To rectify this, we must manually specify the axis boundaries in Scilab using additional commands in Plot2d. This issue brings out a major concern that despite the existence of the translator, there are some functions that require manual conversion.

Section 3 introduces a test problem, given by the Poisson equation with homogeneous Dirichlet boundary conditions, and discusses the finite discretization for the problem in two dimensional space. In the process of finding a solution, we use a direct method, Gaussian elimination, and an iterative method, the conjugate gradient method. To be able to analyze the performance of these methods, we solve the problem on progressively finer meshes. The Gaussian elimination method built into the backslash operator successfully solves the problem up to a mesh resolution of $4,096 \times 4,096$ in both Matlab and Octave, while the Gaussian elimination method built into the backslash operator in FreeMat successfully solves the problem up to a mesh resolution of $2,048 \times 2,048$. However, in Scilab and R, it is only able to solve up to a mesh resolution of $1,024 \times 1,024$. The conjugate gradient method is implemented in the pcg function, which is stored in Matlab and Octave as a m-file. This function is also available in Scilab as a sci file. The matrix free implementation of the conjugate gradient method allows us to solve a mesh resolution up to $8,192 \times 8,192$ for Matlab and Octave. Scilab is able to solve the system for a resolution up to $4,096 \times 4,096$. In FreeMat and R, we wrote our own cg function because they do not have a built in pcg function. In FreeMat we were able to solve the system for a resolution of $2,048 \times 2,048$ and in R for a resolution of
$4,096 \times 4,096$. The existence of the mesh in Matlab, Octave, and Scilab allows us to include the three-dimensional plots of the numerical solution and the numerical error. However, in FreeMat, mesh does not exist and we had to use plot3, which results in a significantly different looking plot. In R too, mesh does not exist and we had to use the perp function.

The syntax of Octave is identical to that of Matlab in our tests. However, we found during our tests that FreeMat lacks a number of functions, such as kron for Kronecker products, pcg for the conjugate gradient method, and mesh for three-dimensional plotting. Otherwise, FreeMat is very much compatible with Matlab. Even though Scilab is designed for Matlab users to smoothly utilize the package and has a m-file translator, it often still requires manual conversions. The tests focused on usability lead us to conclude that the packages Octave and FreeMat are most compatible with Matlab, since they use the same syntax and have the native capability of running $m$-files. Among these two packages, Octave is a significantly more mature software and has significantly more functions available for use.

The results of the numerical experiments on the complex test problem in Section 3 are clearly somewhat surprising. Fundamentally, Matlab and Octave turn out to be able to solve problems of the same size, as measured by the mesh resolution of the problem they were able to solve. More precisely, in the tests of the backslash operator, it becomes clear that Matlab is able to solve the problem faster, which points to more sophisticated internal optimization of this operator. However, considering absolute run times, Octave's handling of the problem is acceptable, except for very large cases. What is surprising is that FreeMat and also Scilab exhibit significant limitations on the problem size they are able to solve by the backslash operator. The tests using the iterative conjugate gradient method also give some surprising results. Whether the function for this method is supplied by the package or hand-written, only basic matrix-vector operations are used by the method. Therefore, it is not clear why FreeMat and Scilab are less effective at solving the problems. R too exhibited limitations on the size of problems it could solve for. Specifically, R could only solve for resolutions until 1,024 using Gaussian elimination as it ran out of the allocated memory. Matlab and Octave both took about the same time to solve the problem, thus confirming the conclusions from the other tests. The data that these conclusions are based on are summarized in Tables 3.1, $3.2,3.3,3.4$, and 3.5.

The next section contains some additional remarks on features of the software packages that might be useful, but go beyond the tests performed in Sections 2 and 3. Sections 1.3 and 1.4 describe the four numerical computational packages in more detail and specify the computing environment used for the computational experiments, respectively.

### 1.2 Additional Remarks

### 1.2.1 Ordinary Differential Equations

One important feature to test would be the ODE solvers in the packages under consideration. For non-stiff ODEs, Matlab has three solvers: ode113, ode23, and ode45 implement an Adams-Bashforth-Moulton PECE solver and explicit Runge-Kutta formulas of orders 2 and 4, respectively. For stiff ODEs, Matlab has four ODE solvers: ode15s, ode23s, ode23t, and
ode23tb implement the numerical differentiation formulas, a Rosenbrock formula, a trapezoidal rule using a "free" interpolant, and an implicit Runge-Kutta formula, respectively.

According to their documentations, Octave and Scilab solve non-stiff ODEs using the Adams methods and stiff equations using the backward differentiation formulas. These are implemented in lsode in Octave and ode in Scilab. The only ODE solver in FreeMat is ode45 which solves the initial value problem probably by a Runge-Kutta method. In R, two packages can be used to solve ODEs: deSolve for initial value problems and bvpSolve for boundary value problems. Many of the functions in the deSolve package use both Adams and Runge-Kutta methods. These functions are implemented based on FORTAN's LSODA implementation that switches automatically between stiff and non-stiff systems [11].

It becomes clear that all software packages considered have at least one solver. Matlab, Octave, Scilab and R have state-of-the-art variable-order, variable-timestep methods for both non-stiff and stiff ODEs available, with Matlab's implementation being the richest and its stiff solvers being possibly more efficient. FreeMat is clearly significantly weaker than the other packages in that it does not provide a state-of-the-art ODE solver, particularly not for stiff problems.

### 1.2.2 Parallel Computing

Parallel Computing is a well-established method today. It takes in fact two forms: sharedmemory and distributed-memory parallelism.

On multi-core processors or on multi-core compute nodes with shared memory among all computational cores, software such as the numerical computational packages considered here automatically use all available cores, since the underlying libraries such as BLAS, LAPACK, etc. use them; studies for Matlab in [9] demonstrated the effectiveness of using two cores in some but not all cases, but the generally very limited effectiveness of using more than two cores. Since the investigations in [9], the situation has changed such that the number of cores used by Matlab, for instance, cannot even be controlled by the user any more, since that control was never respected by the underlying libraries anyway. ${ }^{1}$ Thus, even the 'serial' version for Matlab and the other numerical computational packages considered here are parallel on the shared memory of a compute node, this has at least potential for improved performance, and this feature is included with the basic license fee for Matlab. More recently, Matlab has started to provide built-in support for graphics processing units (GPUs). This is cutting-edge and should give a very significant advantage of Matlab over any other packages.

Still within the first form of parallel computing, Matlab offers the Parallel Computing Toolbox for a reasonable, fixed license fee (a few hundred dollars). This toolbox provides commands such as parfor as an extension of the for loop that allow for the farming out of independent jobs in a master-worker paradigm. This is clearly useful for parameter studies, parametrized by the for loop. The parfor extension uses all available compute nodes assigned to the Matlab job, thus it does go beyond using one node, but each job is 'serial' and lives only on one node.

[^0]The shared-memory parallelism discussed so far limits the size of any one job also that of a worker in a master-worker paradigm, to the memory available on one compute node. The second form of parallelism given by distributed-memory parallelism by contrast pools the memory of all nodes used by a job and thus enables the solution of much larger problems. It clearly has the potential to speed up the execution of jobs over 'serial' jobs, even if they use all computational cores on one node. Matlab offers the MATLAB Distributed Computing Server, which allows exactly this kind of parallelism. However, this product requires a substantial additional license fee that - moreover - is not fixed, but scales with the number of compute nodes (in total easily in the many thousands of dollars for a modest number of nodes).

The potential usefulness of distributed-memory parallelism, the historically late appearance of the Matlab product for it, and its very substantial cost has induced researchers to create alternatives. These include for instance pMatlab ${ }^{2}$ that can be run on top of either Matlab or other "clones" such as Octave. Particularly for pMatlab, a tutorial documentation and introduction to parallel computing is available [6]. In turn, since this distributed-memory parallel package is known to run in conjunction with Octave, this is one more reason to stir the reader to this numerical computational package to consider as alternative to Matlab. The statistical software R also supports parallel computing with the addition of optional open-source packages ${ }^{3}$. In particular, the SNOW (Simple Network of Workstations) package provides a master-worker paradigm. Functions such as clusterCall (evaluate a function on every worker with identical arguments) and parApply (apply a function to the rows or columns of a matrix) are used to distribute computations to workers in a cluster. Another package Rmpi allows MPI communications to be used within R, either in master-worker or single program multiple data (SPMD) paradigms. In addition, in recent years R has added packages ${ }^{4}$ that target GPUs too.

### 1.2.3 Applicability of this Work

Numerical computational packages see usage both in research and in teaching.
In a research context, an individual researcher is often very concerned with the portability of research code and reproducibility of research results obtained by that code. This concern applies over long periods of time, as the researcher changes jobs and affiliations. The software Matlab, while widely available at academic institutions, might not be available at some others. Or even if it is available, it is often limited to a particular computer (as fixed-CPU licenses tend to be cheaper than floating license keys). Freely downloadable packages are an important alternative, since they can be downloaded to the researchers own desktop for convenience or to any other or to multiple machines for more effective use. The more complex test case in Section 3 is thus designed to give a feel for a research problem. Clearly, the use of alternatives assumes that the user's needs are limited to the basic functionalities of Matlab itself; Matlab does have a very rich set of toolboxes for a large variety of applications or for certain areas with more sophisticated algorithms. If the use of one of them is profitable or

[^1]integral to the research, the other packages are likely not viable alternatives.
In the teaching context, two types of courses should be distinguished: There are courses in which Matlab is simply used to let the student solve larger problems (e.g., to solve eigenvalue problems with larger matrices than $4 \times 4$ ) or to let the student focus on the application instead of mathematical algebra (e.g., solve linear system quickly and reliably in context of a physics or a biology problem). We assume here that the instructor or the textbook (or its supplement) give instructions on how to solve the problem using Matlab. The question for the present recommendation is then, whether instructions written for Matlab would be applicable nearly word-for-word in the alternative package. Thus, you would want function names to be the same (e.g., eig for the eigenvalue calculation) and syntax to behave the same. But it is not relevant if the underlying numerical methods used by the package are the same as those of Matlab. And due to the small size of problems in a typical teaching context, efficiency differences between Matlab and its alternatives are not particularly crucial.

Another type of course, in which Matlab is used very frequently, is courses on numerical methods. In those courses, the point is to explain - at least for a simpler version - the algorithms actually implemented in the numerical package. It becomes thus somewhat important what the algorithm behind a function actually is, or at least its behavior needs to be the same as the algorithm discussed in class. Very likely in this context, the instructor needs to evaluate himself/herself on a case-by-case basis to see if the desired learning goal of each programming problem is met. We do feel that our conclusions here apply to most homework encountered in typical introductory numerical analysis courses, and the alternatives should work fine to that end. But as the above discussion on ODE solver shows, there are limitations or restrictions that need to be incorporated in the assignments. For instance, to bring out the difference in behavior between stiff and non-stiff ODE solvers, the ones available in Octave are sufficient, even if their function names do not agree with those in Matlab and their underlying methods are not exactly the same; but FreeMat's single ODE solver is not sufficient to conduct the desired computational comparison between methods from two classes of solvers.

### 1.3 Description of the Packages

### 1.3.1 Matlab

"MATLAB is a high-level language and interactive environment that enables one to perform computationally intensive tasks faster than with traditional programming languages such as C, C++, and Fortran." The web page of the MathWorks, Inc. at www.mathworks. com states that Matlab was originally created by Cleve Moler, a Numerical Analyst in the Computer Science Department at the University of New Mexico. The first intended usage of Matlab, also known as Matrix Laboratory, was to make LINPACK and EISPACK available to students without facing the difficulty of learning to use Fortran. Steve Bangert and Jack Little, along with Cleve Moler, recognized the potential and future of this software, which led to establishment of MathWorks in 1983. As the web page states, the main features of Matlab include high-level language; 2-D/3-D graphics; mathematical functions for various fields;
interactive tools for iterative exploration, design, and problem solving; as well as functions for integrating MATLAB-based algorithms with external applications and languages. In addition, Matlab performs the numerical linear algebra computations using for instance Basic Linear Algebra Subroutines (BLAS) and Linear Algebra Package (LAPACK).

### 1.3.2 GNU Octave

"GNU Octave is a high-level language, primarily intended for numerical computations," as the reference for more information about GNU Octave is www.octave.org. This package was developed by John W. Eaton and named after Octave Levenspiel, a professor at Oregon State University, whose research interest is chemical reaction engineering. At first, it was intended to be used with an undergraduate-level textbook written by James B. Rawlings of the University of Wisconsin-Madison and John W. Ekerdt of the University of Texas. This book was based on chemical reaction design and therefore the prime usage of this software was to solve chemical reactor problems. Due to complexity of other softwares and Octave's interactive interface, it was necessary to redevelop this software to enable the usage beyond solving chemical reactor design problems. The first release of this package, primarily created by John W. Eaton, along with the contribution of other resources such as the users, was on January 4, 1993.

Octave, written in C++ using the Standard Template Library, uses an interpreter to execute the scripting language. It is a free software available for everyone to use and redistribute with certain restrictions. Similar to Matlab, GNU Octave also uses for instance the LAPACK and BLAS libraries. The syntax of Octave is very similar to Matlab, which allows Matlab users to easily begin adapting to the package. Octave is available to download on different operating systems like Windows, Mac OS, and Linux. To download Octave go to http://sourceforge.net/projects/octave. A unique feature included in this package is that we can create a function by simply entering our code on the command line instead of using the editor.

### 1.3.3 FreeMat

FreeMat is a numerical computational package designed to be compatible with other commercial packages such as Matlab and Octave. The supported operating systems for FreeMat include Windows, Linux, and Mac OS X. Samit Basu created this program with the hope of constructing a free numerical computational package that is Matlab friendly. The web page of FreeMat at www.freemat.org. states that some of features for FreeMat include eigenvalue and singular value decompositions, $2 \mathrm{D} / 3 \mathrm{D}$ plotting, parallel processing with MPI, handle-based graphics, function pointers, etc. To download FreeMat, go to http: //sourceforge.net/projects/freemat.

### 1.3.4 Scilab

"Scilab is an open source, cross-platform numerical computational package as well as a high-level, numerically oriented programming language." Scilab was written by INRIA, the

French National Research Institution, in 1990. The web page for Scilab at www.scilab.org states the syntax is largely based on Matlab language. This software is also intended to allow Matlab users to smoothly utilize the package. In order to help in this process, there exists a built in code translator which assists the user in converting their existing Matlab codes into a Scilab code. According to Scilab's web page, the main features of Scilab include hundreds of mathematical functions; high-level programming language; 2-D/3-D visualization; numerical computation; data analysis; and interface with Fortran, C, C++, and Java. Just like Octave, Scilab is also a free software distributed under CeCILL licenses.

Scilab is fully compatible with Linux, Mac OS, and Windows platforms. Like Octave, the source code is available for usage as well as for editing. To download Scilab go to www.scilab.org/products/scilab/download Scilab also uses, for instance, the numerical libraries, LAPACK and BLAS. Unlike Octave, the syntax and built-in Scilab functions may not entirely agree with Matlab.

### 1.3.5 R

$R$ is an open source, cross-platform numerical computational and statistical package as well as a high-level, numerically oriented programming language used for developing statistical software and data analytics. The R programming language is based on the S programming language, which was developed at Bell Laboratories. R was not developed as a Matlab clone, and therefore has a different syntax than the numerical analysis software packages previously introduced. It is part of the GNU project and can be downloaded from www.r-project.org. Add-on packages contributed by the R user community can be downloaded from CRAN (The Comprehensive R Archive Network) at cran.r-project.org.

### 1.4 Description of the Computing Environment

The computations for this study are performed using Matlab R2011a, GNU Octave 3.0.4, FreeMat v4.0, Scilab-5.3.1, and R 2.13.0 under the Linux operating system Redhat Enterprise Linux 5. The Cluster tara in the UMBC High Performance Computing Facility (www.umbc. edu/hpcf) is used to carry out the computations and has a total of 86 nodes, with 82 used for computation. Each node features two quad-core Intel Nehalem X5550 processors (2.66 GHz, $8,192 \mathrm{kB}$ cache per core) with 24 GB of memory.

## 2 Basic Operations Test

This section examines a collection of examples inspired by some basic mathematics courses. This set of examples was originally developed for Matlab by the Center for Interdisciplinary Research and Consulting (CIRC). More information about CIRC can be found at www.umbc. edu/circ. This section focuses on the testing of basic operations using Matlab, Octave, FreeMat, Scilab, and R. We will first begin by solving a linear system; then finding eigenvalues and eigenvectors of a square matrix; and finally, 2-D functional plotting.

### 2.1 Basic operations in Matlab

This section discusses the results obtained using Matlab operations. To run Matlab on the cluster tara, enter matlab at the Linux command line. This starts up Matlab with its complete Java desktop interface. Useful options to Matlab on tara include -nodesktop, which starts only the command window within the shell, and -nodisplay, which disables all graphics output. For complete information on options, use matlab -h.

### 2.1.1 Solving Systems of Equations

The first example that we will consider in this section is solving a linear system. Consider the following system of equations:

$$
\begin{aligned}
-x_{2}+x_{3} & =3 \\
x_{1}-x_{2}-x_{3} & =0 \\
-x_{1}-x_{3} & =-3
\end{aligned}
$$

where the solution to this system is $(1,-1,2)^{T}$. In order to use Matlab to solve this system, let us express this linear system as a single matrix equation

$$
\begin{equation*}
A x=b \tag{2.1}
\end{equation*}
$$

where $A$ is a square matrix consisting of the coefficients of the unknowns, $x$ is the vector of unknowns, and $b$ is the right-hand side vector. For this particular system, we have

$$
A=\left[\begin{array}{ccc}
0 & -1 & 1 \\
1 & -1 & -1 \\
-1 & 0 & -1
\end{array}\right], \quad b=\left[\begin{array}{c}
3 \\
0 \\
-3
\end{array}\right] .
$$

To find a solution for this system in Matlab, left divide (2.1) by $A$ to obtain $x=A \backslash b$. Hence, Matlab use the backslash operator to solve this system. First, the matrix $A$ and vector $b$ are entered using the following:

```
A = [0 -1 1; 1 -1 -1; -1 0 -1]
b = [3;0;-3].
```

Now use $\mathrm{x}=\mathrm{A} \backslash \mathrm{b}$ to solve this system. The resulting vector which is assigned to $x$ is:
$\mathrm{x}=$
1
-1
2
Notice the solution is exactly what was expected based on our hand computations.

### 2.1.2 Calculating Eigenvalues and Eigenvectors

Here, we will consider another important function: computing eigenvalues and eigenvectors. Finding the eigenvalues and eigenvectors is a concept first introduced in a basic Linear Algebra course and we will begin by recalling the definition. Let $A \in \mathbb{C}^{n \times n}$ and $v \in \mathbb{C}^{n}$. A vector $v$ is called the eigenvector of $A$ if $v \neq 0$ and $A v$ is a multiple of $v$; that is, there exists a $\lambda \in \mathbb{C}$ such that

$$
A v=\lambda v
$$

where $\lambda$ is the eigenvalue of $A$ associated with the eigenvector $v$. We will use Matlab to compute the eigenvalues and a set of eigenvectors of a square matrix. Let us consider a matrix

$$
A=\left[\begin{array}{cc}
1 & -1 \\
1 & 1
\end{array}\right]
$$

which is a small matrix that we can easily compute the eigenvalues to check our results. Calculating the eigenvalues using $\operatorname{det}(A-\lambda I)=0$ gives $1+i$ and $1-i$. Now we will use Matlab's built in function eig to compute the eigenvalues. First enter the matrix $A$ and then calculate the eigenvalues using the following commands:
$\mathrm{A}=[1-1 ; 11] ;$
$\mathrm{v}=\operatorname{eig}(\mathrm{A})$
The following are the eigenvalues that are obtained for matrix $A$ using the commands stated above:

```
v =
\[
\begin{aligned}
& 1.0000+1.0000 \mathrm{i} \\
& 1.0000-1.0000 \mathrm{i}
\end{aligned}
\]
```

To check if the components of this vector are identical to the analytic eigenvalues, we can compute

```
v - [1+i;1-i]
```

and it results in
ans =

This demonstrates that the numerically computed eigenvalues have in fact the exact integer values for the real and imaginary parts, but Matlab formats the output for general real numbers.

In order to calculate the eigenvectors in Matlab, we will still use the eig function by slightly modifying it to $[\mathrm{P}, \mathrm{D}]=\operatorname{eig}(\mathrm{A})$ where $P$ will contain the eigenvectors of the square matrix $A$ and $D$ is the diagonal matrix containing the eigenvalues on its diagonals. In this case, the solution is:

```
P =
    0.7071 0.7071
        0-0.7071i 0 + 0.7071i
```

and
D =

```
1.0000 + 1.0000i 0
    0 1.0000 - 1.0000i
```

Calculating the eigenvector enables us to express the matrix $A$ as

$$
\begin{equation*}
A=P D P^{-1} \tag{2.2}
\end{equation*}
$$

where $P$ is the matrix of eigenvectors and D is a diagonal matrix as stated above. To check our solution, we will multiply the matrices generated using eig(A) to reproduce $A$ as suggested in (2.2).
$\mathrm{A}=\mathrm{P} * \mathrm{D} * \operatorname{inv}(\mathrm{P})$
produces
$A=$
1 -1
11
where $\operatorname{inv}(P)$ is used to obtain the inverse of matrix $P$. Notice that the commands above lead to the expected solution, $A$.

### 2.1.3 2-D Plotting

2-D plotting is a very important feature as it appears in all mathematical courses. Since this is a very commonly used feature, let us examine the 2-D plotting feature of Matlab by plotting $f(x)=x \sin \left(x^{2}\right)$ over the interval $[-2 \pi, 2 \pi]$. The data set for this function is given in a data file matlabdata.dat that is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications. Noticing that the data is given in two columns, we will first store the data in a matrix $A$. Second, we will create two vectors, $x$ and $y$, by extracting the data from the columns of $A$. Lastly, we will plot the data.


Figure 2.1: Plots of $f(x)=x \sin \left(x^{2}\right)$ in Matlab using (a) 129 and (b) 1025 equally spaced data points.
$\mathrm{A}=$ load ('matlabdata.dat');
$\mathrm{x}=\mathrm{A}(:, 1)$;
$y=A(:, 2)$;
plot( $x, y$ )
The commands stated above result in the Figure 2.1 (a). Looking at this figure, it can be noted that our axes are not labeled; there are no grid lines; and the peaks of the curves are rather coarse. The title, grid lines, and axes labels can be easily created. Let us begin by labeling the axes using xlabel('x') to label the $x$-axis and ylabel('f(x)') to label the $y$-axis. axis on and grid on can be used to create the axis and the grid lines. The axes are on by default and we can turn them off if necessary using axis off. Let us also create a title for this graph using title ('Graph of $f(x)=x \sin \left(x^{\wedge} 2\right)$ '). We have taken care of the missing annotations and lets try to improve the coarseness of the peaks in Figure 2.1 (a). We use length ( x ) to determine that 129 data points were used to create the graph of $f(x)$. To improve this outcome, we can begin by improving our resolution using
$\mathrm{x}=[-2 * \mathrm{pi}: 4 * \mathrm{pi} / 1024$ : $2 * \mathrm{pi}] ;$
to create a vector 1025 equally spaced data points over the interval $[-2 \pi, 2 \pi]$. In order to create vector $y$ consisting of corresponding $y$ values, use
$\mathrm{y}=\mathrm{x} . * \sin \left(\mathrm{x} .{ }^{\wedge} 2\right)$;
where .* performs element-wise multiplication and .^ corresponds to element-wise array power. Then, simply use plot $(x, y)$ to plot the data. Use the annotation techniques mentioned earlier to annotate the plot. In addition to the other annotations, use $x \lim ([-2 * \mathrm{pi} 2 * \mathrm{pi}])$ to set limit is for the x -axis. We can change the line width to 2 by plot (x,y, 'LineWidth', 2). Finally, Figure 2.1 (b) is the resulting figure with higher resolution as well as the annotations. Observe that by controlling the resolution in Figure 2.1 (b),
we have created a smoother plot of the function $f(x)$. The Matlab code used to create the annotated figure is as follows:

```
x = [-2*pi : 4*pi/1024 : 2*pi];
y = x.*sin(x. - 2);
H = plot(x,y);
set(H,'LineWidth',2)
axis on
grid on
title ('Graph of f(x)=x sin( ( ^ 2)
xlabel ('x')
ylabel ('f(x)')
xlim ([-2*pi 2*pi])
```


### 2.1.4 Programming

Here we will look at a basic example of Matlab programming using a script file. Let's try to plot Figure 2.1 (b) using a script file called plotxsinx.m. The extension .m indicates to Matlab that this is an executable m-file. Instead of typing multiple commands in Matlab, we will collect these commands into this script. The result is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications. Now, call plotxsinx (without the extension) on the command-line to execute it and create the plot with the annotations for $f(x)=x \sin \left(x^{2}\right)$. The plot obtained in this case is Figure 2.1 (b). This plot can be printed to a graphics file using the command

```
print -djpeg file_name_here.jpg
```


### 2.2 Basic operations in Octave

In this section, we will perform the basic operations on GNU Octave. To run Octave on the cluster tara, enter octave at the command line. For more information on Octave and available options, use man octave and octave -h.

Let us begin by solving a system of linear equations. Just like Matlab, Octave defines the backslash operator to solve equations of the form $A x=b$. Hence, the system of equations mentioned in Section 2.1.1 can also be solved in Octave using the same commands:

```
A = [0 -1 1; 1 -1 -1; -1 0 -1];
b = [3;0;-3];
x=A\b
```

which results in
$\mathrm{x}=$
1
-1
2

Clearly the solution is exactly what was expected. Hence, the process of solving the system of equations is identical to Matlab.

Now, let us consider the second operation of finding eigenvalues and eigenvectors. To find the eigenvalues and eigenvectors for matrix $A$ stated in Section 2.1.2, we will use Octave's built in function eig and obtain the following result:

```
v =
    1 + 1i
    1 - 1i
```

This shows exactly the integer values for the real and imaginary parts. To calculate the corresponding eigenvectors, use $[P, D]=\operatorname{eig}(A)$ and obtain:

```
P =
    0.70711 + 0.00000i 0.70711-0.00000i
    0.00000-0.70711i 0.00000 + 0.70711i
D =
    1+1i 0
        0 1 - 1i
```

After comparing this to the outcome generated by Matlab, we can conclude that the solutions are same but they are formatted slightly differently. For instance, matrix $P$ displays an extra decimal place when generated by Octave. The eigenvalues in Octave are reported exactly same as the calculated solution, where as Matlab displays them using four decimal places for real and imaginary parts. Hence, the solution is the same but presented slightly differently from each other. Before moving on, let us determine whether $A=P D P^{-1}$ still holds. Keeping in mind that the results were similar to Matlab's, we can expect this equation to hold true. Let us compute $P D P^{-1}$ by entering $\mathrm{P} * \mathrm{D} * \operatorname{inv}(\mathrm{P})$. Without much surprise, the outcome is

```
ans =
    1 -1
    1
```

An important thing to notice here is that to compute the inverse of a matrix, we use the inv command. Thus, the commands for computing the eigenvalues, eigenvectors, inverse of a matrix, as well as solving a linear system, are the same for Octave and Matlab.

Now, we will look at plotting $f(x)=x \sin \left(x^{2}\right)$ using the given data file. The load command is used to store the data in the file into a matrix $A$. use $\mathrm{x}=\mathrm{A}(:, 1)$ to store the first column as vector $x$ and $\mathrm{y}=\mathrm{A}(:, 2)$ to store the second column as vector $y$. We can create a plot using these vectors via entering $\operatorname{plot}(x, y)$ command in the prompt. Note that to check the number of data points, we can still use the length command. It is clear that this process is identical to the process in Section 2.1.3 that was used to generate Figure 2.1 (a). It would not be incorrect to assume that the figure generated in Octave could be identical to Figure 2.1 (a).


Figure 2.2: Plots of $f(x)=x \sin \left(x^{2}\right)$ in Octave using (a) 129 and (b) 1025 equally spaced data points.

Clearly, the Figure 2.2 (a) is not labeled at all; the grid is also not on; as well as the coarseness around the peaks exists. Therefore, the only difference between the two graphs is that in Figure 2.2 (a) the limits of the axes are different than in Figure 2.1 (a). The rest appears to be same in both of the plots. Let us try to label the axes of this figure using the label command and create the title using the title command. In order to create a smooth graph, like before; we will consider higher resolution. Hence, x = [-2*pi : 4*pi/1024 : 2*pi]; can be used to create a vector of 1025 points and $\mathrm{y}=\mathrm{x} . * \sin \left(\mathrm{x} .^{\wedge} 2\right)$; creates a vector of corresponding nfunctional values. By examining the creation of the $y$ vector, we notice that in Octave .* is known as the "element by element multiplication operator" and . ^ is the "element by element power operator." After using the label to label the axes; title to create a title; and grid on to turn on grid. We obtain Figure 2.2 (b).

Clearly, Figure 2.2 (b) and Figure 2.1 (b) are identical. We can simply put together all the commands in a script file exactly the way described in Section 2.1.4 and generate the Figure 2.1 (b). This results in the same m-file plotxsinx.m, which is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications. One additional command we can use to print the plot to a graphics file is

```
print -djpeg file_name_here.jpg
```


### 2.3 Basic operations in FreeMat

In this section, we perform the basic operations in FreeMat. To run FreeMat on tara, you have to load two modules first by entering module load qt/4.5.2 and module load freemat at the command line. Then start FreeMat by FreeMat. A useful option to Scilab on tara is -noX, which starts only the command window within the shell. For complete information on options, use matlab -help.

We will begin by first solving a linear system. Let us consider matrix $A$ as defined in Section 2.1.1. We can use the same commands a Matlab to produce a result.

```
A = [0 -1 1; 1 -1 -1; -1 0 -1];
b = [3;0;-3];
x = A\b
```

which results in
$\mathrm{x}=$
1
-1
2
as we had expected. Like Matlab and Octave, FreeMat also uses the backslash operator to solve linear systems.

Now, we will consider the second important operation, computing eigenvalues and eigenvectors. For our computations, let us use matrix $A$ stated in Section 2.1.2. We will use FreeMat's built in function eig and obtain the following result:

```
P =
    0.7071 + 0.0000i 0.7071-0.0000i
    0.0000-0.7071i 0.0000 + 0.7071i
D =
```



The outcome is identical to Matlab's results. Just to confirm, we compute $A=P D P^{-1}$ which results in the matrix $A$ as following:

```
ans =
    1.0000+0.0000i -1.0000+0.0000i
    1.0000+0.0000i 1.0000+0.0000i
```

A key point here is that FreeMat uses inv to compute inverse of matrices. So the commands used to solve systems of operations, calculate eigenvalues and eigenvectors, and computing matrix inverse are same as Matlab.

Now we would hope to see an agreement in the plotting and annotation commands. To examine the plotting feature of FreeMat, we will consider $f(x)=x \sin \left(x^{2}\right)$. Let us begin by examining the load command. Just like Matlab and Octave, we can load the data in a matrix $A$ with $\mathrm{A}=$ load('matlabdata.dat') command and use $\mathrm{x}=\mathrm{A}(:, 1)$ to create vector $x$ and $\mathrm{y}=\mathrm{A}(:, 2)$ to create $y$. Now, use plot $(\mathrm{x}, \mathrm{y})$ to generate Figure 2.3 (a) using vector $x$ and $y$. Clearly, the load command and plot command have same functionality as in Matlab.

Without much surprise, Figure 2.3 (a) and Figure 2.1 (a) are same. To annotate Figure 2.3 (a), we will use the same commands as Matlab. So to label the axes use label


Figure 2.3: Plots of $f(x)=x \sin \left(x^{2}\right)$ in FreeMat using (a) 129 and (b) 1025 equally spaced data points.
command, grid on create grid lines, title command to create title. To create a smooth graph, we will create another vector $x$ consisting of more equally spaced data points and a vector $y$ for the corresponding functional values. Use $\mathrm{x}=[-2 * \mathrm{pi}: 4 * \mathrm{pi} / 1024: 2 * \mathrm{pi}]$; to create $x$ and $\mathrm{y}=\mathrm{x} . * \sin \left(\mathrm{x} .{ }^{\wedge} 2\right)$; to create vector $y$. As in the earlier sections, we hope that higher resolution will improve our plot. Let us plot this data using plot ( $\mathrm{x}, \mathrm{y}$ ) ; . Applying the annotation techniques, we generate Figure 2.3 (b). Like in Matlab, we can also put together these commands in an m-file. This results in the same m-file plotxsinx.m, which is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications. Use
print('file_name_here.jpg'\}
to print the plot to a graphics file.

### 2.4 Basic operations in Scilab

In this section, we will perform the basic operations in Scilab. To run Scilab on tara, enter scilab at the command line, which opens its command window. A useful option to Scilab on tara is -nogui, which starts only the command window within the shell. For complete information on options, use matlab -h .

Once again, let us begin by solving the linear system from Section 2.1.1. Scilab follows the same method as GNU Octave and Matlab in solving the system of equations, i.e., it uses the backslash operator to find the solution using the system mentioned in Section 2.1.1, we use the following commands in Scilab:

```
A = [0 -1 1; 1 -1 -1; -1 0-1];
b = [3;0;-3];
x= A\b
```

to set up the matrix $A$ and vector $b$. Using the backslash operator, we obtain the result:


Figure 2.4: Plots of $f(x)=x \sin \left(x^{2}\right)$ in Scilab using (a) 129 and (b) 1025 equally spaced data points.

$$
\begin{aligned}
& \mathrm{x}= \\
& 1 . \\
&-1 . \\
& 2 .
\end{aligned}
$$

Once again, the result is exactly what is obtained when solving the system using an augmented matrix.

Now, let us determine how to calculate the eigenvalues and eigenvectors for the matrix $A$ stated in Section 2.1.2. Scilab uses the spec command which has the same functionality as eig command to compute eigenvalues. Hence, $v=\operatorname{spec}(A)$ results in $\mathrm{v}=$

1.     + i
2.     - i

Clearly, the outcome is exactly what we had expected but the outcome is formatted slightly different from Matlab. When we calculate the a set of corresponding eigenvectors using $[P, D]=\operatorname{spec}(A)$ and the following result is obtained:

D =

$P=$| $1+i$ | 0 |
| :---: | :---: | :--- |
| 0 | $1-1 i$ |
| 0.7071068 | 0.7071068 |
| $-0.7071068 i$ | $0.7071068 i$ |

By comparing $P$, the matrix of eigenvectors computed in Scilab, to $P$, the matrix in Section 2.1.2, we can see that both packages produce same results but they are formatted differently. Let us check our solution by computing $P D P^{-1}$ using the inv command to compute the inverse of the matrix.

```
P*D*inv(P)
ans =
    1. - 1.
    1. 1.
```

which is our initial matrix $A$. Note that one important factor in computing the eigenvalues and eigenvectors is the command used in these computations, spec, and that the eigenvectors found in Scilab and Matlab agree up to six decimal places.

Now, we will plot $f(x)=x \sin \left(x^{2}\right)$ in Scilab. To load the text file matlabdata.dat into a matrix, we use the Scilab command A = fscanfMat('matlabdata.dat'). This is specifically a command to read text files, while Scilab's load command is only for reading binary files; by contrast, Matlab uses load for both purposes. Then we use $\mathrm{x}=\mathrm{A}(:, 1)$ to store the first column vector as x and $\mathrm{y}=\mathrm{A}(:, 2)$ to store the second column as a vector $y$. We can create a plot using these vectors via entering plot ( $x, y$ ). Notice that the Figure 2.4 (a) is not labeled and it is rather coarse. Let us improve our resolution by creating vector $x$ using
$\mathrm{x}=[-2 * \% \mathrm{pi}: 4 * \% \mathrm{pi} / 1024: 2 * \%$ pi $]$
and let $\mathrm{y}=\mathrm{x} . * \sin \left(\mathrm{x} .{ }^{\wedge} 2\right)$ to create a corresponding $y$ vector. Unlike Matlab and Octave, we have to use \%pi to enter $\pi$ in Scilab. In addition, .* and .^ are still performing the element-wise operations called the "element-wise operators" in Scilab. Another factor that remains unchanged is the length command. We can generate the plot using the plot ( $\mathrm{x}, \mathrm{y}$ ) command which creates the Figure 2.4 (a). Once again, we can use xlabel and ylabel to label the axes; title('Graph of $f(x)=x \sin \left(x^{\wedge} 2\right)$ ') to create a title; and xgrid to turn on grid. To plot and create x -axis bounds, use

```
plot2D(x,y,1,'011','',[-2*%pi,y(1),2,*%pi,y($)])
```

Notice that we can put together these commands into a sci-file in Scilab to generate a plot. The resulting script for creating the a plot is as follows:

```
x = -2*%pi:(4*%pi)/1024:2*%pi;
y = x .* sin(x.^2);
plot2d(x,y,1,'011','',[-2*%pi,y(1),2*%pi,y($)])
set(gca(),"axes_visible","on")
set(gca(),"grid",[1,1])
title("Graph of f(x) = x sin(x^2)")
xlabel("x")
ylabel("f(x)")
```

Notice that some of the Matlab commands are not compatible with Scilab. One easier approach to handle this issue is to use the "Matlab to Scilab translator" under the Applications menu or by using mfile2sci command. The translator is unable to convert $x \lim ([-2 * \mathrm{pi} 2 * \mathrm{pi}])$; which we can take care of replacing the plot with plot2d command
stated earlier. The converted code using the translator is referred to as plotxsinx.sci and is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications. Using this script file, we obtain Figure 2.4 (b) which is similar to Figure 2.1 (b). To send these graphics to jpg file, we can use

```
xs2jpg(gcf(),'file_name_here.jpg')
```


### 2.5 Basic operations in $R$

In this section, we will perform the basic operations in $R$. To run $R$ on the cluster tara, enter $R$ at the Linux command line. This starts up the command window of $R$. For the complete list of options available with the command $R$, type $R$--help at the Linux command line. One useful command in R is help(), which provides a detailed description and usage of commands. For example, to view the usage of the command source in R, type help("source"). To quit the R interface, type quit(), which will ask if you wish to save the workspace you have been working on. Typing y will save all the commands you have typed in the session. These saved commands will be available next time you open R. Please refer to http://cran.r-project. org/manuals.html to learn about the R language.

Once again, let us begin by solving the linear system from Section 2.1.1. R has a command called solve to solve the linear system of equations $A x=b$. The command takes two matrices $A$ and $b$ as arguments and returns $x$ as a matrix. To solve the system mentioned in Section 2.1.1, we use the following commands in R:
$\mathrm{A}=\operatorname{array}(\mathrm{c}(0,1,-1,-1,-1,0,1,-1,-1), c(3,3))$
$b=c(3,0,-3)$
to set up the matrix $A$ and vector $b$, respectively. The c() function is prevalent in R code, and concatenates its arguments into a vector. The vector may be used to enumerate the elements of an array, as we have done here. In R, vectors are not the same as matrices. In order to instruct R to treat the above vector $b$ as a column vector, we use the following command:

```
dim(b) = c(3,1)
```

This command sets the dimension of the vector $b$ to 3 rows and 1 column. We have used the array () command to create the $A$ matrix. The first argument to this command is a vector containing the data and the second argument is again a vector containing the number of rows and columns of the matrix we want to create. When creating matrices, R follows the column-ordering scheme, that is, the entries in $c(0,1,-1,-1,-1,0,1,-1,-1)$ list the entries of $A$ along the columns; notice that this is different than the treatment of command-line input in Matlab and other packages. Using the solve command, we obtain the result

```
solve(A, b)
    [,1]
[1,] 1
[2,] -1
[3,] 2
```

Once again, the result is exactly what is obtained when solving the system using an augmented matrix.

In $R$, matrices can also be created using the matrix command. The above system can also be solved using the following commands:

```
A <- matrix(c(0,1,-1,-1,-1,0,1,-1,-1), nrow=3)
b <- matrix(c(3,0,-3))
solve(A, b)
    [,1]
[1,] 1
[2,] -1
[3,] 2
```

Although both the approaches work in this example, it is advisable to use the matrix command when working with matrices in R. Also, notice that we have used = assignment operator in the first method and <- operator in the second method. Both are assignment operators in R. However, = cannot be used to assign values inline unless the assignment is a named parameter. In other words, = can only be used as part of an independent top level expression or a sub-expression in a list of braced expressions. <-, on the other hand, can be used for inline assignments too, but has a potentially undesirable effect of either overwriting an existing variable with the same name in the workspace or creating a new variable with the same name if it does not exist ${ }^{5}$. The common practice is to use <- for assignments, $=$ for named parameters and to altogether avoid inline assignments. Since the features of R language is outside the scope of this report, we will not focus on exploring such subtle differences. Interested users should refer to R manuals for language specific details.

Now, let us determine how to calculate the eigenvalues and eigenvectors for the matrix $A$ stated in Section 2.1.2. R uses the eigen command to compute eigenvalues and eigenvectors. The command returns a data structure that holds both eigenvalues and eigenvectors. The return value can be captured into a variable as eig $=$ eigen(A). Eigenvalues can then be accessed as eig\$values and eigenvectors can be accessed as eig\$vectors. So, eigenvalues and eigenvectors can be calculated and displayed as

```
A <- matrix(c(1,1,-1,1),nrow=2)
eig <- eigen(A)
eig$values
[1] 1+1i 1-1i
eig$vectors
```

| $[$ [,1] | $[, 2]$ |
| ---: | :--- |

[1,] 0.7071068+0.0000000i 0.7071068+0.0000000i
[2,] 0.0000000-0.7071068i $0.0000000+0.7071068 i$

[^2]Clearly, the outcome is exactly what we had expected but the format of the outcome is slightly different from that of Matlab. Notice that Matlab's eig command returns both eigenvalues and eigenvectors in matrix form, with diagonal matrix $D$ and an invertible matrix $P$, where $D$ consists of eigenvalues as diagonal elements and $P$ consists of eigenvectors as columns. By contrast, R returns a heterogeneous list data structure encapsulating both eigenvalues and eigenvectors that can be accessed using the operator $\$$. It is common practice in R to return composite data structures from functions that need to return multiple values and to use the access operator $\$$ as shown above to access individual components in the data structure returned. But matrices $D$ and $P$ can be easily formed from these data structures, for example to verify that the diagonalization yields $A$ again:

```
D <- diag(eig$values)
P <- eig$vectors
P %*% D %*% solve(P)
    [,1] [,2]
[1,] 1+0i -1+0i
[2,] 1+0i 1+0i
```

Notice that the output of $P D P^{-1}$ is same as the matrix $A$. In the above command, $\%$ * $\%$ is used to perform matrix multiplications and the solve() command is used to calculate the inverse of the non-singular matrix $P$ (note that in R,the solve() command is also used to solve a linear system when called using two arguments $A$ and $b$, as shown at the beginning of this section).

Now, we will plot $f(x)=x \sin \left(x^{2}\right)$ in R . To load the text file matlabdata.dat into a matrix, we use the R command $\mathrm{pd}=$ read.table('matlabdata.dat'). There are several ways one can load data into $R$. Some of the commands provided by R are read.csv (to read comma-delimited files), scan, etc. R also has database/source specific commands such as read.dta (to load Stata binary files), read.octave (to load text files saved in the Octave format), etc. We use the command read.table to read rectangular data into an R variable. Rectangular data is read into R as data frames, which look like matrices and their elements can be accessed using [, and ,] operators. Below we plot the data using R's plot command.
$\mathrm{x}<-\operatorname{pd}[, 1]$
$\mathrm{y}<-\mathrm{pd}[, 2]$
plot(x, y, xlab="x", ylab="y")
dev.new()
dev.new() instructs $R$ to generate plots in seperate windows and not to overwrite the current plot with future plots (notice how we have used = operator for named parameters in the plot command). Notice that the Figure 2.5 (a) is not labeled and it is rather coarse. Also notice that by default, R uses circular markers instead of line markers as used in other packages. This behavior can be interpreted as R's distinguishing feature as a statistical package compared to other numerical packages. Let us improve the resolution by creating vector $x$ using


Figure 2.5: Plots of $f(x)=x \sin \left(x^{2}\right)$ in R using (a) 129 and (b) 1025 equally spaced data points.

```
x <- seq(-2*pi, 2*pi, 4*pi/1024)
```

and let $\mathrm{y}<-\mathrm{x} * \sin \left(\mathrm{x}^{\wedge} 2\right)$ to create a corresponding $y$ vector. Unlike Matlab and Octave, we have to use \%pi to enter $\pi$ in R. Unlike Matlab, Octave, and Scilab, we do not need to use .* and . ^ operators to perform element-wise operations. Instead, we simply use $*$ and -

We now plot using the new x and y values again using the plot command, but with additional arguments:

```
plot(x,y, type="l",xlab="x", ylab="f(x)", col="blue",
    main="Graph of f(x)=x sin(x^2)", lwd=2)
grid(nx=NULL,ny=NULL, lty=2, lwd=1, col="gray60", equilogs=TRUE)
```

The grid command above generates a fine grid against the plot. Note that R can handle line breaks in the code, as shown in the plot command above. However, users should ensure that the interpreter knows that the broken line is "to be continued". For example:

```
y<- 1 + 2 + 3 +
    4+5
z<- 1 + 2 + 3
    +4+5
```

would correctly assign the value of 15 to y , but would assign 6 to z which is probably not desired. Notice that the difference between the two assignment statements is that in the first command, the + operator appears at the end of the command, indicating to $R$ that you
intend to continue the command to the following line. The above plot command generates the Figure 2.5 (b) which is similar to Figure 2.1 (b). Notice that compared to Figure 2.5 (a), Figure 2.5 (b) shows solid line markers in the line plot. This is achieved with the argument type="l". To send these graphics to jpg file, we can use
savePlot("file_name_here.jpg", type = "jpeg")
We can save all the above commands to a text file with extension .R or .r. The file can then be executed in R, similar to Matlab's m-files.

## 3 Complex Operations Test

### 3.1 The Test Problem

Starting in this section, we study a classical test problem given by the numerical solution with finite differences for the Poisson problem with homogeneous Dirichlet boundary conditions [1, $4,5,12$ ], given as

$$
\begin{align*}
-\triangle u=f & \text { in } \Omega, \\
u=0 & \text { on } \partial \Omega . \tag{3.1}
\end{align*}
$$

This problem was studied before in, among other sources, $[1,4,7,8,10,12]$. Here $\partial \Omega$ denotes the boundary of the domain $\Omega$ while the Laplace operator is defined as

$$
\Delta u=\frac{\partial^{2} u}{\partial x^{2}}+\frac{\partial^{2} u}{\partial y^{2}}
$$

This partial differential equation can be used to model heat flow, fluid flow, elasticity, and other phenomena [12]. Since $u=0$ at the boundary in (3.1), we are looking at a homogeneous Dirichlet boundary condition. We consider the problem on the two-dimensional unit square $\Omega=(0,1) \times(0,1) \subset \mathbb{R}^{2}$. Thus, (3.1) can be restated as

$$
\begin{array}{ll}
-\frac{\partial^{2} u}{\partial x^{2}}-\frac{\partial^{2} u}{\partial y^{2}}=f(x, y) & \text { for } 0<x<1, \quad 0<y<1  \tag{3.2}\\
u(0, y)=u(x, 0)=u(1, y)=u(x, 1)=0 & \text { for } 0<x<1, \quad 0<y<1
\end{array}
$$

where the function $f$ is given by

$$
f(x, y)=-2 \pi^{2} \cos (2 \pi x) \sin ^{2}(\pi y)-2 \pi^{2} \sin ^{2}(\pi x) \cos (2 \pi y)
$$

The problem is designed to admit a closed-form solution as the true solution

$$
u(x, y)=\sin ^{2}(\pi x) \sin ^{2}(\pi y)
$$

### 3.2 Finite Difference Discretization

Let us define a grid of mesh points $\Omega_{h}=\left(x_{i}, y_{j}\right)$ with $x_{i}=i h, i=0, \ldots, N+1, y_{j}=j h, j=$ $0, \ldots, N+1$ where $h=\frac{1}{N+1}$. By applying the second-order finite difference approximation to the $x$-derivative at all the interior points of $\Omega_{h}$, we obtain

$$
\begin{equation*}
\frac{\partial^{2} u}{\partial x^{2}}\left(x_{i}, y_{i}\right) \approx \frac{u\left(x_{i-1}, y_{j}\right)-2 u\left(x_{i}, y_{j}\right)+u\left(x_{i+1}, y_{j}\right)}{h^{2}} \tag{3.3}
\end{equation*}
$$

If we also apply this to the $y$-derivative, we obtain

$$
\begin{equation*}
\frac{\partial^{2} u}{\partial y^{2}}\left(x_{i}, y_{i}\right) \approx \frac{u\left(x_{i}, y_{j-1}\right)-2 u\left(x_{i}, y_{j}\right)+u\left(x_{i}, y_{j+1}\right)}{h^{2}} \tag{3.4}
\end{equation*}
$$

Now, we can apply (3.3) and (3.4) to (3.2) and obtain

$$
\begin{align*}
& -\frac{u\left(x_{i-1}, y_{j}\right)-2 u\left(x_{i}, y_{j}\right)+u\left(x_{i+1}, y_{j}\right)}{h^{2}}  \tag{3.5}\\
& -\frac{u\left(x_{i}, y_{j-1}\right)-2 u\left(x_{i}, y_{j}\right)+u\left(x_{i}, y_{j+1}\right)}{h^{2}} \approx f\left(x_{i}, y_{j}\right) .
\end{align*}
$$

Hence, we are working with the following equations for the approximation $u_{i, j} \approx u\left(x_{i}, y_{j}\right)$ :

$$
\begin{array}{r}
-u_{i-1, j}-u_{i, j-1}+4 u_{i, j}-u_{i+1, j}-u_{i, j+1}=h^{2} f_{i, j} \quad i, j=1, \ldots, N  \tag{3.6}\\
u_{0, j}=u_{i, 0}=u_{N+1, j}=u_{i, N+1}=0
\end{array}
$$

The equations in (3.6) can be organized into a linear system $A u=b$ of $N^{2}$ equations for the approximations $u_{i, j}$. Since we are given the boundary values, we can conclude there are exactly $N^{2}$ unknowns. In this linear system, we have

$$
A=\left[\begin{array}{ccccc}
S & -I & & & \\
-I & S & -I & & \\
& \ddots & \ddots & \ddots & \\
& & -I & S & -I \\
& & & -I & S
\end{array}\right] \in \mathbb{R}^{N^{2} \times N^{2}}
$$

where

$$
S=\left[\begin{array}{cccccc}
4 & -1 & & & & \\
-1 & 4 & -1 & & \\
& \ddots & \ddots & \ddots & \\
& & -1 & 4 & -1 \\
& & & & -1 & 4
\end{array}\right] \in \mathbb{R}^{N \times N} \text { and } I=\left[\begin{array}{cccc}
1 & & & \\
& 1 & & \\
& & & \\
& & \ddots & \\
& & & 1 \\
& & & \\
& & &
\end{array}\right] \in \mathbb{R}^{N \times N}
$$

and the right-hand side vector $b_{k}=h^{2} f_{i, j}$ where $k=i+(j-1) N$. The matrix $A$ is symmetric and positive definite $[4,12]$. This implies that the linear system has a unique solution and it guarantees that the iterative conjugate gradient method converges.

To create the matrix $A$ in Matlab effectively, we use the observation that it is given by a sum of two Kronecker products [4, Section 6.3.3]: Namely, $A$ can be interpreted as the sum

$$
A=\left[\begin{array}{lllll}
T & & & & \\
& T & & & \\
& & \ddots & & \\
& & & T & \\
& & & & T
\end{array}\right]+\left[\begin{array}{ccccc}
2 I & -I & & & \\
-I & 2 I & -I & & \\
& \ddots & \ddots & \ddots & \\
& & -I & 2 I & -I \\
& & & -I & 2 I
\end{array}\right] \in \mathbb{R}^{N^{2} \times N^{2}},
$$

where

$$
T=\left[\begin{array}{ccccc}
2 & -1 & & & \\
-1 & 2 & -1 & & \\
& \ddots & \ddots & \ddots & \\
& & -1 & 2 & -1 \\
& & & -1 & 2
\end{array}\right] \in \mathbb{R}^{N \times N}
$$

and $I$ is the $N \times N$ identity matrix, and each of the matrices in the sum can be computed by Kronecker products involving $T$ and $I$, so that $A=I \otimes T+T \otimes I$. These formulas are the basis for the code in the function setupA shown in driver_ge.m, which is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications.

One of the things to consider to confirm the convergence of the finite difference method is the finite difference error. The finite difference error is defined as the difference between the true solution $u(x, y)$ and the numerical solution $u_{h}$ defined on the mesh points by $u_{h}\left(x_{i}, y_{j}\right)=$ $u_{i, j}$. Since the solution $u$ is sufficiently smooth, we expect the finite difference error to decrease as $N$ gets larger and $h=\frac{1}{N+1}$ gets smaller. Specifically, the finite difference theory predicts that the error will converge as $\left\|u-u_{h}\right\|_{L^{\infty}(\Omega)} \leq C h^{2}$, as $h \rightarrow 0$, where $C$ is a constant independent of $h[2,5]$. For sufficiently small $h$, we can then expect that the ratio of errors on consecutively refined meshes behaves like

$$
\begin{equation*}
\text { Ratio }=\frac{\left\|u-u_{2 h}\right\|}{\left\|u-u_{h}\right\|} \approx \frac{C(2 h)^{2}}{C h^{2}}=4 \tag{3.7}
\end{equation*}
$$

Thus, we will print this ratio in the following tables in order to confirm convergence of the finite difference method.

### 3.3 Matlab Results

### 3.3.1 Gaussian Elimination

Let us begin solving the linear system arising from the Poisson problem by Gaussian elimination in Matlab. We know that this is easiest approach for solving linear systems for the user of Matlab, although it may not necessarily be the best method for large systems. To create matrix $A$, we make use of the Kronecker tensor product, as described in Section 3.2. This can be easily implemented in Matlab using the kron function. The system is then solved using the backslash operator. Figure 3.1 shows the results of this for a mesh with $N=32$. Figure 3.1 (a) shows the mesh plot of the numerical solution vs. $(x, y)$. The error at each mesh point is computed by subtracting the numerical solution from the analytical solution and is plotted in Figure 3.1 (b). Notice that the maximum error occurs at the center. The code to solve this system for $N=32$ and produce the plots is contained in driver_ge.m, which is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications.

Table 3.1 (a) shows the results of a study for this problem using Gaussian elimination with mesh resolutions $N=2^{\nu}$ for $\nu=1,2,3, \ldots, 13$. The table lists the mesh resolution $N$, the number of degrees of freedom (DOF) $N^{2}$, the norm of the finite difference error $\left\|u-u_{h}\right\|$, the ratio of consecutive error norms (3.7), and the observed wall clock time in HH:MM:SS. To create this table, we use a version of driver_ge.m with the graphics commands commented out.

The norms of the finite difference errors clearly go to zero as the mesh resolution $N$ increases. The ratios between error norms for consecutive rows in the table tend to 4 , which confirms that the finite difference method for this problem is second-order convergent with errors behaving like $h^{2}$, as predicted by the finite difference theory. By looking at this table,


Figure 3.1: Mesh plots for $N=32$ in Matlab (a) of the numerical solution and (b) of the numerical error.
it can be concluded that Gaussian elimination runs out of memory for $N=8,192$. Hence, we are unable to solve this problem for $N$ larger than 4,096 via Gaussian elimination. This leads to the need for another method to solve larger systems. Thus, we will use an iterative method known as the conjugate gradient method to solve this linear system.

### 3.3.2 Conjugate Gradient Method

Now, we use the conjugate gradient method to solve the Poisson problem [1, 12]. This iterative method is an alternative to using Gaussian elimination to solve a linear system and is accomplished by replacing the backslash operator by a call to the pcg function. We use the zero vector as the initial guess and a tolerance of $10^{-6}$ on the relative residual of the iterates. This is implemented in a Matlab code driver_cg.m that is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications. This code is equivalent to driver_ge.m from above, with the backslash operator replaced by a call to the pcg function.

The system matrix $A$ accounts for the largest amount of memory used by the conjugate gradient method. Hence, to solve the problem for larger meshes, we can use a so-called matrix-free implementation of the method that avoids storing $A$. The only place, where $A$ enters into the iterative method is the matrix-vector multiplication $q=A p$ in each iteration. Hence, for a matrix-free implementation, instead of supplying $A$ itself, the user supplies a function that computes the matrix-vector product $q=A p$ directly for a given input vector $p$ without ever storing a system matrix $A$. Thus, the matrix-free implementation returns a vector $q$ for the input vector $p$ by performing the component-wise product on $p$ and the matrix $A$ using the knowledge of the components of $A$. Thus, we can expect to save memory significantly. This is implemented in the funtion Ax at the bottom of the code driver_cg.m.

Table 3.1 (b) shows results of a study using the conjugate gradient method with this matrix-free implementation of the matrix-vector multiplication. The column \#iter lists the number of iterations taken by the iteration method to converge. The finite difference error

Table 3.1: Convergence results for the test problem in Matlab using (a) Gaussian elimination and (b) the conjugate gradient method. The tables list the mesh resolution $N$, the number of degrees of freedom (DOF), the finite difference norm $\left\|u-u_{h}\right\|_{L^{\infty}(\Omega)}$, the ratio of consecutive errors, and the observed wall clock time in HH:MM:SS.

| (a) Gaussian Elimination |  |  |  |  |
| ---: | ---: | ---: | ---: | ---: |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | Time |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | N/A | $<00: 00: 01$ |
| 64 | 4,096 | $7.7812 \mathrm{e}-4$ | 3.8719 | $<00: 00: 01$ |
| 128 | 16,384 | $1.9766 \mathrm{e}-4$ | 3.9366 | $<00: 00: 01$ |
| 256 | 65,536 | $4.9807 \mathrm{e}-5$ | 3.9685 | $<00: 00: 01$ |
| 512 | 262,144 | $1.2501 \mathrm{e}-5$ | 3.9843 | $00: 00: 01$ |
| 1,024 | $1,048,576$ | $3.1313 \mathrm{e}-6$ | 3.9923 | $00: 00: 31$ |
| 2,048 | $4,194,304$ | $7.8362 \mathrm{e}-7$ | 3.9959 | $00: 00: 27$ |
| 4,096 | $16,777,216$ | $1.9610 \mathrm{e}-7$ | 3.9960 | $00: 02: 07$ |
| 8,192 |  | out of memory |  |  |


| (b) Conjugate Gradient Method |  |  |  |  |  |
| ---: | ---: | ---: | ---: | ---: | ---: |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | \#iter | Time |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | N/A | 48 | $<00: 00: 01$ |
| 64 | 4,096 | $7.7811 \mathrm{e}-4$ | 3.8719 | 96 | $<00: 00: 01$ |
| 128 | 16,384 | $1.9765 \mathrm{e}-4$ | 3.9368 | 192 | $<00: 00: 01$ |
| 256 | 65,536 | $4.9797 \mathrm{e}-5$ | 3.9690 | 387 | $00: 00: 02$ |
| 512 | 262,144 | $1.2494 \mathrm{e}-5$ | 3.9856 | 783 | $00: 00: 40$ |
| 1,024 | $1,048,576$ | $3.1266 \mathrm{e}-6$ | 3.9961 | 1,581 | $00: 05: 47$ |
| 2,048 | $4,194,304$ | $7.8019 \mathrm{e}-7$ | 4.0075 | 3,192 | $00: 50: 26$ |
| 4,096 | $16,777,216$ | $1.9353 \mathrm{e}-7$ | 4.0313 | 6,452 | $07: 46: 24$ |
| 8,192 | $67,108,864$ | $4.7400 \mathrm{e}-8$ | 4.0829 | 13,033 | $69: 42: 45$ |

shows the same behavior as in Table 3.1 (a) with ratios of consecutive errors approaching 4 as for Gaussian elimination; this confirms that the tolerance on the relative residual of the iterates is tight enough. To create this table, we use a version of driver_cg.m with the graphics commands commented out.

Tables 3.1 (a) and (b) indicate that Gaussian elimination is faster than the conjugate gradient method in Matlab, whenever it does not run out of memory. For problems greater than 4,096 , the results show that the conjugate gradient method is able to solve for larger mesh resolutions.


Figure 3.2: Mesh plots for $N=32$ in Octave (a) of the numerical solution and (b) of the numerical error.

### 3.4 Octave Results

### 3.4.1 Gaussian Elimination

In this section, we will solve the Poisson problem discussed in Section 3.2 via Gaussian elimination method in Octave. Just like Matlab, we can solve the equation using the backslash operator. Using the same m-file we used in Matlab, driver_ge.m,which is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications, we create Figure 3.2 which is identical to Figure 3.1.

The numerical results in Table 3.2 (a) are identical to the results in the Table 3.1 (a), but the timing results show that Matlab is significantly faster than Octave when solving a system of linear equations using the method of Gaussian elimination.

### 3.4.2 Conjugate Gradient Method

Now, let us try to solve the problem using the conjugate gradient method in Octave. Just like Matlab, there exists a pcg.m function. The input requirements for the pcg function are identical to the Matlab pcg function. Once again, we use the zero vector as the initial guess and the tolerance is $10^{-6}$ on the relative residual of the iterations. We can use the m -file driver_cg.m which, is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications.

Just like in Matlab, Tables 3.2 (a) and (b) indicate that Gaussian elimination is faster than the conjugate gradient method in Octave, whenever it does not run out of memory. For problems greater than 4,096 , the results show that the conjugate gradient method is able to solve for larger mesh resolutions.

Comparing Table 3.2 (b) to Table 3.1 (b), we see that the conjugate gradient method in Octave solves a problem faster than Matlab.

Table 3.2: Convergence results for the test problem in Octave using (a) Gaussian elimination and (b) the conjugate gradient method. The tables list the mesh resolution $N$, the number of degrees of freedom (DOF), the finite difference norm $\left\|u-u_{h}\right\|_{L^{\infty}(\Omega)}$, the ratio of consecutive errors, and the observed wall clock time in HH:MM:SS.

| (a) Gaussian Elimination |  |  |  |  |
| ---: | ---: | ---: | ---: | ---: |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | Time |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | N/A | $<00: 00: 01$ |
| 64 | 4,096 | $7.7812 \mathrm{e}-4$ | 3.8719 | $<00: 00: 01$ |
| 128 | 16,384 | $1.9766 \mathrm{e}-4$ | 3.9366 | $<00: 00: 01$ |
| 256 | 65,536 | $4.9807 \mathrm{e}-5$ | 3.9685 | $<00: 00: 01$ |
| 512 | 262,144 | $1.2501 \mathrm{e}-5$ | 3.9843 | $00: 00: 02$ |
| 1,024 | $1,048,576$ | $3.1313 \mathrm{e}-6$ | 3.9922 | $00: 00: 16$ |
| 2,048 | $4,194,304$ | $7.8362 \mathrm{e}-7$ | 3.9959 | $00: 01: 57$ |
| 4,096 | $16,777,216$ | $1.9610 \mathrm{e}-7$ | 3.9960 | $00: 15: 37$ |
| 8,192 |  | out of memory |  |  |


| (b) Conjugate gradient method |  |  |  |  |  |
| ---: | ---: | ---: | ---: | ---: | ---: |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | \#iter | Time |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | N/A | 48 | $<00: 00: 01$ |
| 64 | 4,096 | $7.7811 \mathrm{e}-4$ | 3.8719 | 96 | $<00: 00: 01$ |
| 128 | 16,384 | $1.9765 \mathrm{e}-4$ | 3.9368 | 192 | $<00: 00: 01$ |
| 256 | 65,536 | $4.9797 \mathrm{e}-5$ | 3.9690 | 387 | $00: 00: 03$ |
| 512 | 262,144 | $1.2494 \mathrm{e}-5$ | 3.9856 | 783 | $00: 00: 27$ |
| 1,024 | $1,048,576$ | $3.1266 \mathrm{e}-6$ | 3.9961 | 1,581 | $00: 04: 23$ |
| 2,048 | $4,194,304$ | $7.8019 \mathrm{e}-7$ | 4.0075 | 3,192 | $00: 40: 07$ |
| 4,096 | $16,777,216$ | $1.9353 \mathrm{e}-7$ | 4.0313 | 6,452 | $05: 52: 11$ |
| 8,192 | $67,108,864$ | $4.6797 \mathrm{e}-8$ | 4.1355 | 13,033 | $40: 44: 32$ |

### 3.5 FreeMat Results

### 3.5.1 Gaussian Elimination

Once again, we will solve the Poisson problem via Gaussian elimination method, but using FreeMat this time. To create the system matrix $A$ using setupA, we need to write our own implementation of the kron function, since it is not implemeneted in FreeMat. Furthermore, mesh does not exist in FreeMat, so we used the plot3 command to create Figure 3.3 (a) the plot3 of the numerical solution vs. $(x, y)$ and Figure $3.3(\mathrm{~b})$ the plot3 of the error vs. $(x, y)$. Despite the view adjustments for these figures using the view command, Figure 3.3 (b) is still very hard to study. The code to solve this system for $N=32$ and produce the plots is contained in freemat_ge.m, which is posted along with the tech. report [3] at www.umbc. edu/hpcf under Publications.


Figure 3.3: Plot3 plots for $N=32$ in FreeMat (a) of the numerical solution and (b) of the numerical error.

The numerical results in Table 3.3 (a) are identical to the results in Tables 3.1 (a) and 3.2 (a) for the mesh resolutions that FreeMat can solve for. However, Gaussian elimination in FreeMat ran significantly slower than Octave, which in turn ran slower than Matlab, in those cases. Moreover, Gaussian elimination in FreeMat could only solve the problem up to $N=2,048$ and ran out of memory already for $N=4,096$.

### 3.5.2 Conjugate Gradient Method

Unlike Matlab and Octave, a pcg function does not exist in FreeMat. To address this issue, we wrote our own cg function to use in FreeMat. The code to solve this system for $N=32$ is contained in freemat_cg.m, which is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications. To create the table of results for larger $N$, we used a version of this function with graphics commands commented out.

We encountered some issues in FreeMat when we were solving for a mesh resolution of $4,096 \times 4,096$, namely the time it took to solve the problem was excessively long. We predicted that the code would run for approximatley 195 hours or 8 days. We determined this by having the code display the error every 100 iterations. After fifteen hours of running the code only 500 iterations had been completed. Using this knowledge we calculated that an iteration occured every 1.8 minutes. So to solve the system for a mesh resolution of $4,096 \times 4,096$ it would take approximately eight days to compelete the calculations. Due to the length of time needed to complete the calculations we did not collect the all of the results, but because the error was printing in the correct trend and appearing every 100 iterations we were able to guarantee that the code was working correctly and the computer was not running out of memory.

The numerical results in Table 3.3 (b) are identical to the results in Tables 3.1 (b) and 3.2 (b) for the mesh resolutions that FreeMat can solve for. However, the conjugate gradient method in FreeMat ran significantly slower than Octave and Matlab in those cases. As explained in the previous paragraph, simulations with $N>2,048$ would take an excessive amount of time, and we did not run them to completion.

Table 3.3: Convergence results for the test problem in FreeMat using Gaussian elimination. The tables list the mesh resolution $N$, the number of degrees of freedom (DOF), the finite difference norm $\left\|u-u_{h}\right\|_{L^{\infty}(\Omega)}$, the ratio of consecutive errors, and the observed wall clock time in HH:MM:SS.

| (a) Gaussian Elimination |  |  |  |  |  |
| ---: | ---: | ---: | ---: | ---: | ---: |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | Time |  |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | N/A | $<00: 00: 01$ |  |
| 64 | 4,096 | $7.7812 \mathrm{e}-4$ | 3.8719 | $<00: 00: 01$ |  |
| 128 | 16,384 | $1.9766 \mathrm{e}-4$ | 3.9366 | $<00: 00: 01$ |  |
| 256 | 65,536 | $4.9807 \mathrm{e}-5$ | 3.9685 | $00: 00: 04$ |  |
| 512 | 262,144 | $1.2501 \mathrm{e}-5$ | 3.9843 | $00: 00: 28$ |  |
| 1,024 | $1,048,576$ | $3.1313 \mathrm{e}-6$ | 3.9922 | $00: 03: 15$ |  |
| 2,048 | $4,194,304$ | $7.8362 \mathrm{e}-7$ | 3.9959 | $00: 14: 29$ |  |
| 4,096 | out of memory |  |  |  |  |
| 8,192 | out of memory |  |  |  |  |
| Conjugate Gradient Method |  |  |  |  |  |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | $\#$ iter | Time |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | N/A | 48 | $<00: 00: 01$ |
| 64 | 4,096 | $7.7810 \mathrm{e}-4$ | 3.8719 | 96 | $00: 00: 03$ |
| 128 | 16,384 | $1.9765 \mathrm{e}-4$ | 3.9368 | 192 | $00: 00: 23$ |
| 256 | 65,536 | $4.9797 \mathrm{e}-5$ | 3.9689 | 387 | $00: 03: 07$ |
| 512 | 262,144 | $1.2494 \mathrm{e}-5$ | 3.9856 | 783 | $00: 15: 21$ |
| 1,024 | $1,048,576$ | $3.1266 \mathrm{e}-6$ | 3.9961 | 1,581 | $03: 24: 25$ |
| 2,048 | $4,194,304$ | $7.8019 \mathrm{e}-7$ | 4.0075 | 3,192 | $03: 52: 24$ |
| 4,096 | excessive time requirement |  |  |  |  |
| 8,192 | excessive time requirement |  |  |  |  |

### 3.6 Scilab Results

### 3.6.1 Gaussian Elimination

Once again, we will solve the Poisson equation via Gaussian elimination, this time using Scilab. Scilab also uses the backslash operator to solve the linear system. To compute the Kronecker tensor product of matrix $X$ and $Y$ in Scilab, we can use the X.*.Y command when setting up the system matrix $A$. Figure 3.4 (a) is a mesh plot of numerical solution for a mesh resolution of $N=32$, and Figure 3.4 (b) is a plot of the error associated with the numerical solution. The mesh plots are equivalent to the Matlab mesh plots in Figure 3.1. The code to solve this system for $N=32$ and produce the plots is contained in driver_ge.sci, which is posted along with the tech. report [3] at www.umbc.edu/hpcf under Publications. The initial version of this Scilab sci-file is obtained using the "Matlab to Scilab translator" under


Figure 3.4: Mesh plots for $N=32$ in Scilab (a) of the numerical solution and (b) of the numerical error.
the Applications menu or by using mfile2sci command in Scilab.
To create the tables in Scilab we used a version driver_ge.sci with the graphics commands commented out. The Scilab code used to create the tables also utilizes the Scilab command stacksize("max"). The stacksize("max") command allows Scilab to use all available memory when running the code. The numerical results in Table 3.4 (a) are identical to the results in Tables 3.1 (a) and 3.2 (a) for the mesh resolutions that Scilab can solve for. However, Gaussian elimination in Scilab ran significantly slower than Octave, which in turn ran slower than Matlab, in those cases. Moreover, Gaussian elimination in Scilab can only solve the problem up to $N=1,024$ and runs out of memory already for $N=2,048$, despite the use of stacksize("max"). In fact, without stacksize("max"), Scilab can only solve the problem up to $N=256$ and runs out of memory already for $N=512$.

We attempted to see if we could get better results by switching to UMFPack or TAUCS as alternative linear solvers available in Scilab. These solvers can be accessed by replacing the backslash operator in the linear solve. This is the only difference in the files driver_UMFPack_solver.sci driver_TAUCS_solver.sci, respectively, compared to the driver_ge.sci. These files are posted along with the tech. report. The results for UMFPack and TAUCS in Tables 3.4 (c) and (d), respectively, show that neither UMFPack nor TAUCS were able to solve for $N=2,048$ either, but they were both much faster than the default backslash operator that was used in Table 3.4 (a). At this point, we hope to solve a larger system using the conjugate gradient method.

### 3.6.2 Conjugate Gradient Method

Let us use the conjugate gradient method to solve the Poisson problem in Scilab. Here, we will use Scilab's pcg function. In order to solve, the initial guess is the zero vector and the tolerance is $10^{-6}$ on the relative residual of the iterates. We can use the sci-file driver_cg.sci which is posted along with the tech. report [3] at www.umbc.edu/hpcf under

Table 3.4: Convergence results for the test problem in Scilab using (a) Gaussian elimination and (b) the conjugate gradient method. The tables list the mesh resolution $N$, the number of degrees of freedom (DOF), the finite difference norm $\left\|u-u_{h}\right\|_{L^{\infty}(\Omega)}$, the ratio of consecutive errors, and the observed wall clock time in HH:MM:SS.

| (a) Gaussian elimination |  |  |  |  |
| ---: | ---: | ---: | ---: | ---: |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | Time |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | N/A | $<00: 00: 01$ |
| 64 | 4,096 | $7.7812 \mathrm{e}-4$ | 3.8719 | $<00: 00: 01$ |
| 128 | 16,384 | $1.9766 \mathrm{e}-4$ | 3.9366 | $00: 00: 11$ |
| 256 | 65,536 | $4.9807 \mathrm{e}-5$ | 3.9685 | $00: 03: 19$ |
| 512 | 262,144 | $1.2500 \mathrm{e}-5$ | 3.9846 | $00: 39: 04$ |
| 1,024 | $1,048,576$ | $3.1313 \mathrm{e}-6$ | 3.9920 | $08: 32: 20$ |
| 2,048 |  | out of memory |  |  |
| 4,096 |  | out of memory |  |  |
| 8,192 |  | out of memory |  |  |


| (b) Conjugate gradient method |  |  |  |  |  |
| ---: | ---: | ---: | ---: | ---: | ---: |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | \#iter | Time |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | $\mathrm{~N} / \mathrm{A}$ | 48 | $<00: 00: 01$ |
| 64 | 4,096 | $7.7811 \mathrm{e}-4$ | 3.8719 | 96 | $<00: 00: 01$ |
| 128 | 16,384 | $1.9765 \mathrm{e}-4$ | 3.9368 | 192 | $<00: 00: 01$ |
| 256 | 65,536 | $4.9797 \mathrm{e}-5$ | 3.9690 | 387 | $00: 00: 04$ |
| 512 | 262,144 | $1.2494 \mathrm{e}-5$ | 3.9856 | 783 | $00: 00: 36$ |
| 1,024 | $1,048,576$ | $3.1266 \mathrm{e}-6$ | 3.9960 | 1,581 | $00: 06: 57$ |
| 2,048 | $4,194,304$ | $7.8018 \mathrm{e}-7$ | 4.0075 | 3,192 | $00: 48: 12$ |
| 4,096 | $16,777,216$ | $1.9353 \mathrm{e}-7$ | 4.0313 | 6,452 | $06: 03: 43$ |
| 8,192 |  | out of memory |  |  |  |


| (c) UMFPack Gaussian elimnation |  |  |  |  |
| ---: | ---: | ---: | ---: | ---: |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | Time |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | N/A | $<00: 00: 01$ |
| 64 | 4,096 | $7.7812 \mathrm{e}-4$ | 3.8719 | $<00: 00: 01$ |
| 128 | 16,384 | $1.9766 \mathrm{e}-4$ | 3.9366 | $<00: 00: 01$ |
| 256 | 65,536 | $4.9807 \mathrm{e}-5$ | 3.9685 | $<00: 00: 01$ |
| 512 | 262,144 | $1.2500 \mathrm{e}-5$ | 3.9846 | $00: 00: 03$ |
| 1,024 | $1,048,576$ | $3.1313 \mathrm{e}-6$ | 3.9920 | $00: 00: 27$ |
| 2,048 |  | out of memory |  |  |
| 4,096 |  | out of memory |  |  |
| 8,192 |  | out of memory |  |  |


| (d) TAUCS Gaussian elimination |  |  |  |  |
| ---: | ---: | ---: | ---: | ---: |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | Time |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | N $/ \mathrm{A}$ | $<00: 00: 01$ |
| 64 | 4,096 | $7.7812 \mathrm{e}-4$ | 3.8719 | $<00: 00: 01$ |
| 128 | 16,384 | $1.9766 \mathrm{e}-4$ | 3.9366 | $<00: 00: 01$ |
| 256 | 65,536 | $4.9807 \mathrm{e}-5$ | 3.9685 | $<00: 00: 01$ |
| 512 | 262,144 | $1.2500 \mathrm{e}-5$ | 3.9846 | $00: 00: 01$ |
| 1,024 | $1,048,576$ | $3.1313 \mathrm{e}-6$ | 3.9920 | $00: 00: 09$ |
| 2,048 |  | out of memory |  |  |
| 4,096 |  | out of memory |  |  |
| 8,192 |  | out of memory |  |  |

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The numerical results in Table 3.4 (b) are identical to the results in Tables 3.1 (b) and 3.2 (b) for the mesh resolutions that Scilab can solve for. Also the run times are comparable for these cases. However, the conjugate gradient method in Scilab ran out of memory for $N=8,192$, despite the use of stacksize("max").


Figure 3.5: Mesh plots for $N=32$ in R (a) of the numerical solution and (b) of the numerical error.

### 3.7 R Results

### 3.7.1 Gaussian Elimination

Once again, we will solve the Poisson equation via Gaussian elimination, this time using R. As mentioned earlier, $R$ uses the solve command to solve the linear system. To compute the Kronecker tensor product of matrix $X$ and $Y$ in R, we have used the kronecker() command while setting up the system matrix $A$. Figure 3.5 (a) is a mesh plot of numerical solution for a mesh resolution of $N=32$, and Figure $3.5(\mathrm{~b})$ is a plot of the error associated with the numerical solution. The mesh plots are equivalent to the Matlab mesh plots in Figure 3.1.

To create the tables in $R$, we have used a version of the driver script driver_ge.r with the graphics commands commented out. Comparing the results in Table 3.5 (a) with Tables 3.1 (a), 3.2 (a), 3.3 (a), and 3.4 (a), one can say that R's performance was similar to Freemat's and better than Scilab's but worse than Matlab's and Octave's performances. Also, R could not solve for the resolutions $N=2,048,4,096$, and 8,192 , as it ran out of allocated memory.

### 3.7.2 Conjugate Gradient Method

Unlike Matlab and Octave, $R$ does not have a pcg function. As a result, we have written our own cg function in R . The code is divided into two files: cg.r contains the CG method, and driver_cg.r calls the CG function in cg.r by either setting up the sparse matrix $A$ or by using a function called Ax() to calculate the product of the matrix $A$ and the column vector $x$ directly from the vector without setting up a matrix.

Table 3.5: Convergence results for the test problem in R using (a) Gaussian elimination and (b) the conjugate gradient method. The tables list the mesh resolution $N$, the number of degrees of freedom (DOF), the finite difference norm $\left\|u-u_{h}\right\|_{L^{\infty}(\Omega)}$, the ratio of consecutive errors, and the observed wall clock time in HH:MM:SS.

| (a) Gaussian elimination |  |  |  |  |
| ---: | ---: | ---: | ---: | ---: |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | Time |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | N/A | $<00: 00: 01$ |
| 64 | 4,096 | $7.7812 \mathrm{e}-4$ | 3.8719 | $<00: 00: 01$ |
| 128 | 16,384 | $1.9766 \mathrm{e}-4$ | 3.9366 | $<00: 00: 01$ |
| 256 | 65,536 | $4.9807 \mathrm{e}-5$ | 3.9685 | $00: 00: 03$ |
| 512 | 262,144 | $1.2500 \mathrm{e}-5$ | 3.9843 | $00: 00: 26$ |
| 1,024 | $1,048,576$ | $3.1313 \mathrm{e}-6$ | 3.9921 | $00: 04: 22$ |
| 2,048 |  | out of memory |  |  |
| 4,096 |  | out of memory |  |  |
| 8,192 |  | out of memory |  |  |


| (b) Conjugate gradient method |  |  |  |  |  |
| ---: | ---: | ---: | ---: | ---: | ---: |
| $N$ | DOF | $\left\\|u-u_{h}\right\\|$ | Ratio | \#iter | Time |
| 32 | 1,024 | $3.0128 \mathrm{e}-3$ | N/A | 48 | $<00: 00: 01$ |
| 64 | 4,096 | $7.7810 \mathrm{e}-4$ | 3.8719 | 96 | $<00: 00: 01$ |
| 128 | 16,384 | $1.9765 \mathrm{e}-4$ | 3.9368 | 192 | $<00: 00: 01$ |
| 256 | 65,536 | $4.9797 \mathrm{e}-5$ | 3.9690 | 387 | $00: 00: 07$ |
| 512 | 262,144 | $1.2494 \mathrm{e}-5$ | 3.9856 | 783 | $00: 00: 53$ |
| 1,024 | $1,048,576$ | $3.1266 \mathrm{e}-6$ | 3.9961 | 1,581 | $00: 06: 57$ |
| 2,048 | $4,194,304$ | $7.8019 \mathrm{e}-7$ | 4.0075 | 3,192 | $01: 05: 52$ |
| 4,096 | $16,777,216$ | $1.9366 \mathrm{e}-7$ | 4.0286 | 6,452 | $11: 16: 23$ |
| 8,192 | $67,108,864$ | $4.7401 \mathrm{e}-8$ | 4.0855 | 13,033 | $109: 27: 24$ |

The numerical results in Table 3.5 (b) are identical to the results in Tables 3.1 (b), 3.2 (b), 3.3 (b), and 3.4 (b). While the conjugate gradient method in R ran significantly slower than Octave and Matlab, the performance is similar to Scilab and better than FreeMat.

After investigating the results, we found out that a significant amount of run-time was being spent in performing matrix algebra (specifically, subtractions on sparse matrices). There might be more efficient alternate ways in R to perform such matrix operations and there might be a significant scope for improvement.

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[^0]:    ${ }^{1}$ Cleve Moler, plenary talk at the SIAM Annual Meeting 2009, Denver, CO, and ensuing personal communication.

[^1]:    ${ }^{2}$ http://www.ll.mit.edu/mission/isr/pmatlab/pmatlab.html
    ${ }^{3}$ http://cran.r-project.org/web/views/HighPerformanceComputing.html
    ${ }^{4}$ http://gpgpu.org/2009/06/14/r-gpgpu

[^2]:    ${ }^{5}$ Another assignment operator <<- in R assigns values in global scope if the variable does not exist or overwrites the global variable if it exists. Keeping these operator behaviors in mind, one can argue that the $=$ operator is probably safer to use. However, we will continue to use <- as it is the common practice to assign values in $R$

