

# Project 1

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We present our Ferrari algorithm for solving linear equations. We wrote the one-dimensional Poisson equation, utilizing Dirichlet boundary conditions, as a linear set of equations and as a tridiagonal matrix. We compared a specialized algorithm for solving the tridiagonal matrix to an LU-decomposition of said matrix. Our best algorithm, the specialized solver, runs as  $4n$  FLOPS with  $n$  the dimensionality of the matrix.

## INTRODUCTION

As an introduction to the central ideas of the class, we studied the one-dimensional Poisson equation with Dirichlet boundary conditions. Namely, transforming the differential equation into a set of linear equations, and consequently, a matrix. We implemented a general algorithm, a specialized algorithm, and a LU-decomposition algorithm. In the course of this report, we introduce the theoretical model and the different algorithms we developed, then discuss the results of the different methods. Important points of comparison lie with relative error and relative speed of the calculation due to FLOPS.

## THEORY

### Theoretical solution of the one dimensional Poisson equation

In general, the one dimensional Poisson equation reads as follows:

$$-u''(x) = f(x) \quad (1)$$

Through discretized approximation of  $u$ , we can solve for  $f$  using a set of linear equations:

$$f(x) = -\frac{v_{i+1} + u_{i-1} - 2u_i}{h^2}; i = 1, \dots, n \quad (2)$$

Using Dirichlet boundary conditions,  $u_0 = u_{n+1} = 0$ , We can then rewrite this as a set of linear equations in the form of a tridiagonal matrix:

$$\hat{A} \cdot \hat{u} = \hat{f} \quad (3)$$

Consequently, we can solve these linear equations through forward and backward substitution.

### Specific Poisson equation

For our purposes of solving the Poisson equation, we assume a function

$$f(x) = 100e^{-10x} \quad (4)$$

and a closed form solution with the Dirichlet boundary conditions:

$$u(x) = 1 - (1 - e^{-10})x - e^{-10x} \quad (5)$$

### Solving a Tridiagonal Matrix

This is general tridiagonal matrix:

$$\begin{bmatrix} d_1 & e_1 & 0 & 0 \\ e_1 & d_2 & e_2 & 0 \\ 0 & e_2 & d_3 & e_{i-n} \\ 0 & 0 & e_{i-n} & d_n \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \\ u_3 \\ u_n \end{bmatrix} = \begin{bmatrix} v_1 \\ v_2 \\ v_3 \\ v_n \end{bmatrix}$$

We can reduce this generalized matrix into an upper triangular matrix from forward and backward substitution.

$$\tilde{f}_i = f_i - \frac{\tilde{f}_{i-1}e_{i-1}}{\tilde{d}_{i-1}}; \tilde{d}_i = d_i - e_{i-1}^2/\tilde{d}_{i-1} \quad (6)$$

$$u_i = (\tilde{f}_i - e_i u_{i+1} + 1)/\tilde{d}_i \quad (7)$$

where  $d_i$  are the diagonal matrix elements and  $e_i$  are the off diagonal matrix elements.

$$\begin{bmatrix} d_1 & e_1 & 0 & 0 \\ 0 & \tilde{d}_2 & e_2 & 0 \\ 0 & 0 & \tilde{d}_3 & e_{n-1} \\ 0 & 0 & 0 & \tilde{d}_n \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \\ u_3 \\ u_n \end{bmatrix} = \begin{bmatrix} v_1 \\ v_2 \\ v_3 \\ v_n \end{bmatrix}$$

We were able to create another program that performed matrix mathematics to solve for the set of linear equations. Taking a tridiagonal matrix, we can solve it using LU decomposition.

### LU-decomposition

$$\mathbf{A} = \mathbf{L}\mathbf{U} \quad (8)$$

$\mathbf{A}$ , the matrix can be decomposed into the product of  $\mathbf{U}$ , upper triangular matrix and  $\mathbf{L}$ , a lower triangular matrix

$$\mathbf{L}\mathbf{U}\mathbf{x} = \mathbf{b} \quad (9)$$

Which allows you to replace  $\mathbf{A}$  and then use either triangular matrix to solve the problem.

$$\mathbf{L}\mathbf{y} = \mathbf{b} \quad (10)$$

$$\mathbf{U}\mathbf{x} = \mathbf{x} \quad (11)$$

### ALGORITHMS

The matrix that represents the set of linear equations referred to in equation (2) appears as:

$$\begin{bmatrix} 2 & -1 & 0 & 0 \\ -1 & 2 & -1 & 0 \\ 0 & -1 & 2 & -1 \\ 0 & 0 & -1 & 2 \end{bmatrix}$$

This form can be generalized to any size  $n \times n$  matrix, where the diagonal elements  $d_i$  each equal 2 and the off diagonal elements  $e_i$  each equal -1.

Equations (6) and (7), representative of forward and backwards substitution, can be generalized for this specific matrix, yielding:

### METHODS

### RESULTS AND DISCUSSIONS

#### Relative error

In order to properly test the effectiveness of our algorithm, we are measuring the closeness, or relative error of our analytic solution:

$$\epsilon_i = \log_{10}\left(\left|\frac{v_i - u_i}{u_i}\right|\right); i = 1, \dots, n \quad (12)$$

$u_i$  is the analytic solution, and  $v_i$  is the numerical solution.

We ran the program with the step sizes of  $n = 1, 2$ , and 3, corresponding to a  $10 \times 10$ ,  $100 \times 100$ , and  $1000 \times 1000$  matrix. We also optimized our program through specialization. Diagonal matrix elements were set to 2 and off-diagonal elements were set to -1. FLOPS were reduced from 9 in the general case to 4 in the specialized case. The program was much more efficient, and faster as a consequence to the precalculation.

Analyzing step size, there is a clear difference from  $N=10^1$  and  $N=10^2$ , while  $N=10^2$  to  $N=10^3$  does not seem to superficially change as much as is demonstrated in the graphs below.

Further investigation shows that the overabundance of data points does give us the full curve, it also results in a much larger error, as shown in the table below. There seems to be a saddle point situated around

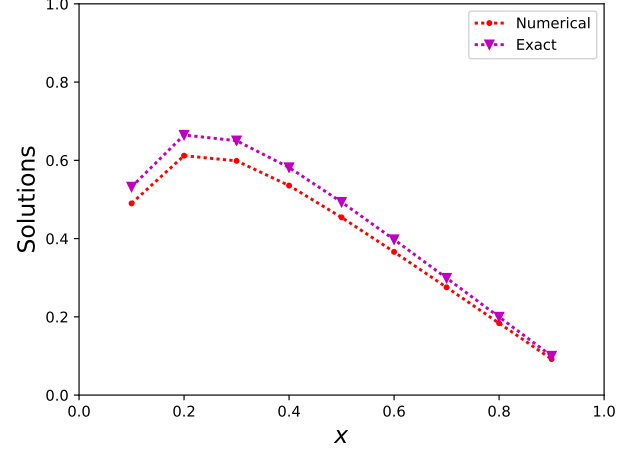


FIG. 1.  $n = 1$ . The x axis is the x values in the range  $[0,1]$  and the y axis is the resulting u values.

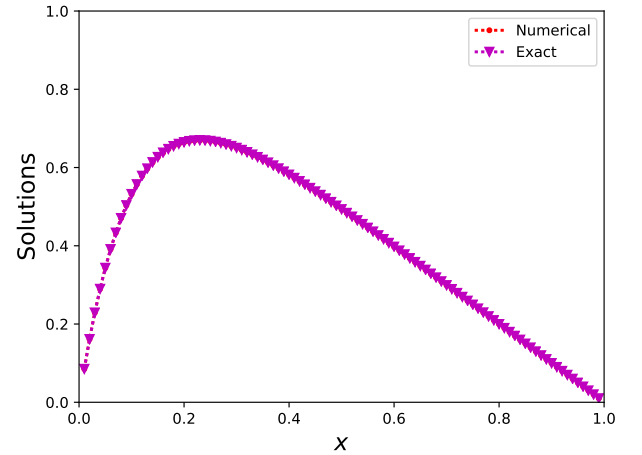


FIG. 2.  $n = 2$ . The x axis is the x values in the range  $[0,1]$  and the y axis is the resulting u values.

$N=10^2$ , wherein the calculation is optimized in regard to completeness of the graph and uncertainty from the numerical solution.

Step Size ( $n$ )	Max Relative Error ( $ \epsilon $ )
1	1.101
2	3.079
3	5.079
4	7.079
5	9.079
6	11.50
7	12.27

The above table shows us the absolute valued relative error results from varying step sizes of  $N = 10^n$  from 1 to

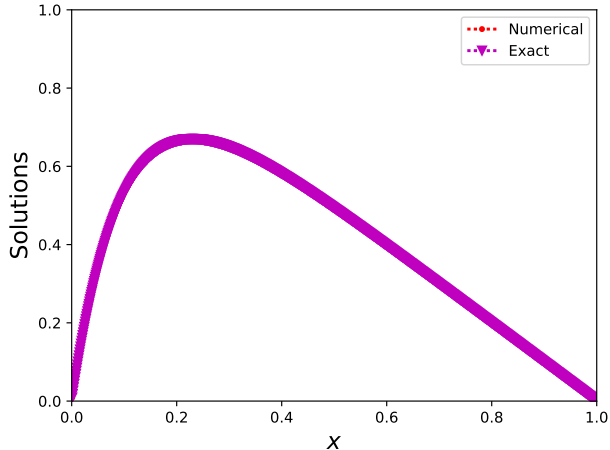


FIG. 3.  $n=3$ . The x axis is the x values in the range  $[0,1]$  and the y axis is the resulting u values.

7. The step size error for  $n=7$  seems to be the breaking point, the error starts to widely fluctuate. Therefore, the largest step size we would recommend for an accurate result is  $n=6$ .

For further analysis, we included a timer in both the specialized program and the LU program, and compared the timing values for, for as large square matrices as either could reasonably calculate.

Column Size	Specialized Time (s)	LU Time (s)
$10^1$	$1.7 \cdot 10^{-5}$	$1.48 \cdot 10^{-4}$
$10^2$	$2.4 \cdot 10^{-5}$	$3.79 \cdot 10^{-4}$
$10^3$	$2.12 \cdot 10^{-4}$	$9.25 \cdot 10^{-3}$
$10^4$	$1.24 \cdot 10^{-3}$	1.18
$10^5$	$1.28 \cdot 10^{-2}$	—
$10^6$	$1.26 \cdot 10^{-1}$	—
$10^7$	1.26	—

We found that specialized time was around an order of magnitude faster than LU time, as expected. When we reached powers of  $10^4$ , the difference jumped several orders of magnitude. We were unable to compute a LU matrix at  $N = 10^5$ , initializing the program with this input would immediately seize the program up. For the specialized solver,  $N = 10^7$  was the most reasonable limit, guessing at the size of the output file for  $N = 10^7$ , 400MB, we can assume that  $N = 10^8$  would not only take 10 or more seconds to compute but would result in a file size well over a GB, if not several GB. Orders above this would only increase file size and computation time.

## CONCLUSIONS

### Results

We investigated a few ways of solving a tridiagonal matrix problem. As expected, specialization outperforms generalized "brute force" methods. Specialization is faster, and in this case, can be taken out into greater orders of magnitude. Although there is improvement in this realm, we found that the propagation of numerical errors becomes unmanageable after a certain point. Our sweet spot was around the order of  $10^2$ .

### Future Prospects

There may be a method to optimize the LU decomposition program, to allow for us to solve problems greater than  $10^4$ , it might be useful to look into this.

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- [1] G. A. Miller, A. K. Oppen, and E. J. Stephenson, Annu. Rev. Nucl. Sci. **56**, 253 (2006).