EMTH211-19S2 Engineering Linear Algebra and Statistics

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LINEAR SYSTEMS

PROBLEM

Solve a **linear** system of equations

 $A \mathbf{x} = \mathbf{b}$.

GAUSSIAN ELIMINATION - A REFRESHER

Given a system of linear equations $A \mathbf{x} = \mathbf{b}$, Gaussian elimination solves this system in **two** steps.

Q ROW REDUCTION: Row operations are used to reduce the augmented matrix $\begin{bmatrix} A & \mathbf{b} \end{bmatrix}$ to an **upper triangular** form; that is,

$$[A \mid \mathbf{b}] \longrightarrow \begin{bmatrix} * & * & * & \cdots & * & * & * \\ 0 & * & * & \cdots & * & * & * \\ 0 & 0 & * & \cdots & * & * & * \\ \vdots & \vdots & \vdots & \ddots & \vdots & \vdots & \vdots \\ 0 & 0 & 0 & \cdots & 0 & * & * \end{bmatrix}.$$

BACK SUBSTITUTION is then used on the row reduced matrix to solve the system of equations.

Gaussian elimination - Row Reduction

- Elementary row operations are operations that leave the solution unchanged.
- Since a row of the augmented matrix corresponds to an equation in the system, there are three types of elementary row operations.
 - Interchange two rows

$$R_i \leftrightarrow R_j$$
.

ullet Add a multiple lpha of one row to another row

$$R_i \rightarrow R_i + \alpha R_i$$
.

• Multiply a row by a non-zero constant β

$$R_i \rightarrow \beta R_i$$
.

• It is the first two operations that we use in row reduction.

Gaussian elimination - Row Reduction

$$[A \mid \mathbf{b}] \longrightarrow \begin{bmatrix} a_{11} & a_{12} & a_{13} & \cdots & a_{1n} & b_1 \\ a_{21} & a_{22} & a_{32} & \cdots & a_{2n} & b_2 \\ a_{31} & a_{32} & a_{33} & \cdots & a_{3n} & b_3 \\ \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \end{bmatrix}$$

• We wish to eliminate a_{21} . We can do this by

$$R_2 \to R_2 - \frac{\alpha_{21}}{\alpha_{11}} \frac{R_1}{R_1}$$
.

This will update the entire second row.

• We now wish to eliminate a_{31} . Thus

$$R_3 \rightarrow R_3 - \frac{\alpha_{31}}{\alpha_{11}} \, \textcolor{red}{R_1}$$

and so on down the first column.

Gaussian elimination - Row Reduction

$$[A \mid \mathbf{b}] \longrightarrow \begin{bmatrix} \boxed{a_{11}} & a_{12} & a_{13} & \cdots & a_{1n} & b_1 \\ 0 & \tilde{a}_{22} & \tilde{a}_{23} & \cdots & \tilde{a}_{2n} & \tilde{b}_2 \\ 0 & \tilde{a}_{32} & \tilde{a}_{33} & \cdots & \tilde{a}_{3n} & \tilde{b}_3 \\ \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \end{bmatrix}$$

• We wish to eliminate \tilde{a}_{32} . We do this by

$$R_3 \rightarrow R_3 - \frac{\tilde{a}_{32}}{\tilde{a}_{22}} \frac{R_2}{R_2}$$

- and so on down the second column.
- Repeating this procedure on the remaining columns of A, we eventually obtained the row reduced form of the augmented matrix.

Gaussian elimination – Row Reduction

- At each step an **entire** row of the augmented matrix is updated.
- When working on the jth column, we add multiplies of the jth row to eliminate the entry.
- The elements on the diagonal are called **pivots**. If any pivot is zero we must use a row interchange to try to obtain a non-zero pivot.
- If it is not possible to get a non-zero pivot then the system has either no solution or infinitely many solutions.
- At the end of row reduction, the matrix has the form

GAUSSIAN ELIMINATION - BACK SUBSTITUTION

- If all the pivots are non-zero, the system has an unique solution.
- After row reduction, the linear system has the form

$$\begin{bmatrix} \mathbf{r}_{11} & * & * & \cdots & * & * \\ 0 & \mathbf{r}_{22} & * & \cdots & * & * \\ \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & 0 & 0 & \cdots & \mathbf{r}_{n-1,n-1} & * \\ 0 & 0 & 0 & \cdots & 0 & \mathbf{r}_{nn} \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_{n-1} \\ x_n \end{bmatrix} = \begin{bmatrix} \beta_1 \\ \beta_2 \\ \vdots \\ \beta_{n-1} \\ \beta_n \end{bmatrix}$$

• The last row gives

$$x_n = \frac{\beta_n}{r_{nn}}.$$

• The second last row gives

$$x_{n-1} = \frac{\beta_{n-1} + r_{n-1,n} x_n}{r_{n-1,n-1}} = \frac{r_{nn} \beta_{n-1} + r_{n-1,n} \beta_n}{r_{n-1,n-1} r_{nn}}$$

and so on.

Gaussian elimination – What can go wrong?

Consider the system

$$0.01x_1 + x_2 = 1$$
$$x_1 - x_2 = 0.$$

- Note that each coefficient and the right hand side is a two-digit number (for example, 0.01 is represented as 0.10×10^{-1})
- EXACT ARITHMETIC:
 - Row Reduction:

$$\begin{bmatrix} 0.01 & 1 & 1 \\ 1 & -1 & 0 \end{bmatrix} \rightarrow \begin{bmatrix} 0.01 & 1 & 1 \\ 0 & -101 & -100 \end{bmatrix}$$

Back Substitution:

$$\mathbf{x} = \begin{bmatrix} \frac{100}{101} \\ \frac{100}{101} \end{bmatrix} = \begin{bmatrix} 0.990099 \dots \\ 0.990099 \dots \end{bmatrix}.$$

 The best we could expect to obtain from a two-digit computer would be the closest two-digit numbers to the exact solution; namely

$$\mathbf{x} \approx \begin{bmatrix} 0.99 \\ 0.99 \end{bmatrix}$$
.

So what happens on our hypothetical two-digit computer?

GAUSSIAN ELIMINATION - WHAT CAN GO WRONG?

- FINITE PRECISION ARITHMETIC: (on a two-digit computer)
 - Row Reduction:

$$\begin{bmatrix} 0.01 & 1 & | & 1 \\ 1 & -1 & | & 0 \end{bmatrix} \xrightarrow{\text{two-digit}} \begin{bmatrix} 0.01 & 1 & | & 1 \\ 0 & 100 & | & -100 \end{bmatrix}$$
digits.

since, to two digits,

$$-1 - 100 = -(0.10 \times 10^{1}) - (0.10 \times 10^{3}) = -0.10 \times 10^{3} = -100.$$

Back Substitution:

$$\mathbf{x} \approx \begin{bmatrix} 0 \\ 1 \end{bmatrix}$$
.

- While we still have a good approximation for x_2 , $x_1 = 0$ which is a poor (to say the least) approximation for 0.99.
- However if we first do a row swap then the computation becomes
 - Row Reduction:

$$\begin{bmatrix} 1 & -1 & | & 0 \\ 0.01 & 1 & | & 1 \end{bmatrix} \xrightarrow{\text{two-digit}} \begin{bmatrix} 1 & -1 & | & 0 \\ 0 & 1 & | & 1 \end{bmatrix} \xrightarrow{\text{mey improve}} \text{rove}$$

$$\mathbf{x} \approx \begin{bmatrix} 1 \\ 1 \end{bmatrix}! \qquad \text{metrix}$$

Back Substitution:

$$\mathbf{x} \approx \begin{bmatrix} 1 \\ 1 \end{bmatrix}!$$
 accuracy of the matrix

GAUSSIAN ELIMINATION - THE MORAL

- This example demonstrates that, with finite precision arithmetic, different pivots may lead to dramatically different answers through Gaussian elimination. Some of these answers may be acceptable but some can be totally unacceptable.
- We need to develop a pivot selection strategy to avoid (if possible) bad answers!
- But first we need to understand what is happening in this example.

GAUSSIAN ELIMINATION - BACKWARD ERROR ANALYSIS

- What system of equations do we need to solve with exact arithmetic in order to obtain the same row reduced matrix as we computed with finite precision arthmetic?
- Note that the first row of a matrix is unchanged by Gaussian elimination.
 In the second (that is, good) case, we have

$$\begin{bmatrix} 1 & -1 & 0 \\ 0.01 & 1 & 1 \end{bmatrix} \xrightarrow{\text{two-digit}} \begin{bmatrix} 1 & -1 & 0 \\ 0 & 1 & 1 \end{bmatrix} \xleftarrow{\text{exact}} \begin{bmatrix} 1 & -1 & 0 \\ 0.01 & 0.99 & 1 \end{bmatrix}$$

• We see that the two systems are close:

$$x_1 - x_2 = 0$$
 $x_1 - x_2 = 0$ $0.01x_1 + x_2 = 1$ two-digit $0.01x_1 + 0.99x_2 = 1$

- Just to reiterate, the second system is the system that we would need to solve using exact arithmetic to obtain the same solution as solving the first system by two-digit arithmetic.
- Therefore the effect of using a two-digit computer to perform Gaussian elimination is to solve a slightly different (that is perturbed) system.

GAUSSIAN ELIMINATION - BACKWARD ERROR ANALYSIS

• What about the first case which gave us such a bad answer?

$$\begin{bmatrix} 0.01 & 1 & 1 \\ 1 & -1 & 0 \end{bmatrix} \xrightarrow{\text{two-digit}} \begin{bmatrix} 0.01 & 1 & 1 \\ 0 & -100 & -100 \end{bmatrix} \xleftarrow{\text{exact}} \begin{bmatrix} 0.01 & 1 & 1 \\ 1 & 0 & 0 \end{bmatrix}$$

• In this case the two systems are not close:

$$0.01x_1 + x_2 = 1$$
 $0.01x_1 + x_2 = 1$ $x_1 - x_2 = 0$ $x_1 = 0$ two-digit exact

- In both cases, the computed solution may be viewed as the exact solution to a perturbed system.
- This approach is called backward error analysis. The error in the computation is pushed back from the computation and placed on the data.
- It is an useful approach since frequently data is already inaccurate through experimental or modelling errors.

GAUSSIAN ELIMINATION - BACKWARD ERROR ANALYSIS

- Ideally we would like that errors created by the computer solution would be no larger than the errors inherent in the data.
- Thus we would like the computer solution to be an exact solution to a slightly perturbed problem.
- However this example demonstrates that Gaussian elimination, *unmodified*, may result in the solution to a **greatly perturbed** problem.
- In both row reduction and backwards substitution we are dividing by pivots. Therefore if a pivot is small, any errors will be magnified.
- In the bad case, the pivot was 0.01. Therefore, each time we divide by the pivot, errors could be magnified by a factor of 1/0.01 = 100 (and this does not allow for errors in the pivot itself).
- In the good case, the pivot was 1. Now errors are not magnified.
- This suggests the strategy avoid small pivots.

Gaussian elimination - Small Pivots

• But (and there is always a but!) what about the rescaled system

$$10x_1 + 1000x_2 = 1000$$
$$x_1 - x_2 = 0?$$

In this case we obtain

$$\begin{bmatrix} 10 & 1000 & 1000 \\ 1 & -1 & 0 \end{bmatrix} \xrightarrow{\text{two-digit}} \begin{bmatrix} 10 & 1000 & 1000 \\ 0 & -100 & -100 \end{bmatrix}$$

the same bad result.

- Here the pivot is 10 so the bad result is not a consequence of a small pivot.
- If we do a row swap and use the smaller pivot, we obtain

$$\begin{bmatrix} 1 & -1 & 0 \\ 10 & 1000 & 1000 \end{bmatrix} \xrightarrow{\text{two-digit}} \begin{bmatrix} 1 & -1 & 0 \\ 0 & 1000 & 1000 \end{bmatrix}$$

and thus the good result.

GAUSSIAN ELIMINATION - SCALING

- However, in a properly scaled system, we should avoid small pivots.
- Ideally, the system should be scaled before it is given to the computer to solve.
 - We can scale an *equation* by multiplying by a non-zero constant. This is equivalent to an elementary row operation.
 - We can also scale the variables of a problem

$$\chi_i \rightarrow \alpha_i \chi_i$$

where α_i are non-zero constants. This does not correspond to an elementary row operation (in fact, it is an elementary *column* operation).

• There is *no* strategy that will work in all situations. However one should use scaling, if possible, to avoid "large" and very small non-zero numbers in the system. The errors in the example above arose from adding a large number to a small number and the consequent loss of precision when using finite precision arithmetic.

GAUSSIAN ELIMINATION - SCALING

• For example, consider

$$x_1 - 0.0001x_2 = 0$$
$$100x_1 + x_2 = 100.$$

- One might
 - scale the second equation

$$x_1 - 0.0001x_2 = 0$$
$$x_1 + 0.01x_2 = 1$$

• and scale $x_2 \rightarrow 0.01x_2 = \tilde{x}_2$ say

$$x_1 - 0.01\tilde{x}_2 = 0$$
$$x_1 + \tilde{x}_2 = 1.$$

• In the original system, the coefficients varied by six orders of magnitude; in the scaled system, they varied by only two orders of magnitude.

GAUSSIAN ELIMINATION - PARTIAL PIVOTING

- Again there is *no* pivoting strategy that will work in all situations.
- From a practical point of view, partial pivoting is the strategy most commonly used (MATLAB uses it by default). It balances low computational overhead with accuracy of the final answer.
- With partial pivoting, we swap rows so that the pivot always has the largest absolute value of all remaining entries in that column.
- For example, in the third column, we search the boxed area for the element with largest absolute value. We then use a row operation to move this element to the pivot position.

Gaussian elimination - Complete Pivoting

- An alternative is complete pivoting.
- In this case we search the submatrix

$$\begin{bmatrix} * & * & * & * & * & \cdots & * & * \\ 0 & * & * & * & \cdots & * & * \\ 0 & 0 & & * & * & \cdots & * & * \\ 0 & 0 & & * & * & \cdots & * & * \\ \vdots & \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & 0 & & * & * & \cdots & * & * \end{bmatrix}$$

for the largest (in absolute value) entry. We then use row *and* column swaps to bring this element to the pivot position.

• Whilst this strategy is better than partial pivoting, it's use is rarely justified since the computational overhead is *significantly* higher.

Gaussian elimination - Example of Partial Pivoting

Consider the system whose augmented matrix is

- Column 1:
 - We want -4 moved to the pivot position so

$$R_1 \leftrightarrow R_3$$
.

• Gaussian elimination row operations

$$R_2 \rightarrow R_2 + \frac{1}{2} R_1$$

 $R_3 \rightarrow R_3 + \frac{1}{4} R_1$
 $R_4 \rightarrow R_4 - \frac{1}{4} R_1$.

Gaussian elimination - Example of Partial Pivoting

$$\begin{bmatrix} -4 & 0 & -1 & 0 & 1 \\ 0 & 2 & 2.25 & 0 & 0.75 \\ 0 & 0 & 2.75 & -1 & 1.25 \\ 0 & 1 & -0.5 & 0 & 1.5 \end{bmatrix}$$

- Column 2:
 - We want 2 moved to the pivot position so

$$R_2 \leftrightarrow R_4$$
.

• Gaussian elimination row operation

$$R_4 \rightarrow R_4 - \frac{1}{2} R_2$$

which completes the second column.

GAUSSIAN ELIMINATION - EXAMPLE OF PARTIAL PIVOTING

$$\begin{bmatrix} -4 & 0 & -1 & 0 & 1 \\ 0 & 2 & 2.25 & 0 & 0.75 \\ 0 & 0 & 2.75 & -1 & 1.25 \\ 0 & 0 & -1.625 & 0 & 1.125 \end{bmatrix}$$

- Column 3:
 - No row swaps required.
 - Gaussian elimination row operation

$$R_4 \to R_4 + \tfrac{1.625}{2.75} \; R_3$$

which gives the row reduced form.

GAUSSIAN ELIMINATION - ANOTHER EXAMPLE

Solve the system whose augmented matrix is

$$\begin{bmatrix} 1 & 2 & \left| \frac{1}{3} \right| \\ 2 & 4.01 & \left| \frac{2}{3} \right] \end{bmatrix}$$

using a three-digit computer.

EXACT SOLUTION is

$$\mathbf{x} = \begin{bmatrix} \frac{1}{3} \\ 0 \end{bmatrix}.$$

With three-digit arithmetic, the augmented matrix becomes

- To obtain a three-digit representation of this system, we are forced to introduce inaccuracies (albeit small) in the vector **b**.
- So what happens?

GAUSSIAN ELIMINATION - ANOTHER EXAMPLE

Gaussian elimination with partial pivoting gives

$$\begin{bmatrix} 1 & 2 & 0.333 \\ 2 & 4.01 & 0.667 \end{bmatrix} \rightarrow \begin{bmatrix} 2 & 4.01 & 0.667 \\ 1 & 2 & 0.333 \end{bmatrix} \xrightarrow{\text{three-digit}} \begin{bmatrix} 2 & 4.01 & 0.667 \\ 0 & -0.005 & -0.0005 \end{bmatrix}$$

- Note that the outcome of the row operation $R_2 \to R_2 \frac{1}{2} R_1$ depends on the order in which the arithmetic operations are performed.
 - If we compute $\frac{1}{2} (2R_2 R_1)$ we obtain

$$\frac{4 - 4.01}{2} = \frac{-0.01}{2} = -0.005$$

• whereas if we compute $R_2 - \frac{1}{2} R_1$ we obtain

$$2 - \frac{4.01}{2} = 2 - 2.01 = -0.010.$$

Finally the computed solution is

$$\mathbf{x} \approx \begin{bmatrix} 0.133 \\ 0.100 \end{bmatrix}$$
.

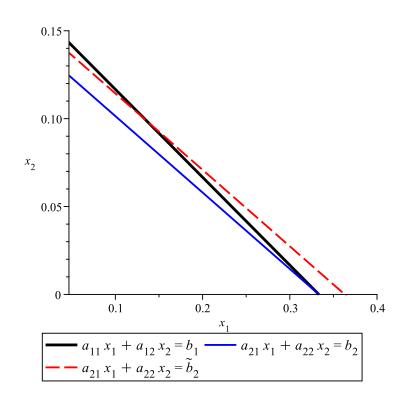
Gaussian elimination - Another Example

• Partial pivoting does not help in this case. The inaccuracies in the data are less than 0.1%. However

 $\begin{bmatrix} 0.133 \\ 0.100 \end{bmatrix} \qquad \begin{bmatrix} 0.333 \\ 0.000 \end{bmatrix}$ computed solution exact solution (to 3 significant figures)

- This is an example of an ill-conditioned system.
- Geometrically, this system represents two lines that are *almost parallel*.
- We will revisit ill-conditioned systems later in this course.

GAUSSIAN ELIMINATION - ILL-CONDITIONED SYSTEMS



How fast is fast?

- We have discussed accuracy issues for Gaussian elimination but what about speed? Is it a *fast* algorithm or do we need another algorithm?
- First we need a measure on how fast an algorithm is.
- Given that an algorithm is used on a computer (rather than by us), we measure the speed of a algorithm in terms of the number of slow operations that a computer has to perform.
- The slowest operations are the *floating point* operations; multiplication/division and addition/subtraction.
- Floating point operations are measured in terms of flops.
- Flops per second is one measure that is used to benchmark computer performance.
 - The fastest supercomputers today have performance in the 1 exaflop/sec $(10^{18} \text{ flop/sec}) \text{ range.}$
 - The fastest Intel processors have performance around 1 teraflop (10¹²) flop)/sec.

 Graphic processing units (GPUs) used on dedicated graphics cards are, for
 - double precision floating point, in the 10 teraflop range.
- So what is the flop count for Gaussian elimination and should it concern us?

Gaussian elimination – flop count

Two well-known (!) formulae that will help us in the flop count are

$$\sum_{i=1}^{n} i = \frac{1}{2} (n^2 + n)$$

$$\sum_{i=1}^{n} i^2 = \frac{1}{3} n^3 + \frac{1}{2} n^2 + \frac{1}{6} n$$

Row Reduction:

• FIRST COLUMN: We use row operations of the form

$$R_i \rightarrow R_i - a_{i1} \times \frac{R_1}{a_{11}}$$

- We need to divide (that is multiply!) each element in R_1 by α_{11} . However we do not need to do this for the first element in this row and so there are $n \times to$ perform (remember though there are only n-1 remaining elements in the coefficient matrix, we also have the element b_1 in the augmented matrix). We update the first row so that the pivot element is now 1.
- Each row we need to perform one addition and one multiplication. Again we do not need to do this for the first element since we *know* that this element is 0 by construction. Thus there are $n \times and n + b$ to perform.
- Since there are n-1 rows we need to update we have a total count

$$\lceil (n-1)n+n \rceil = n^2 \times \text{ and } (n-1)n+$$

operations to perform. Thus $2n^2 - n$ flops are required.

GAUSSIAN ELIMINATION - FLOP COUNT

Row Reduction:

• Remaining Columns: The count is the same for remaining columns except that, for the j^{th} column, we are performing the operations on a $(n-j)\times(n-j+1)$ submatrix. Thus the flop count is

$$2(n-j+1)^2 - (n-j+1)$$

for
$$j = 2, \ldots, n$$
.

• The total flop count for row reduction is

$$\sum_{j=1}^{n} [2(n-j+1)^2 - (n-j+1)] = \sum_{i=1}^{n} [2i^2 - i] \qquad (i = n-j+1)$$
$$= \frac{2}{3} n^3 + \frac{1}{2} n^2 - \frac{1}{6} n$$
$$\approx \frac{2}{3} n^3$$

for large n. We are only interested in large n, so we write

Row Reduction =
$$O(\frac{2}{3} n^3)$$
.

After the above row reduction, the linear system has the form

$$\begin{bmatrix} 1 & * & * & \cdots & * & * \\ 0 & 1 & * & \cdots & * & * \\ \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & 0 & 0 & \cdots & 1 & * \\ 0 & 0 & 0 & \cdots & 0 & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_{n-1} \\ x_n \end{bmatrix} = \begin{bmatrix} \beta_1 \\ \beta_2 \\ \vdots \\ \beta_{n-1} \\ \beta_n \end{bmatrix}$$

The last row gives

$$x_n = \beta_n$$
.

• The second last row gives

$$x_{n-1} = \beta_{n-1} - r_{n-1,n} x_n$$

and so on.

GAUSSIAN ELIMINATION - FLOP COUNT

• BACK SUBSITUTION: In the $(n-j+1)^{\text{th}}$ row (that is, the j^{th} row from the bottom) we need $(j-1) \times$ and (j-1) + (check; note after the above row reduction, the pivot elements are all 1). Therefore the flop count is

$$\sum_{j=1}^{n} \left[2(j-1) \right] = n^2 - n$$

We write

Back Substitution =
$$O(n^2)$$
.

• The total flop count for Gaussian elimination without pivoting is

Gaussian Elimination =
$$O(\frac{2}{3} n^3)$$

since $n^2 << n^3$ for large n.

- What about pivoting? It takes one addition to compare two numbers (computers compute the difference and then check the sign).
- Partial Pivoting: In the j^{th} column we need to compare n-j+1 elements to obtain the pivot. This involves n-j comparisons. Therefore the flop count is

$$\sum_{j=1}^{n} [n-j] = \sum_{i=0}^{n-1} i = \frac{1}{2} n^2 - \frac{1}{2} n$$

or

Partial Pivoting = $O(\frac{1}{2} n^2)$

GAUSSIAN ELIMINATION - FLOP COUNT

• Complete Pivoting: In the jth column we need to compare $(n-j+1)^2$ elements to obtain the pivot. This involves $(n-j+1)^2-1$ comparisons. Therefore the flop count is

$$\sum_{j=1}^{n} \left[(n-j+1)^2 - 1 \right] = \sum_{i=1}^{n} \left[i^2 - 1 \right] = \frac{1}{3} n^3 + \frac{1}{2} n^2 - \frac{5}{6} n$$

or

Complete Pivoting =
$$O(\frac{1}{3} n^3)$$

 Whereas partial pivoting does not increase the flop count significantly, complete pivoting does.

> Gaussian Elimination with partial pivoting = $O(\frac{2}{3} n^3)$ Gaussian Elimination with complete pivoting = $O(n^3)$

As a comparison we have (check if you like)

Gauss-Jordan Elimination = $\mathbb{O}(n^3)$ Computing $A^{-1} = \mathbb{O}(2n^3)$ Multiplying two $n \times n$ matrices = $\mathbb{O}(2n^3)$

- A generic Gaussian elimination routine would take $O(\frac{8}{3} n^3)$ to compute the inverse. The algorithm has to use the fact that there are many zeros in the identity matrix to obtain the above speed. Computing the inverse is the **least efficient** method of solving a system of linear equations.
- The fact that row reduction grows as *cube* of the number of variables is a very big issue. If we double the size of the problem, then the problem will take at least 8 times longer to solve.

IS SPEED AN ISSUE?

- These days most computer monitors have at least 10^6 pixels. Thus, in updating such a monitor, one could get a system of equations that have 10^6 variables! If we used Gaussian elimination, how long would it take?
- The flop count would be

$$0.67 \times 10^{18}$$
.

That is, on a

supercomputer 0.67 sec state of the art graphics processor 67,000 sec (18 hours) state of the art PC 670,000 sec (8 days)

- Therefore Gauss elimination is NOT a fast algorithm.
- More precisely, generic Gauss elimination is not fast. As seen in the first lab, it can be fast for certain problems.