Lecture 7 Gaussian Elimination with Pivoting

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Naive Gaussian Elimination Algorithm

- Forward Elimination
- + Backward substitution
- = Naive Gaussian Elimination

- Identify why our basic GE method is "naive": identify where the errors come from?
 - division by zero, near-zero
- Propose strategies to eliminate the errors
 - partial pivoting, complete pivoting, scaled partial pivoting
- Investigate the cost: does pivoting cost too much?
- Try to answer "How *accurately* can we solve a system with or without pivoting?"
 - Analysis tools: norms, condition number, …

Why is our basic GE "naive"?

Example

$$A = \begin{bmatrix} 0 & 2 & 3 \\ 4 & 5 & 6 \\ 7 & 8 & 9 \end{bmatrix}$$

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Example

$$A = \begin{bmatrix} 1e - 10 & 2 & 3 \\ 4 & 5 & 6 \\ 7 & 8 & 9 \end{bmatrix}$$

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Solve:

$$A = \begin{bmatrix} 2 & 4 & -2 & -2 \\ 1 & 2 & 4 & -3 \\ -3 & -3 & 8 & -2 \\ -1 & 1 & 6 & -3 \end{bmatrix} \qquad b = \begin{bmatrix} -4 \\ 5 \\ 7 \\ 7 \end{bmatrix}$$

Note that there is nothing "wrong" with this system. A is full rank. The solution exists and is unique.

Form the augmented system.

$$\begin{bmatrix} 2 & 4 & -2 & -2 & | & -4 \\ 1 & 2 & 4 & -3 & | & 5 \\ -3 & -3 & 8 & -2 & | & 7 \\ -1 & 1 & 6 & -3 & | & 7 \end{bmatrix}$$

Subtract 1/2 times the first row from the second row, add 3/2 times the first row to the third row, add 1/2 times the first row to the fourth row. The result of these operations is:

$$\begin{bmatrix} 2 & 4 & -2 & -2 & | & -4 \\ 0 & 0 & 5 & -2 & | & 7 \\ 0 & 3 & 5 & -5 & | & 1 \\ 0 & 3 & 5 & -4 & | & 5 \end{bmatrix}$$

The *next* stage of Gaussian elimination will not work because there is a zero in the *pivot* location, \tilde{a}_{22} .

Swap second and fourth rows of the augmented matrix.

$$\begin{bmatrix} 2 & 4 & -2 & -2 & | & -4 \\ 0 & 3 & 5 & -4 & | & 5 \\ 0 & 3 & 5 & -5 & | & 1 \\ 0 & 0 & 5 & -2 & | & 7 \end{bmatrix}$$

Continue with elimination: subtract (1 times) row 2 from row 3.

$$\begin{bmatrix} 2 & 4 & -2 & -2 & | & -4 \\ 0 & 3 & 5 & -4 & | & 5 \\ 0 & 0 & 0 & -1 & | & -4 \\ 0 & 0 & 5 & -2 & | & 7 \end{bmatrix}$$

Another zero has appear in the pivot position. Swap row 3 and row 4.

[2]	4	-2	-2	-4
$\begin{bmatrix} 2\\0\\0\\0 \end{bmatrix}$	3	5	-4	$\begin{array}{c} -4\\5\\7\end{array}$
0	0	5	-2	7
0	0	0	-1	-4

The augmented system is now ready for backward substitution.

another example

$$\begin{bmatrix} \varepsilon & 1 \\ 1 & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} 1 \\ 2 \end{bmatrix}$$

Example

With Naive GE,

$$\begin{bmatrix} \varepsilon & 1 \\ 0 & (1 - \frac{1}{\varepsilon}) \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} 1 \\ 2 - \frac{1}{\varepsilon} \end{bmatrix}$$

Solving for x_1 and x_2 we get

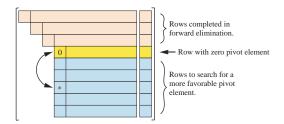
$$x_2 = \frac{2 - 1/\varepsilon}{1 - 1/\varepsilon}$$
$$x_1 = \frac{1 - x_2}{\varepsilon}$$

For $\varepsilon \approx 10^{-20}$, $x_1 \approx 0$, $x_2 \approx 1$

Partial Pivoting: Exchange only rows

- Exchanging rows does not affect the order of the x_i
- For increased numerical stability, make sure the largest possible pivot element is used. This requires searching in the partial column below the pivot element.
- Partial pivoting is usually sufficient.

To avoid division by zero, swap the row having the zero pivot with one of the rows below it.



To minimize the effect of roundoff, always choose the row that puts the largest pivot element on the diagonal, i.e., find i_p such that $|a_{i_p,i}| = \max(|a_{k,i}|)$ for k = i, ..., n

Partial Pivoting: Usually sufficient, but not always

• Partial pivoting is usually sufficient

Consider

$$\begin{bmatrix} 2 & 2c & | & 2c \\ 1 & 1 & | & 2 \end{bmatrix}$$

With Partial Pivoting, the first row is the pivot row:

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

and for large *c*:

$$\begin{bmatrix} 2 & 2c & | & 2c \\ 0 & -c & | & -c \end{bmatrix}$$

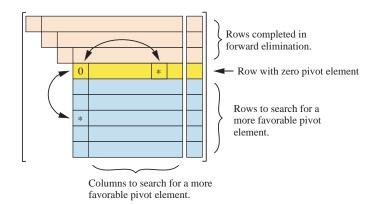
so that y = 1 and x = 0. (exact is x = y = 1)

 The pivot is selected as the largest in the column, but it should be the largest <u>relative</u> to the full submatrix.

Full (or Complete) Pivoting: Exchange both rows and columns

- Column exchange requires changing the order of the x_i
- For increased numerical stability, make sure the largest possible pivot element is used. This requires searching in the pivot row, *and* in all rows below the pivot row, starting the pivot column.
- Full pivoting is less susceptible to roundoff, but the increase in stability comes at a cost of more complex programming (not a problem if you use a library routine) and an increase in work associated with searching and data movement.

Full Pivoting



We simulate full pivoting by using a scale with partial pivoting.

- pick pivot element as the largest relative entry in the column (relative to the other entries in the row)
- do not swap, just keep track of the order of the pivot rows
- call this vector $\ell = [\ell_1, \ldots, \ell_n]$.

SPP Process

Determine a scale vector s. For each row

$$s_i = \max_{1 \leqslant j \leqslant n} |a_{ij}|$$

- 2 initialize $\ell = [\ell_1, \dots, \ell_n] = [1, \dots, n]$.
- select row i to be the row with the largest ratio

$$\frac{|a_{\ell_i 1}|}{s_{\ell_i}} \qquad 1 \leqslant i \leqslant n$$

- swap ℓ_i with ℓ_1 in ℓ
- Solution Now we need n-1 multipliers for the first column:

$$m_1 = \frac{a_{\ell_i 1}}{a_{\ell_1 1}}$$

- So the index to the rows are being swapped, NOT the actual row vectors which would be expensive
- If finally use the multiplier m_1 times row ℓ_1 to subtract from rows ℓ_i for $2 \leq i \leq n$ Image: A math a math

SPP Process continued

For the second column in forward elimination, we select row *j* that yields the largest ratio of

$$\frac{|a_{\ell_i,2}|}{s_{\ell_i}} \qquad 2 \leqslant i \leqslant n$$

2 swap ℓ_i with ℓ_2 in ℓ

(a) Now we need n - 2 multipliers for the second column:

$$m_2 = \frac{a_{\ell_i,2}}{a_{\ell_2 2}}$$

- finally use the multiplier m_2 times row ℓ_2 to subtract from rows ℓ_i for $3 \le i \le n$
- **(**) the process continues for row k
- note: scale factors are not updated

An Example

Consider

$$\begin{bmatrix} 2 & 4 & -2 \\ 1 & 3 & 4 \\ 5 & 2 & 0 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix} = \begin{bmatrix} 6 \\ -1 \\ 2 \end{bmatrix}$$

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• The first equation corresponds to the last index ℓ_n :

$$a_{\ell_n n} x_n = b_{\ell_n} \Rightarrow x_n = \frac{b_{\ell_n}}{a_{\ell_n n}}$$

2 The second equation corresponds to the second to last index ℓ_{n-1} :

$$x_{n-1} = \frac{1}{a_{\ell_{n-1}n-1}} \left(b_{\ell_{n-1}} - a_{\ell_{n-1}n} x_n \right)$$

The Algorithms

Listing 1: (forward) GE with SPP

```
Initialize \ell = [1, \ldots, n]
1
      Set s to be the max |a_{ij}| in each row
2
      for k=1 to n
3
         rmax = 0
4
         for i = k to n
5
             r = |a_{\ell,k}/s_{\ell_i}|
6
             if(r > rmax)
7
                rmax = r
8
                i = i
9
         end
10
         swap \ell_i and \ell_k
11
         for i = k + 1 to n
12
             xmult = a_{\ell_i k}/a_{\ell_k k}
13
             a_{\ell_{ik}} = xmult
14
             for j = k + 1 to n
15
                a_{\ell,i} = a_{\ell,i} - xmult \cdot a_{\ell,i}
16
17
             end
         end
18
      end
19
```

See Gauss algorithm on page 267 of NMC6 (p285 NMC5) < D > < D >

The Algorithms

Note: the multipliers are stored in the location $a_{\ell,k}$ in the text

Listing 2: (back solve) GE with SPP

```
for k=1 to n-1
1
          for i = k+1 to n
2
             b_{\ell_i} = b_{\ell_i} - a_{\ell_i k} b_{\ell_k}
3
          end
4
       end
5
      x_n = b_{\ell_n}/a_{\ell_n n}
6
      for i = n - 1 down to 1
7
          sum = b_{\ell_i}
8
          for j = i + 1 to n
9
             sum = sum - a_{\ell_i j} x_j
10
          end
11
          x_i = sum/a_{\ell_i,i}
12
       end
13
```

See Gauss algorithm on page 269 of NMC6 (p267 NMC5)

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Geometric Interpretation of Singularity

Consider a 2×2 system describing two lines that intersect

$$y = -2x + 6$$
$$y = \frac{1}{2}x + 1$$

The matrix form of this equation is

$$\begin{bmatrix} 2 & 1 \\ -1/2 & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} 6 \\ 1 \end{bmatrix}$$

The equations for two parallel but not intersecting lines are

$$\begin{bmatrix} 2 & 1 \\ 2 & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} 6 \\ 5 \end{bmatrix}$$

Here the coefficient matrix is singular (rank(A) = 1), and the system is inconsistent

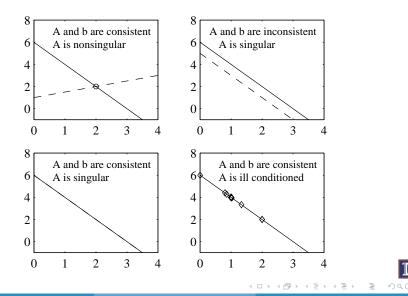
The equations for two parallel and coincident lines are

$$\begin{bmatrix} 2 & 1 \\ 2 & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} 6 \\ 6 \end{bmatrix}$$

The equations for two nearly parallel lines are

$$\begin{bmatrix} 2 & 1 \\ 2+\delta & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \end{bmatrix} = \begin{bmatrix} 6 \\ 6+\delta \end{bmatrix}$$

Geometric Interpretation of Singularity



Consider the solution of a 2×2 system where

$$b = \begin{bmatrix} 1\\ 2/3 \end{bmatrix}$$

One expects that the exact solutions to

$$Ax = \begin{bmatrix} 1\\ 2/3 \end{bmatrix}$$
 and $Ax = \begin{bmatrix} 1\\ 0.6667 \end{bmatrix}$

will be different. Should these solutions be a lot different or a little different?

Norms

Vectors:

$$||x||_{p} = (|x_{1}|^{p} + |x_{2}|^{p} + \dots + |x_{n}|^{p})^{1/p}$$
$$||x||_{1} = |x_{1}| + |x_{2}| + \dots + |x_{n}| = \sum_{i=1}^{n} |x_{i}|$$
$$||x||_{\infty} = \max(|x_{1}|, |x_{2}|, \dots, |x_{n}|) = \max_{i}(|x_{i}|)$$

Matrices:

$$\|A\| = \max_{x \neq 0} \frac{\|Ax\|}{\|x\|}$$
$$\|A\|_p = \max_{x \neq 0} \frac{\|Ax\|_p}{\|x\|_p}$$
$$\|A\|_1 = \max_{1 \leq i \leq m} \sum_{i=1}^m |a_{ij}|$$
$$\|A\|_{\infty} = \max_{1 \leq i \leq m} \sum_{i=1}^n |a_{ij}|$$

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Effect of Perturbations to b

Perturb b with δb such that

$$\frac{\|\delta b\|}{\|b\|} \ll 1,$$

The perturbed system is

$$A(x+\delta x_b)=b+\delta b$$

The perturbations satisfy

 $A\delta x_b = \delta b$

Analysis shows (see next two slides for proof) that

$$\frac{\|\delta x_b\|}{\|x\|} \leqslant \|A\| \|A^{-1}\| \frac{\|\delta b\|}{\|b\|}$$

Thus, the effect of the perturbation is small only if $||A|| ||A^{-1}||$ is small.

$$\frac{\|\delta x_b\|}{\|x\|} \ll 1$$
 only if $\|A\| \|A^{-1}\| \sim 1$

Let $x + \delta x_b$ be the *exact* solution to the perturbed system

$$A(x + \delta x_b) = b + \delta b \tag{1}$$

Expand

 $Ax + A\delta x_b = b + \delta b$

Subtract *Ax* from left side and *b* from right side since Ax = b

 $A\delta x_b = \delta b$

Left multiply by A^{-1}

$$\delta x_b = A^{-1} \delta b \tag{2}$$

Effect of Perturbations to *b* (Proof, p. 2)

Take norm of equation (2)

$$\left\|\delta x_{b}\right\| = \left\|A^{-1}\,\delta b\right\|$$

Applying consistency requirement of matrix norms

$$\|\delta x_b\| \leqslant \|A^{-1}\| \|\delta b\| \tag{3}$$

Similarly, Ax = b gives ||b|| = ||Ax||, and

$$\|b\| \leqslant \|A\| \|x\|$$

Rearrangement of equation (4) yields

$$\frac{1}{\|x\|} \leqslant \frac{\|A\|}{\|b\|}$$

(4)

(5)

Multiply Equation (4) by Equation (3) to get

$$\frac{\|\delta x_b\|}{\|x\|} \le \|A\| \|A^{-1}\| \frac{\|\delta b\|}{\|b\|}$$
(6)

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Summary:

If $x + \delta x_b$ is the *exact* solution to the perturbed system

$$A(x+\delta x_b)=b+\delta b$$

then

$$\frac{\|\delta x_b\|}{\|x\|} \leqslant \|A\| \|A^{-1}\| \frac{\|\delta b\|}{\|b\|}$$

Effect of Perturbations to A

Perturb A with δA such that

$$\frac{\|\delta A\|}{\|A\|} \ll 1,$$

The perturbed system is

$$(A + \delta A)(x + \delta x_A) = b$$

Analysis shows that

$$\frac{\|\delta x_A\|}{\|x+\delta x_A\|} \leqslant \|A\| \|A^{-1}\| \frac{\|\delta A\|}{\|A\|}$$

Thus, the effect of the perturbation is small *only if* $||A|| ||A^{-1}||$ is small.

$$\frac{\|\delta x_A\|}{\|x+\delta x_A\|} \ll 1 \quad \text{only if} \quad \|A\| \|A^{-1}\| \sim 1$$

Effect of Perturbations to both *A* and *b*

Perturb both A with δA and b with δb such that

$$\frac{\|\delta A\|}{\|A\|} \ll 1 \quad \text{and} \quad \frac{\|\delta b\|}{\|b\|} \ll 1$$

The perturbation satisfies

$$(A + \delta A)(x + \delta x) = b + \delta b$$

Analysis shows that

$$\frac{\|\delta x\|}{\|x+\delta x\|} \ \leqslant \ \frac{\|A\|\|A^{-1}\|}{1-\|A\|\|A^{-1}\|\frac{\|\delta A\|}{\|A\|}} \left[\frac{\|\delta A\|}{\|A\|}+\frac{\|\delta b\|}{\|b\|}\right]$$

Thus, the effect of the perturbation is small *only if* $||A|| ||A^{-1}||$ is small.

$$\frac{\|\delta x\|}{\|x+\delta x\|} \ll 1 \quad \text{only if} \quad \|A\| \|A^{-1}\| \sim 1$$

The condition number

 $\kappa(A) \equiv \|A\| \|A^{-1}\|$

indicates the sensitivity of the solution to perturbations in A and b. The condition number can be measured with any p-norm. The condition number is always in the range

 $1\leqslant \kappa(A)\leqslant \infty$

- $\kappa(A)$ is a mathematical property of A
- Any algorithm will produce a solution that is sensitive to perturbations in A and b if κ(A) is large.
- In exact math a matrix is either singular or non-singular. $\kappa(A) = \infty$ for a singular matrix
- $\kappa(A)$ indicates how close A is to being numerically singular.
- A matrix with large κ is said to be ill-conditioned

In Practice, applying Gaussian elimination with partial pivoting and back substitution to Ax = b gives the **exact solution**, \hat{x} , to the **nearby problem**

 $(A+E)\hat{x} = b$ where $||E||_{\infty} \leq \varepsilon_m ||A||_{\infty}$

Gaussian elimination with partial pivoting and back substitution "gives exactly the right answer to nearly the right question." — Trefethen and Bau An algorithm that gives the exact answer to a problem that is near to the original problem is said to be **backward stable**. Algorithms that are not backward stable will tend to amplify roundoff errors present in the original data. As a result, the solution produced by an algorithm that is not backward stable will not necessarily be the solution to a problem that is close to the original problem.

Gaussian elimination without partial pivoting is *not* backward stable for arbitrary *A*.

If *A* is symmetric and positive definite, then Gaussian elimination without pivoting in backward stable.

Let \hat{x} be the numerical solution to Ax = b. $\hat{x} \neq x$ (x is the exact solution) because of roundoff.

The **residual** measures how close \hat{x} is to satisfying the original equation

$$r = b - A\hat{x}$$

It is not hard to show that

$$\frac{\hat{x} - x\|}{\|\hat{x}\|} \leqslant \kappa(A) \frac{\|r\|}{\|b\|}$$

Small ||r|| does not guarantee a small $||\hat{x} - x||$. If $\kappa(A)$ is large the \hat{x} returned by Gaussian elimination and back substitution (or any other solution method) is *not* guaranteed to be anywhere near the true solution to Ax = b.

- Applying Gaussian elimination with partial pivoting and back substitution to Ax = b yields a numerical solution \hat{x} such that the residual vector $r = b A\hat{x}$ is small *even if* the $\kappa(A)$ is large.
- If A and b are stored to machine precision ε_m , the numerical solution to Ax = b by any variant of Gaussian elimination is correct to d digits where

$$d = |\log_{10}(\varepsilon_m)| - \log_{10}(\kappa(A))$$

$$d = |\log_{10}(\varepsilon_m)| - \log_{10}(\kappa(A))$$

Example:

NUMPY computations have $\varepsilon_m \approx 2.2 \times 10^{-16}$. For a system with $\kappa(A) \sim 10^{10}$ the elements of the solution vector will have

$$d = |\log_{10}(2.2 \times 10^{-16})| - \log_{10} (10^{10})$$

= 16 - 10
= 6

correct (decimal) digits

Summary of Limits to Numerical Solution of Ax = b

- $\kappa(A)$ indicates how close *A* is to being numerically singular
- **2** If $\kappa(A)$ is "large", *A* is **ill-conditioned** and *even the best* numerical algorithms will produce a solution, \hat{x} that cannot be guaranteed to be close to the true solution, *x*
- In practice, Gaussian elimination with partial pivoting and back substitution produces a solution with a small residual

$$r = b - A\hat{x}$$

even if $\kappa(A)$ is large.

Consider the scalar equation

$$5x = 20 \qquad \Longrightarrow \qquad x = (5)^{-1}20$$

The extension to a system of equations is, of course

$$Ax = b \implies x = A^{-1}b$$

where $A^{-1}b$ is the formal solution to Ax = bIn numpy notation the system is solved with

x = np.linalg.solve(A,b)

Given an $n \times n$ matrix A, and an $n \times 1$ vector b the \ operator performs a sequence of tests on the A matrix. MATLAB attempts to solve the system with the method that gives the least roundoff and the fewest operations. When A is an $n \times n$ matrix the Python statement:

x = np.linalg.solve(A,b)

solves the linear system by using the LINPACK routines DGESV or ZGESVThese routines solve a dense linear system using gausian elimination with partial pivoting.