Numerical Linear Algebra for Computational Science and Information Engineering

> Lecture 08 Classical Iterative Methods for Linear Systems

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Towards iterative methods to solve linear systems

- So far, we saw how direct methods can be used to solve linear systems:
 Factorize the matrix (e.g., LU or Cholesky factorization) with cost O(n³)
 - **② Solve** the system using the computed factors with cost $\mathcal{O}(n^2)$
- Direct methods are very stable and accurate.

However, they can have a very high computational cost.

In general, direct methods are not suitable for very large n.

As a potentially more efficient way to solve linear systems, we will explore iterative methods.

Iterative methods are **inexact** in the sense that they rely on the generation of a sequence of approximate solutions which **converges towards the solution**, but the sequence is stopped at finite accuracy.

In general, iterative methods do not require explicit access to the matrix and they rely on the matrix-vector kernel $x \mapsto Ax$.

Iterative methods are particularly **recommended** for cases **where** the **matrix-vector product can be efficiently deployed**, e.g., as it is the case for sparse matrices.

Splitting methods Section 7.1 in Darve & Wootters (2021)

General splitting methods

- > Splitting methods are simple iterative methods to solve linear systems.
- Consider the A = M N splitting of a matrix A, where M is non-singular.
- The linear system Ax = b can be recast as follows:

$$Mx - Nx = b$$
$$x - M^{-1}Nx = M^{-1}b$$
$$x = Gx + M^{-1}b$$

so that x is a fixed point of $f: x \mapsto M^{-1}Nx + M^{-1}b$, i.e., x = f(x), and where $G := M^{-1}N$ is the iteration matrix.

▶ To solve a fixed point problem, one can start with any point x, and compute f(x). Then compute f(f(x)), then f(f(f(x))) and so on, until the sequence converges. In particular, we consider the following

Splitting method update rule

Given a matrix A=M-N where M is non-singular, the update rule for a general splitting method with a given $x^{(0)}$ is

$$x^{(k+1)} := G x^{(k)} + M^{-1} b \quad \text{where} \ \ G := M^{-1} N.$$

General splitting methods, cont'd₁

▶ The error $e^{(k+1)} := x^{(k+1)} - x$ is such that $e^{(k+1)} = Gx^{(k)} + M^{-1}b - Gx - M^{-1}b$ $= Gx^{(k)} - Gx$ $= Ge^{(k)}$

so that $e^{(k)} = G^k e^{(0)}$.

• The convergence theory depends on the iteration matrix $G = M^{-1}N$:

Theorem (Convergence of splitting methods)

Given b and A = M - N with non-singular A and M, the iteration

 $x^{(k+1)} = Gx^{(k)} + M^{-1}b$ where $G := M^{-1}N$

converges for any starting $\boldsymbol{x}^{(0)}$ if and only if

 $\rho(G) < 1$

where $\rho(G)$ is the spectral radius, i.e., the largest modulus of eigenvalue of the iteration matrix G.

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General splitting methods, cont'd₂

- Even though analyzing the spectrum of the iteration matrix G is generally difficult, it is understood that, the smaller the modulus ρ(G), the faster the convergence.
- How should we pick M and N? The selection of M and N may be guided by two desirable properties:
 - Linear systems of the form Mz = d are easy to solve.
 This suggest good choices for M are diagonal or triangular.
 - The spectral radius $\rho(G)$ is less than 1.
- We will see several examples of splitting methods, namely
 - Jacobi method
 - Gauss-Seidel method
 - Over-relaxation method

Jacobi method Section 7.2 in Darve & Wootters (2021)

Jacobi method

- Let A = D L U where
 - D is diagonal
 - L is strictly lower-triangular, i.e., with zeros on the diagonal
 - $\bullet \ U$ is strictly upper-triangular

Then, the splitting is clearly unique given A as we choose M=D and $N=L+U\colon$

$$A = M - M$$

The Jacobi splitting leads to the following iteration

Jacobi iterations

Suppose A = D - U - L as above. The update formula for Jacobi iteration is given by

$$Dx^{(k+1)} = (L+U)x^{(k)} + b.$$

Jacobi method, cont'd

The convergence of Jacobi iterations is as follows:

Theorem (Convergence of Jacobi iterations)

If A is strictly diagonally dominant, i.e.,

$$|a_{ii}| > \sum_{j \neq i} |a_{ij}|$$

for i = 1, 2..., n, then Jacobi iterations converge for any initial guess $x^{(0)}$.

Note that this condition is not necessary to ensure convergence.

The necessary condition to ensure convergence remains that the iteration matrix $G_{\text{Jacobi}} = D^{-1}(L+U)$ has a spectral radius smaller than one.

► The Jacobi method is especially simple to implement.

It is also well-suited for parallel implementation as we have

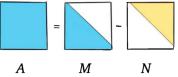
$$x_i^{(k+1)} = \left(b_i + (L+U)[i,:]x^{(k)}\right) / d_{ii}$$
$$x_i^{(k+1)} = \left(b_i - (A-D)[i,:]x^{(k)}\right) / d_{ii}.$$

Gauss-Seidel method Section 7.3 in Darve & Wootters (2021)

Gauss-Seidel method

- Let A = D L U where
 - D is diagonal
 - L is strictly lower-triangular, i.e., with zeros on the diagonal
 - $\bullet \ U$ is strictly upper-triangular

Then, the splitting is clearly unique given A as we choose M=D-L and $N=U\colon$



The Gauss-Seidel splitting leads to the following iteration

Gauss-Seidel iterations

Suppose A = D - U - L as above. The update formula for Gauss-Seidel iteration is given by

$$(D-L)x^{(k+1)} = Ux^{(k)} + b.$$

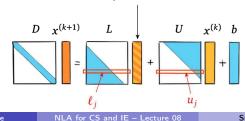
Gauss-Seidel method, $cont'd_1$

- ▶ Intuitively, "putting more information in M" should help with convergence of the method, and this is indeed the case, i.e., $\rho(G_{GS}) = \rho(G_{Jacobi})^2$.
- ▶ On the other hand, solving triangular systems with M = D − L is more involved than solving diagonal systems with M = D.
- We can compare Gauss-Seidel to Jacobi iterations as follows:

$$\begin{aligned} (D-L)x^{(k+1)} &= Ux^{(k)} + b \\ Dx^{(k+1)} &= Lx^{(k+1)} + Ux^{(k)} + b \text{ (Gauss-Seidel)} \\ Dx^{(k+1)} &= Lx^{(k)} + Ux^{(k)} + b \quad \text{(Jacobi)} \end{aligned}$$

 $x^{(k)}$ for Jacobi: $x^{(k+1)}$ for Gauss-Seidel.

and see they are very similar except for one term:



Gauss-Seidel method, $cont'd_2$

The convergence of Gauss-Seidel iterations is as follows:

Theorem (Convergence of Gauss-Seidel iterations) *If A*

- is strictly diagonally dominant, or
- is symmetric positive definite (SPD)

then Gauss-Seidel iterations converge irrespective of the initial guess $x^{(0)}$.

Note that these conditions are not necessary to ensure convergence.

The necessary condition to ensure convergence remains that the iteration matrix $G_{\rm GS} = (D-L)^{-1}U$ has a spectral radius smaller than one.

Successive over-relaxation Section 7.4 in Darve & Wootters (2021)

Successive over-relaxation

- Successive over-relaxation (SOR) consists of introducing a parameter to a splitting method in order to get a handle of the speed of convergence.
- ln particular, we use a parameter ω to boost convergence.

The idea is to start with the Gauss-Seidel update step as follows:

$$\begin{split} Dx_{\rm GS}^{(k+1)} &= Lx_{\rm GS}^{(k+1)} + Ux^{(k)} + b \\ x_{\rm GS}^{(k+1)} &= D^{-1} \left(Lx_{\rm GS}^{(k+1)} + Ux^{(k)} + b \right) \\ x_{\rm GS}^{(k+1)} &= x^{(k)} + \left[D^{-1} \left(Lx_{\rm GS}^{(k+1)} + Ux^{(k)} + b \right) - x^{(k)} \right] \\ \text{so that } x_{\rm GS}^{(k+1)} &= x^{(k)} + \Delta x_{\rm GS}^{(k)} \text{ where} \\ \Delta x_{\rm GS}^{(k)} &= D^{-1} \left(Lx_{\rm GS}^{(k+1)} + Ux^{(k)} + b \right) - x^{(k)} \end{split}$$

is the update to $x^{(k)}$ in a Gauss-Seidel iteration.

In SOR, the idea is to scale this correction by a parameter $0 < \omega < 2$:

$$x_{\text{SOR}}^{(k+1)} = x^{(k)} + \omega \Delta x_{\text{GS}}^{(k)}$$

Successive over-relaxation, cont'd

When ω ≈ 0, we are very cautious and only make small corrections to x^(k).
 When ω = 1, we recover a Gauss-Seidel iteration.
 When ω ≈ 2, we are very confident in the Gauss-Seidel correction and

When $\omega \approx 2$, we are very confident in the Gauss-Seidel correction and apply it twice instead of once.

The update formula of the SOR sequence is given as follows:

$$x_{\text{SOR}}^{(k+1)} = x_{\text{SOR}}^{(k)} + \omega \left[D^{-1} \left(Lx_{\text{SOR}}^{(k+1)} + Ux_{\text{SOR}}^{(k)} + b \right) - x_{\text{SOR}}^{(k)} \right]$$
$$= (1 - \omega) x_{\text{SOR}}^{(k)} + \omega \left[D^{-1} \left(Lx_{\text{SOR}}^{(k+1)} + Ux_{\text{SOR}}^{(k)} + b \right) \right]$$

which yields the following iterations:

SOR iterations

Let $\omega \in (0,2),$ and suppose A=D-L-U as above. The update formula for SOR iterations is

$$(D - \omega L)x_{\text{SOR}}^{(k+1)} = ((1 - \omega)D + \omega U)x_{\text{SOR}}^{(k)} + \omega b.$$

Successive over-relaxation, cont'd

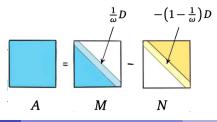
Since the update formula can written as

$$\left(\frac{1}{\omega}D - L\right)x_{\text{SOR}}^{(k+1)} = \left(\left(\frac{1}{\omega} - 1\right)D + U\right)x_{\text{SOR}}^{(k)} + b$$

and that we have

$$\left(\frac{1}{\omega}D - L\right) - \left(\left(\frac{1}{\omega} - 1\right)D + U\right) = D - L - U = A$$

we can say that SOR is a splitting method with $M = \frac{1}{\omega}D - L$ and $N = (\frac{1}{\omega} - 1)D + U$ such that A = M - N:



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Successive over-relaxation, cont'd

Although SOR is a heuristic, it can lead to significant improvements in the convergence rate when ω is chosen appropriately.

Theorem (Convergence of SOR iterations)

If A is symmetric positive definite (SPD), then SOR iterations converge irrespective of the initial guess $x^{(0)}$, for any $\omega \in (0, 2)$.

Note that this condition is not necessary to ensure convergence.

The necessary condition to ensure convergence remains that the iteration matrix

$$G = \left(\frac{1}{\omega}D - L\right)^{-1} \left(\left(\frac{1}{\omega} - 1\right)D + U\right)$$
$$= \left(D - \omega L\right)^{-1} \left(\left(1 - \omega\right)D + \omega U\right)$$

has a spectral radius smaller than one.

Homework problems

Homework problem

Turn in your own solution to the following problem:

Pb. 18 Let
$$A = \begin{bmatrix} 1 & 2 & 0 \\ 0 & 1 & 0 \\ 1 & 0 & 1 \end{bmatrix}$$
. Analyze the spectrum of the iteration matrix

and show whether

- (a) A Jacobi iteration would converge.
- (b) A Gauss-Seidel iteration would converge.
- (c) A SOR iteration would converge with $\omega=1/2.$