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# Computational Problem 2

## Table of Contents

Setup .....	1
EXERCISE 1 .....	1
(a) .....	1
(b) .....	1
(c) .....	2

## Setup

```
close all, clear all, clc
format long, format compact
fs = 16;
set(0, 'defaulttextfontsize', fs);
set(0, 'defaultaxesfontsize', fs);
```

## EXERCISE 1

### (a)

Since matrix  $A_\epsilon$  is a pentadiagonal matrix, so only main diagonal, and the first two upper and two lower diagonals exists nonzero elements.

Thus, for size  $n$ , each row has a maximum of 5 nonzero elements. Because if  $A_\epsilon$  is strictly diagonal dominant, it must satisfy  $|a_{ii}| > \sum_{j=1, j \neq i}^n |a_{ij}|$ , for  $i = 1, \dots, n$ .

So, matrix  $A_\epsilon$  is strictly diagonal dominant if  $1 > 2 \times (\epsilon^2 + \epsilon)$ . And we can get  $\epsilon < 0.366$ .

Therefore,  $\epsilon \in [0, 0.366)$ .

```
syms e
cond1 = e >= 0;
cond2 = e <= 1;
cond3 = 2*(e^2 + e) < 1;
conds = [cond1 cond2 cond3];

sol = solve(conds, e, 'ReturnConditions', true);
vpa(sol.conditions)

ans =
0.0 <= x & x < 0.36602540378443864676372317075294
```

### (b)

```
tol = 1e-10;
```

```

nmax = 1000;
x0 = [0; 0; 0; 0; 0];
[A, b] = matrix(5, 0.3);

% Jacobi method
[x, niter, relresiter, xiter] = itermeth(A, b, x0, nmax, tol, 'J');
niter

% Gauss-Seidel method
[x, niter, relresiter, xiter] = itermeth(A,b,x0,nmax,tol,'G');
niter

    itermeth converged in 50 iterations.
niter =
    50
    itermeth converged in 14 iterations.
niter =
    14

```

For Jacobi method, it needs 50 iterations to convergence; and for Gauss-Seidel method, it requires 14 iterations to convergence.

## (c)

Plot the spectral radius against the value of  $\epsilon$  for both methods

```

x=linspace(0, 1, 101);
figure
grid on
xlabel('epsilon')
ylabel('spectral radius')
hold on

radius_BJ = zeros(size(101));
radius_BGS = zeros(size(101));

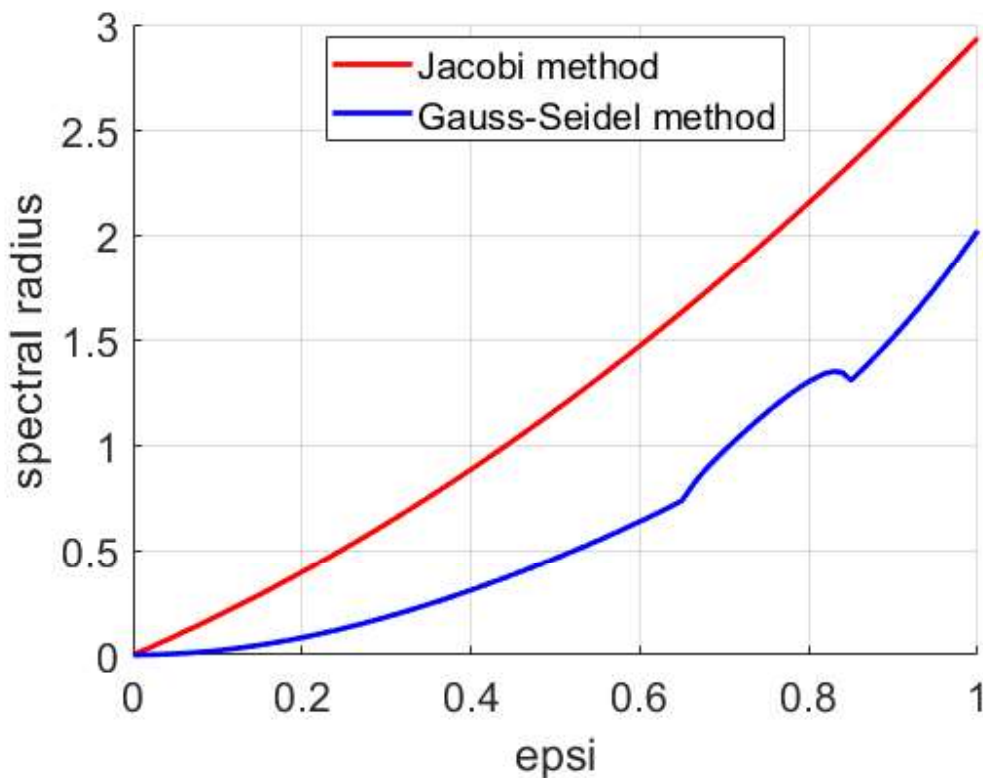
for e = 0:0.01:1
    [A, b] = matrix(5, e);
    D = diag(diag(A)); % diagonal part of A_epsilon
    L = tril(A,-1); % lower triangular part of A_epsilon
    U = triu(A,1); % lower triangular part of A_epsilon

    B_J = -(D^(-1)) * (L + U);
    B_GS = -(D + L)^(-1) * U;

    radius_BJ(round(e*100) + 1) = max(abs(eig(B_J))); % spectral radius is
    the maximum modulus of eigenvalues
    radius_BGS(round(e*100) + 1) = max(abs(eig(B_GS)));
end

plot(x, radius_BJ, 'r','LineWidth', 2);
plot(x, radius_BGS, 'blue','LineWidth', 2);
legend({'Jacobi method', 'Gauss-Seidel method'}, 'Location', 'Best');

```



Since the Jacobi method and the Gauss-Seidel method converge if the spectral radius of both  $B_J$  and  $B_{GS}$  less than 1. Therefore, according to the graph, when  $x \in [0, 0.44]$ , both  $B_J$  and  $B_{GS}$  less than 1, and this result is 0.074 larger than the answer of 0.366 from question (a).

According to the graph, if both methods converge, the Gauss-Seidel method is faster than the Jacobi method, because its spectral radius is smaller than the Jacobi method.

When  $n = 5$  and  $\epsilon = 0.5$ , I recommend using the Gauss-Seidel method.

```
[A, b] = matrix(5, 0.5);
D = diag(diag(A)); % diagonal part of A_epsilon
L = tril(A, -1); % lower triangular part of A_epsilon
U = triu(A, 1); % lower triangular part of A_epsilon

B_J = -(D^(-1)) * (L + U);
B_GS = -(D + L)^(-1) * U;

radius_BJ = max(abs(eig(B_J))) % spectral radius is the maximum modulus of
eigenvalues
radius_BGS = max(abs(eig(B_GS)))

tol = 1e-10;
nmax = 1000;
x0 = [0; 0; 0; 0; 0];
[A, b] = matrix(5, 0.5);
```

```
% Jacobi method
[x, niter, relresiter, xiter] = itermeth(A, b, x0, nmax, tol, 'J');

% Gauss-Seidel method
[x, niter, relresiter, xiter] = itermeth(A,b,x0,nmax,tol,'G');

radius_BJ =
    1.163568373244976
radius_BGS =
    0.460186129658958
itermeth reached the maximum iteration number without converging.
itermeth converged in 27 iterations.
```

Since spectral radius of  $B_J$  larger than 1, the Jacobi method will not convergence. Therefore, it should use the Gauss-Seidel method When  $n = 5$  and  $\epsilon = 0.5$ .

*Published with MATLAB® R2022b*