# ME630A

# Assignment-2

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Assignment - 2

Given poisson equation

as 
$$3^{2}\phi + 3^{2}\phi = S\phi$$
 for  $x \in [0,1]$ 
 $y \in [0,1]$ 

coshore  $S\phi = 2 \sinh[10(x-\frac{1}{2})] + 40(x-\frac{1}{2}) \cosh[10(x-\frac{1}{2})]$ 
 $+ 100(x-\frac{1}{2})^{2} \sinh[10(x-\frac{1}{2})] + 40(y-\frac{1}{2}) \cosh[10(y-\frac{1}{2})]$ 
 $+ 2 \sinh[10(y-\frac{1}{2})] + 40(y-\frac{1}{2}) \cosh[10(y-\frac{1}{2})]$ 
 $+ 4(x^{2}+y^{2})e^{2xy}$ 

Given boundary conditions as

 $\phi(0,y) = \frac{1}{4} \sinh(-5) + (y-\frac{1}{2})^{2} \sinh[10(y-\frac{1}{2})] + 1$ 
 $\phi(1,y) = \frac{1}{4} \sinh(-5) + (x-\frac{1}{2})^{2} \sinh[10(y-\frac{1}{2})] + e^{2y}$ 
 $\phi(x,0) = \frac{1}{4} \sinh(-5) + (x-\frac{1}{2})^{2} \sinh[10(x-\frac{1}{2})] + e^{2y}$ 
 $\phi(x,1) = \frac{1}{4} \sinh(5) + (x-\frac{1}{2})^{2} \sinh[10(x-\frac{1}{2})] + e^{2y}$ 

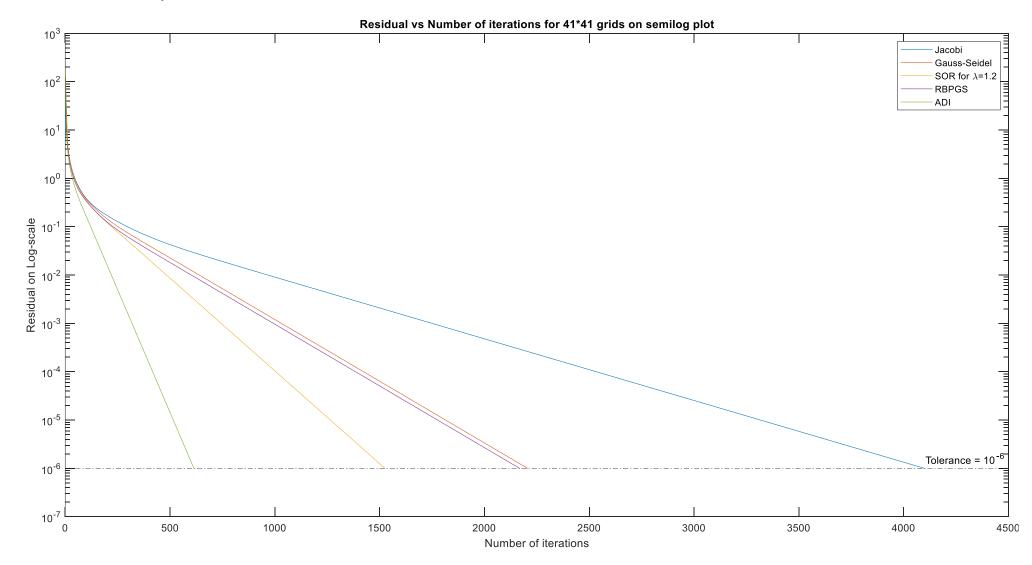
Also given analytical solution φ(x,y)= (x-1/2)<sup>2</sup> sinh [10(x-1/2)] + (y-1/2)<sup>2</sup> sinh [10(y-1/2)] + e<sup>2xy</sup>. We need to find number inumerical solution of above poisson equation using 1. Jacobi 2. Gauss - Seidel 3. SOR for different value of 2 especially 2=1.2 4. RBPGS 5. ADI using 41×41 Ggaids.

and 81×81 Ggaids.

All the above numerical solution technique is implemented through MATLAB code. Please find the code (.m file) attached with this support.

#### Plots for 41\*41 grids

#### 1. Residual plots for different iterative methods



After plotting residual for various iterative method vs number of iterations took for convergence we can make following observation. (Note: Tolerance =  $10^{-6}$ ).

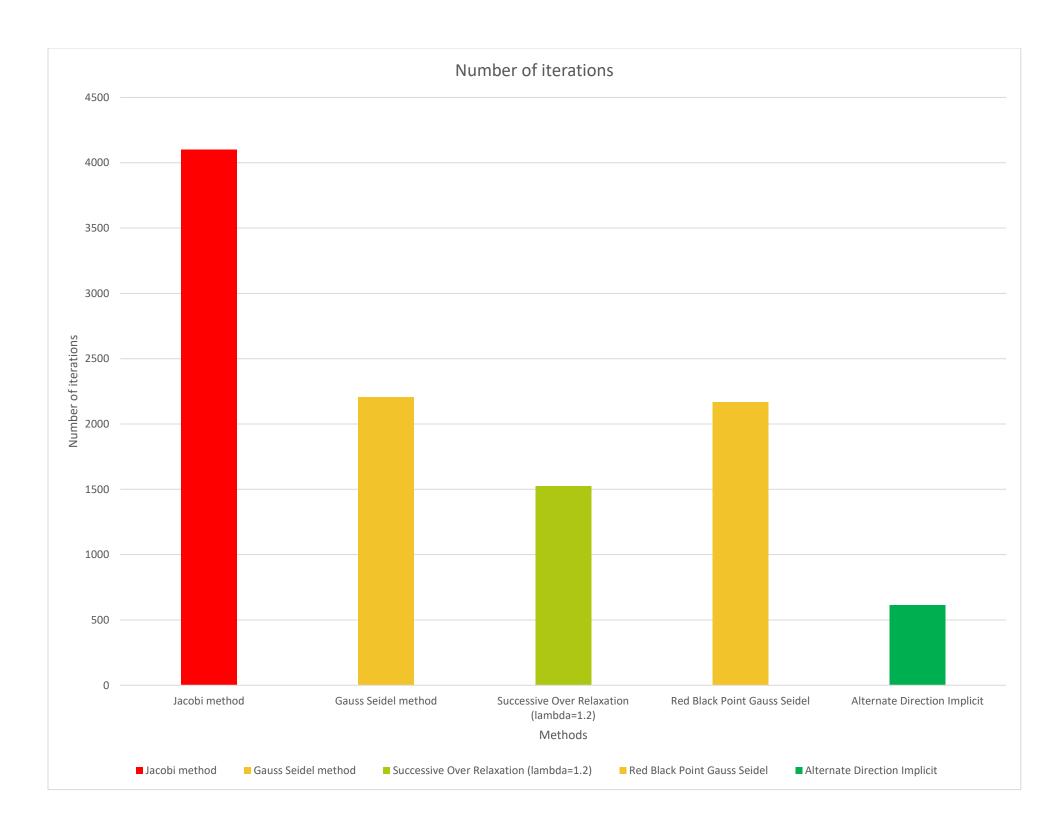
- Jacobi method took 4101 iterations to converge to the tolerance value.
- Gauss Seidel method took 2207 iterations to converge to the tolerance value.
- SOR method with  $\lambda$ =1.2 took 1526 iterations to converge to the tolerance value.
- RBPGS method took 2169 iterations to converge to the tolerance value.
- ADI method took 615 iterations to converge to the tolerance value.

It was also observed during analysis that for  $\lambda$ =1.76 for SOR the Convergence rate increase further.

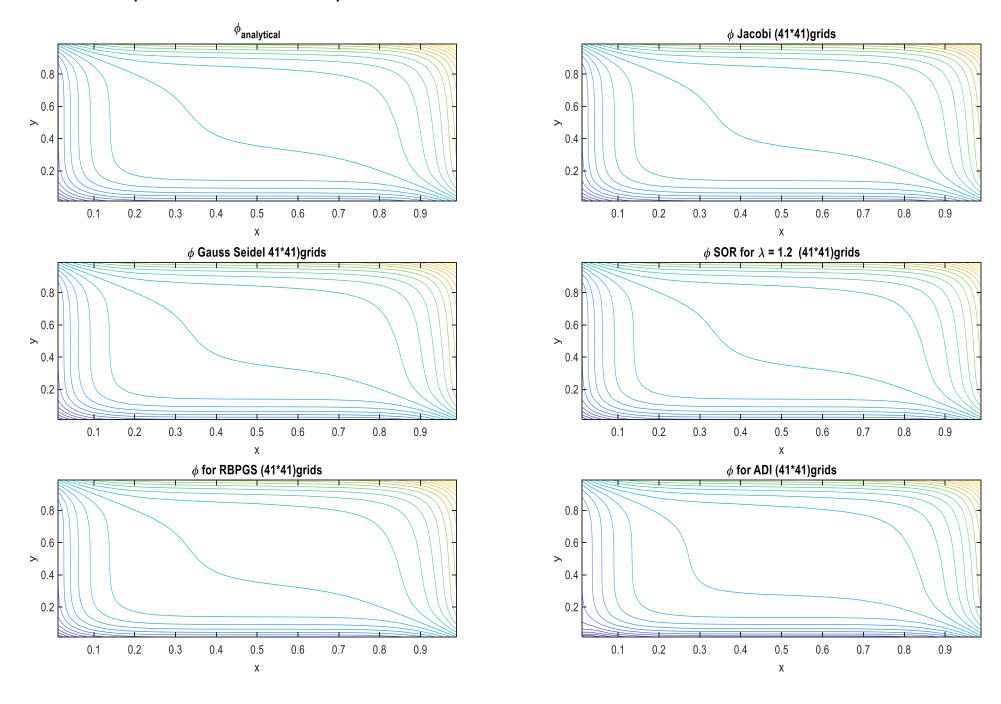
The rate of convergence is slowest for the Jacobi method as it is an explicit method and uses all the  $n^{th}$  value of  $\Phi$  for computing the  $(n+1)^{th}$  value of  $\Phi$ . This is improved by Gauss Seidel method by taking  $(n+1)^{th}$  value of  $\Phi$  (i.e., for (i-1, j), (i, j-1)) which has been computed for that iteration. The Successive over relaxation method further accelerates the rate of convergence by introducing  $\lambda$  ( $0 < \lambda < 2$ ). Here  $\lambda = 1.2$  is used. The Red Black Point Gauss Seidel enable parallel computing. However, here it is implemented with simple MATLAB code. It is observed that its rate of convergence is of the order of Gauss Seidel method but a little less than it. We note that Alterative Direction Implicit method has the fastest convergence rate. This was expected as we solve for entire x-coordinate for some y and entire y-coordinate for some x in one go using TDMA, hence the effect of Boundary Condition propagates faster inside the domain. Therefore, it converges faster.

S.no	Method	Number of iterations	Number of times faster that Jacobi
1	Jacobi method	4101	1 time
2	Gauss Seidel method	2207	1.86 times
3	Successive Over Relaxation ( $\lambda$ =1.2)	1526	2.69 times
4	Red Black Point Gauss Seidel	2169	1.89 times
5	Alternate Direction Implicit	615	6.67 times

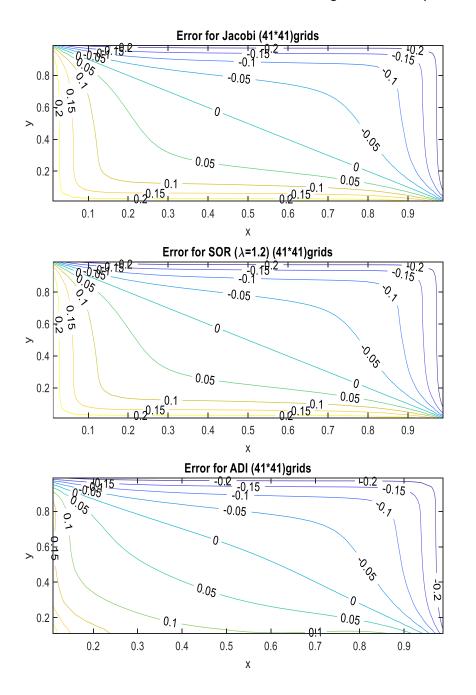
Hence, the rate of convergence is fastest for ADI and slowest for Jacobi method.

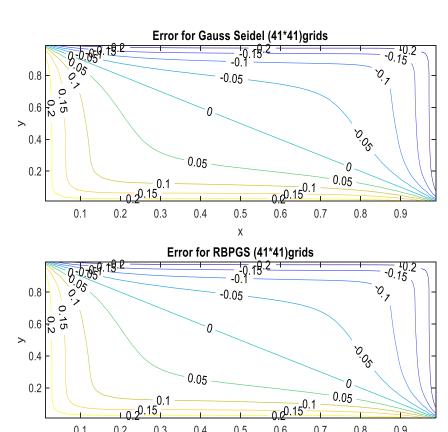


#### 2. Contour plots for $\Phi$ for analytical solution and different iterative methods



# 3. Contour plots of error ( $\Phi_{\text{converged}} - \Phi_{\text{analytical}}$ ) for different iterative methods





0.5

Χ

0.4

0.1

0.2

0.3

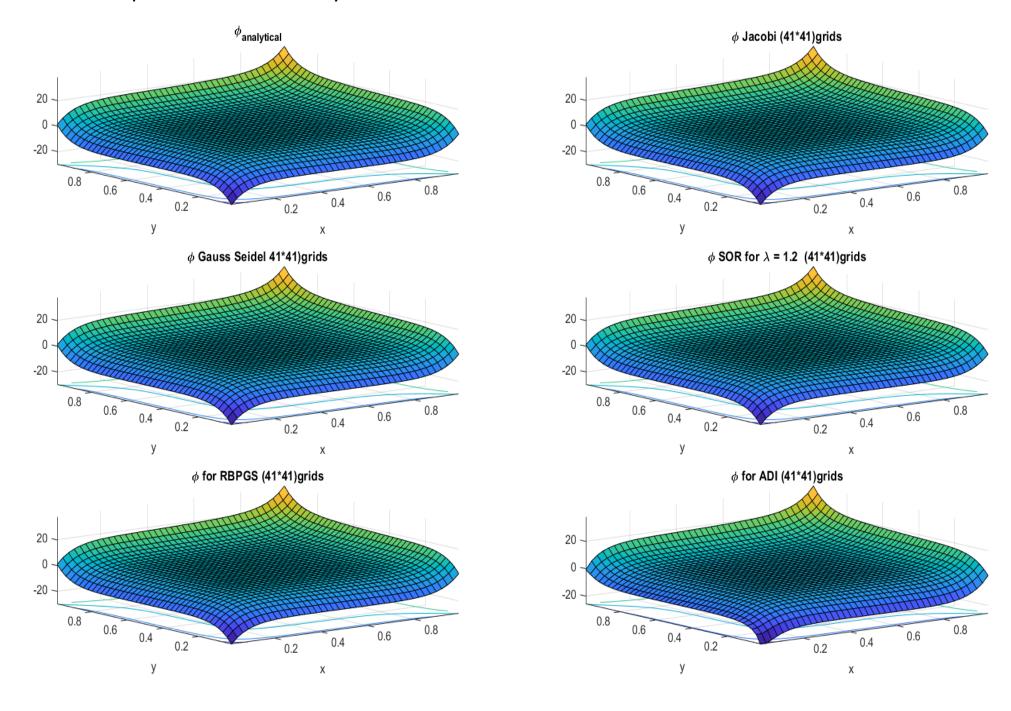
0.6

0.7

8.0

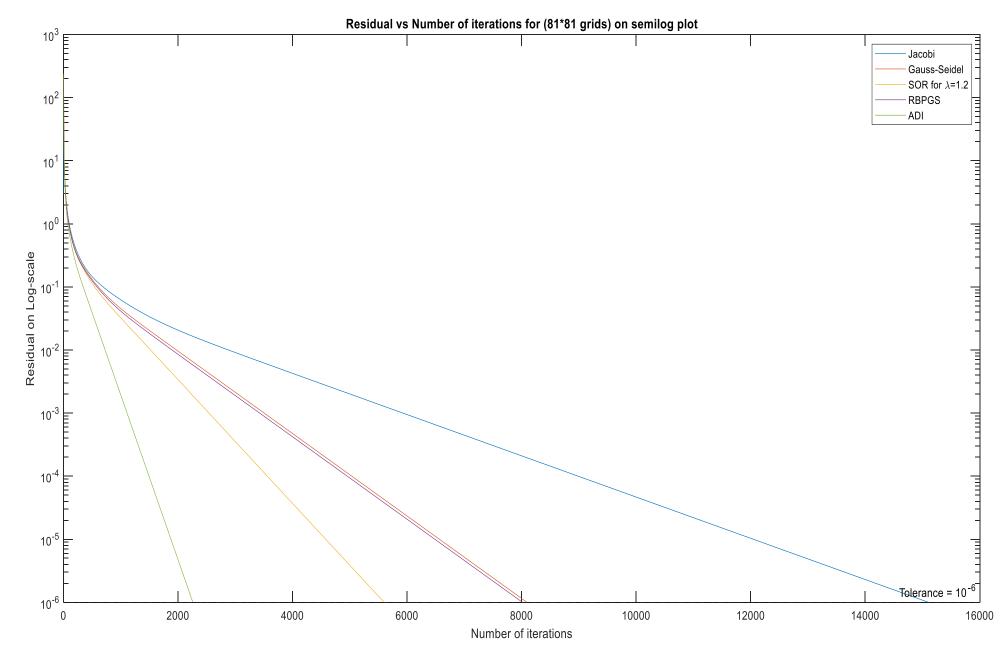
0.9

### 4. Surface plots for $\boldsymbol{\Phi}$ $\,$ for analytical solution and different iterative methods



#### Plots for 81\*81 grids

1. Residual plots for different iterative methods



After plotting residual for various iterative method vs number of iterations took for convergence we can make following observation. (Note: Tolerance =  $10^{-6}$ ).

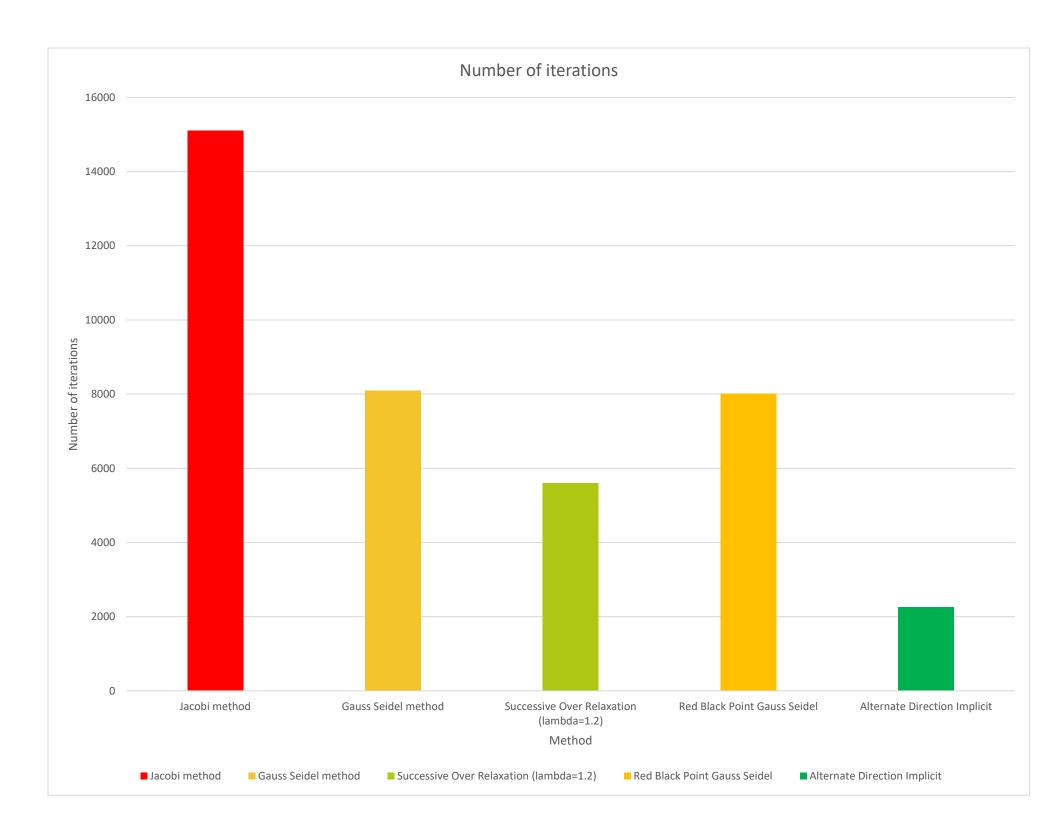
- Jacobi method took 15107 iterations to converge to the tolerance value.
- Gauss Seidel method took 8093 iterations to converge to the tolerance value.
- SOR method with  $\lambda$ =1.2 took 5598 iterations to converge to the tolerance value.
- RBPGS method took 8014 iterations to converge to the tolerance value.
- ADI method took 2258 iterations to converge to the tolerance value.

It was also observed during analysis that for  $\lambda$ =1.76 for SOR the Convergence rate increase further.

We can make following observations and comments from residual plots. The rate of convergence is slowest for the Jacobi method as it is an explicit method and uses all the n<sup>th</sup> value of  $\Phi$  for computing the (n+1)<sup>th</sup> value of  $\Phi$ . This is improved by Gauss Seidel method by taking (n+1)<sup>th</sup> value of  $\Phi$  (i.e., for (i-1, j), (i, j-1)) which has been computed for that iteration. The Successive over relaxation method further accelerates the rate of convergence by introducing  $\lambda$  (0 <  $\lambda$  < 2). Here  $\lambda$ =1.2 is used. The Red Black Point Gauss Seidel enable parallel computing. However, here it is implemented with simple MATLAB code. It is observed that its rate of convergence is of the order of Gauss Seidel method but a little less than it. We note that Alterative Direction Implicit method has the fastest convergence rate. This is because as we solve for entire x-coordinate for some y and entire y-coordinate for some x in one go using TDMA, hence the effect of Boundary Condition propagates faster inside the domain. Therefore, it converges faster.

S.no	Method	Number of iterations	Number of times faster that Jacobi
1	Jacobi method	15107	1 time
2	Gauss Seidel method	8093	1.87 times
3	Successive Over Relaxation ( $\lambda$ =1.2)	5598	2.70 times
4	Red Black Point Gauss Seidel	8014	1.89 times
5	Alternate Direction Implicit	2258	6.69 times

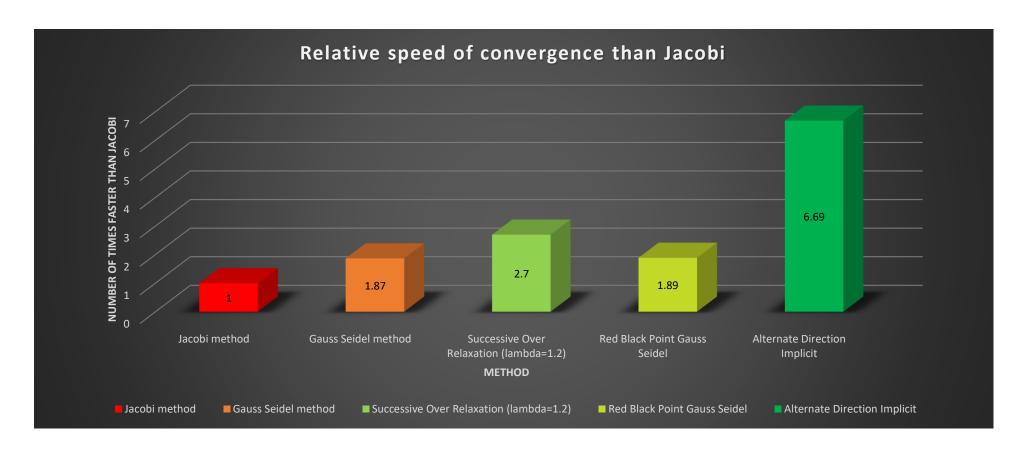
Hence, the rate of convergence is fastest for ADI and slowest for Jacobi method.



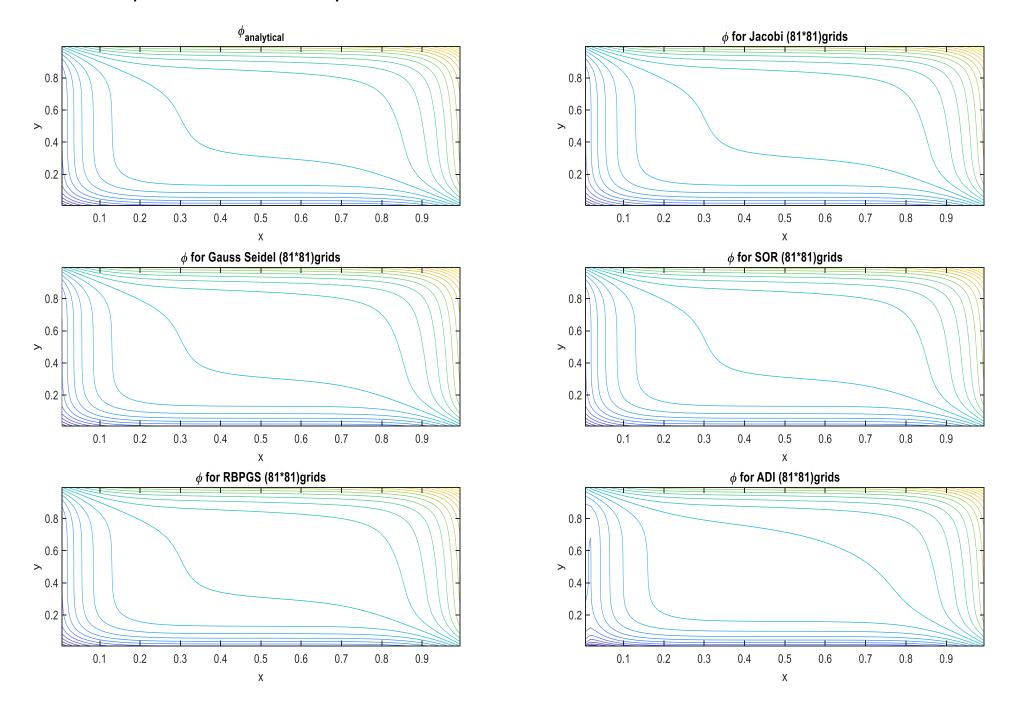
After looking at plots for residual for 41x41 grids and 81x81 grids we finally conclude that with increase in number of grids the number of iterations required for convergence for same tolerance value (here,  $10^{-6}$  )increases rapidly for all iterative methods used here. This is because we have to solve for  $\Phi$  at a greater number of nodes.

Finally, we conclude that

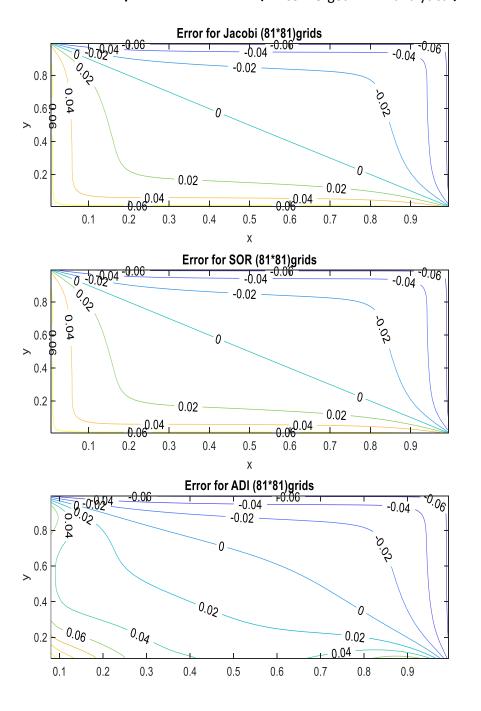
S.no	Method	Number of times faster than Jacobi
1	Jacobi method	1 time
2	Gauss Seidel method	1.87 times
3	Successive Over Relaxation ( $\lambda$ =1.2)	2.70 times
4	Red Black Point Gauss Seidel	1.89 times
5	Alternate Direction Implicit	6.69 times

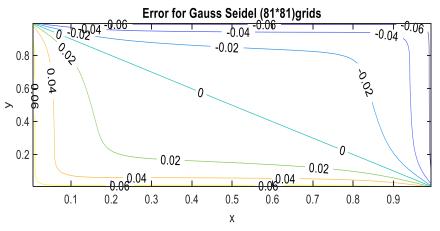


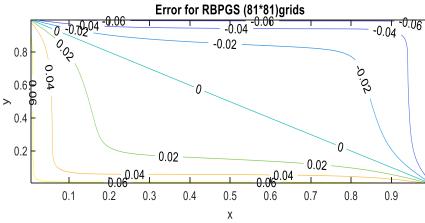
### 2.Contour plots for $\boldsymbol{\Phi}$ $\,$ for analytical solution and different iterative methods



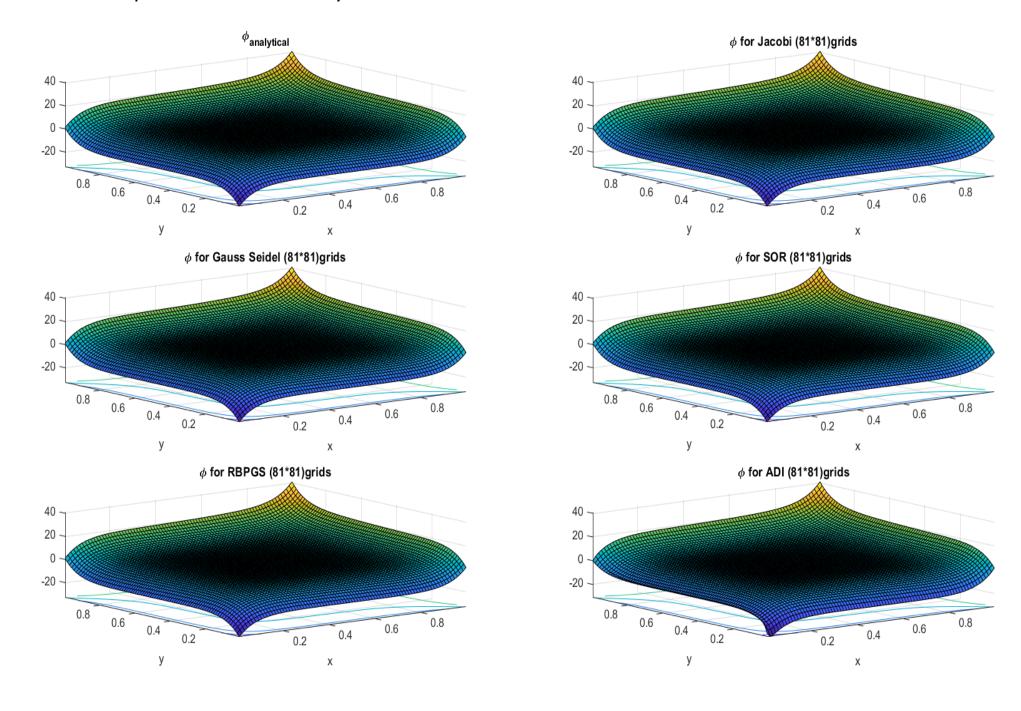
### 3.Contour plots of error ( $\Phi_{\text{converged}} - \Phi_{\text{analytical}}$ ) for different iterative methods







#### 4. Surface plots for $\Phi$ for analytical solution and different iterative methods



- 1. Please find the MATLAB code (\*.m files) for iterative methods and TDMA included with the submission.
- 2. Please find the plots also included with the submission.
- 3. Also find  $\Phi$  solutions and residual variable file (in \*.mat format) for both 41x41 grids and 81x81 grids generated after executing the code included with submission.
- 4. I am also including code (whose \*.m file in point 1 is included in submission) in this \*.pdf for your quick reference.

Thank You

Rikesh Sharma 180606

#### Algorithm for implementing different iterative methods

```
%Rikesh Sharma
%180606
%ME630A (CFD) | < submission date: 22/09/2021 >
function poisson solution()
%For 41*41 GRIDS
%Initialization of domain, grids, coeffients of equations and Boundary Conditions
domainX=1;
domainY=1;
Nx=41;
Ny=41;
nxp2=Nx+2;
nyp2=Ny+2;
deltaX=domainX/Nx;
deltaY=domainY/Ny;
pc = -2*(1/deltaX^2 + 1/deltaY^2);
px = 1/deltaX^2;
py = 1/deltaY^2;
x=(-deltaX/2 : deltaX : domainX+deltaX/2);
y=(-deltaY/2 : deltaY : domainY+deltaY/2);
bcx=zeros(2,nyp2); bcy=zeros(nxp2,2);
bcx(1,:)=0.25.*sinh(-5)+(y-0.5).^2.*sinh(10.*(y-0.5))+1;
bcx(2,:)=0.25.*sinh(5)+(y-0.5).^2.*sinh(10.*(y-0.5))+exp(2.*y);
```

```
bcy(:,1)=0.25.*sinh(-5)+(x-0.5).^2.*sinh(10.*(x-0.5))+1;
bcy(:,2)=0.25.*sinh(5)+(x-0.5).^2.*sinh(10.*(x-0.5))+exp(2.*x);
[x,y] = meshgrid(x,y);
phi=zeros(nxp2,nyp2); phi old=zeros(nxp2,nyp2);
Sphi=2.*sinh(10.*(x-0.5))+40.*(x-0.5).*cosh(10.*(x-0.5))+100.*((x-0.5).^2).*sinh(10.*(x-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5
(0.5) + 40.*(v-0.5).*cosh(10.*(v-0.5)) + 100.*((v-0.5).^2).*sinh(10.*(v-0.5)) + 4.*(x.^2+v.^2).*exp(2.*x.*v);
%Solution for 41*41 Grids%
residual Jacobi=zeros(2,1);
iter Jacobi=zeros(2,1);
maxError = 1e-6;
error=1:
iter=0;
%Jacobi method%
while(error>maxError && iter<100000)</pre>
         iter=iter+1;
         phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
         phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
         phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
         phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
         for i=2:Nx+1
                   for j=2:Ny+1
                            phi(i,j) = (Sphi(i,j) - (px*phi old(i-1,j) + px*phi old(i+1,j) + py*phi old(i,j-1) + py*phi old(i,j+1)))/pc;
                   end
         end
              disp(iter);
         errorMatrix=(phi(2:Nx+1,2:Ny+1)-phi_old(2:Nx+1,2:Ny+1));
          error = norm(errorMatrix,2);
          residual Jacobi(iter,1) = error;
         iter Jacobi(iter,1)=iter;
         phi old=phi;
용
               disp(error);
x=(deltaX/2 : deltaX : domainX-deltaX/2);
y=(deltaY/2 : deltaY : domainY-deltaY/2);
[x,y] = meshgrid(x,y);
phi analytic=(x-0.5).^2.*sinh(10.*(x-0.5))+(y-0.5).^2.*sinh(10.*(y-0.5))+exp(2.*x.*y);
phi Jacobi=phi(2:Nx+1,2:Ny+1);
phi=zeros(nxp2,nyp2); phi old=zeros(nxp2,nyp2);
residual Gauss=zeros(2,1);
iter Gauss=zeros(2,1);
```

```
error=1;
iter=0;
%Gauss Seidel method%
while(error>maxError && iter<100000)</pre>
    iter=iter+1;
    phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
    phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
    phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
    phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
    for i=2:Nx+1
        for j=2:Ny+1
            phi(i,j) = (Sphi(i,j) - (px*phi(i-1,j) + px*phi old(i+1,j) + py*phi(i,j-1) + py*phi old(i,j+1)))/pc;
        end
    end
응
      disp(iter);
    errorMatrix=(phi(2:Nx+1,2:Ny+1)-phi old(2:Nx+1,2:Ny+1));
    error = norm(errorMatrix,2);
    residual Gauss(iter,1) = error;
    iter Gauss(iter,1)=iter;
    phi old=phi;
      disp(error);
end
phi Gauss=phi(2:Nx+1,2:Ny+1);
phi=zeros(nxp2,nyp2); phi old=zeros(nxp2,nyp2);
residual SOR=zeros(2,1);
iter SOR=zeros(2,1);
lambda=1.2;
error=1;
iter=0;
%Succesive Over Relaxation method%
while(error>maxError && iter<100000)</pre>
    iter=iter+1;
    phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
    phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
    phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
    phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
    for i=2:Nx+1
        for j=2:Ny+1
```

```
phi(i,j) = (1-lambda) * phi(i,j) + lambda * (Sphi(i,j) - (px*phi(i-1,j) + px*phi old(i+1,j) + py*phi(i,j-1,j) + px*phi old(i+1,j) + py*phi(i,j-1,j) + px*phi old(i+1,j) + px*phi old(i+1
1) +py*phi old(i,j+1)))/pc;
                        end
            end
                  disp(iter);
            errorMatrix=(phi(2:Nx+1,2:Ny+1)-phi old(2:Nx+1,2:Ny+1));
            error = norm(errorMatrix,2);
            residual SOR(iter,1) = error;
            iter SOR(iter,1)=iter;
            phi old=phi;
                  disp(error);
end
phi SOR=phi(2:Nx+1,2:Ny+1);
phi=zeros(nxp2,nyp2); phi old=zeros(nxp2,nyp2);
residual RBPGS=zeros(2,1);
iter RBPGS=zeros(2,1);
error=1;
iter=0;
%RBPGS method%
while(error>maxError && iter<100000)</pre>
            iter=iter+1;
            phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
            phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
           phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
            phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
            for i=2:Nx+1
                       m = 2 + mod(i, 2);
                        for j=m:2:Ny+1
                                   phi(i,j) = (Sphi(i,j)-(px*phi old(i-1,j)+px*phi old(i+1,j)+py*phi old(i,j-1)+py*phi old(i,j+1)))/pc;
                        end
            end
            for i=2:Nx+1
                       m = 2 + mod(i+1, 2);
                        for j=m:2:Ny+1
                                   phi(i,j) = (Sphi(i,j) - (px*phi(i-1,j) + px*phi(i+1,j) + py*phi(i,j-1) + py*phi(i,j+1)))/pc;
                        end
            end
응
                  disp(iter);
```

```
errorMatrix=(phi(2:Nx+1,2:Ny+1)-phi old(2:Nx+1,2:Ny+1));
    error = norm(errorMatrix,2);
    residual RBPGS(iter,1)=error;
    iter RBPGS(iter,1)=iter;
    phi old=phi;
      disp(error);
end
phi RBPGS=phi(2:Nx+1,2:Ny+1);
%ADI method%
phi=zeros(nxp2,nyp2); phi old=zeros(nxp2,nyp2);
residual ADI=zeros(2,1);
iter ADI=zeros(2,1);
error=1;
iter=0;
phi old(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
phi old(nxp2, 2:Ny+1) = 2*bcx(2, 2:Ny+1) - phi(Nx+1, 2:Ny+1);
phi old(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
phi old(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
ar=zeros(Nx,1); ar(2:Nx,1)=-px;
br=zeros(Nx,1); br(:,1)=-pc;
cr=zeros(Nx,1); cr(2:Nx,1)=-px;
dr=zeros(Nx,1);
ac=zeros(Ny,1); ac(2:Ny,1)=-py;
bc=zeros(Ny,1); bc(:,1)=-pc;
cc=zeros(Ny,1); cc(2:Ny,1)=-py;
dc=zeros(Ny,1);
while(error>maxError && iter<100000)</pre>
    iter=iter+1;
    phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
    phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
    phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
    phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
    for j=2:Ny+1
        dr(1,1) = py.*phi(2,j-1) + py.*phi old(2,j+1) - Sphi(2,j) + px*phi(1,j);
        dr(2:Nx-1,1)=py.*phi(3:Nx,j-1)+py.*phi old(3:Nx,j+1)-Sphi(3:Nx,j);
        dr(Nx,1) = py.*phi(Nx+1,j-1) + py.*phi old(Nx+1,j+1) - Sphi(Nx+1,j) + px*phi(Nx+2,j);
        phi(2:Nx+1,j)=TDMA(ar,br,cr,dr);
    end
    phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
```

```
phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
    phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
    phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
    for i=2:Nx+1
        dc(1,1) = px.*phi(i-1,2) + px.*phi(i+1,2) - Sphi(i,2) + py*phi(i,1);
        dc(2:Ny-1,1)=px.*phi(i-1,3:Ny)+px.*phi(i+1,3:Ny)-Sphi(i,3:Ny);
        dc(Ny,1)=px.*phi(i-1,Ny+1)+px.*phi(i+1,Ny+1)-Sphi(i,Ny+1) + py*phi(i,Ny+2);
        phi(i,2:Ny+1) = TDMA(ac,bc,cc,dc);
    end
    phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
    phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
    phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
    phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
응
      disp(iter);
    errorMatrix=(phi(2:Nx+1,2:Ny+1)-phi old(2:Nx+1,2:Ny+1));
    error = norm(errorMatrix,2);
    residual ADI(iter,1) = error;
    iter ADI(iter,1)=iter;
    phi old=phi;
      disp(error);
end
phi ADI=phi(2:Nx+1,2:Ny+1);
%Plots for 41*41 Grids%
x=(deltaX/2 : deltaX : domainX-deltaX/2);
y=(deltaY/2 : deltaY : domainY-deltaY/2);
[x,y] = meshgrid(x,y);
%Plotting phi contours
figure
subplot(3,2,1)
contour(x,y,phi analytic,25);
xlabel('x');
ylabel('y');
title('{\phi} {analytical}');
subplot(3,2,2)
contour(x,y,phi Jacobi,25);
xlabel('x');
ylabel('y');
title('{\phi} Jacobi (41*41)grids');
subplot(3,2,3)
contour(x,y,phi Gauss,25);
xlabel('x');
```

```
ylabel('y');
title('{\phi} Gauss Seidel 41*41)grids');
subplot(3,2,4)
contour(x,y,phi SOR,25);
xlabel('x');
ylabel('y');
title('\{\phi\}\ SOR for \lambda = 1.2 (41*41)grids');
subplot(3,2,5)
contour(x,y,phi RBPGS,25);
xlabel('x');
ylabel('y');
title('{\phi} for RBPGS (41*41)grids');
subplot(3,2,6)
contour(x,y,phi ADI,25);
xlabel('x');
ylabel('y');
title('{\phi} for ADI (41*41)grids');
%Plotting phi surfaces
figure
subplot(3,2,1)
surfc(x,y,phi analytic);
xlabel('x');
ylabel('y');
title('{\phi} {analytical}');
subplot(3,2,2)
surfc(x,y,phi Jacobi);
xlabel('x');
ylabel('y');
title('{\phi} Jacobi (41*41)grids');
subplot(3,2,3)
surfc(x,y,phi Gauss);
xlabel('x');
ylabel('y');
title('{\phi} Gauss Seidel 41*41)grids');
subplot(3,2,4)
surfc(x,y,phi SOR);
xlabel('x');
ylabel('y');
title('\{\phi\}\ SOR for \lambda = 1.2 (41*41)grids');
subplot(3,2,5)
surfc(x,y,phi RBPGS);
xlabel('x');
ylabel('y');
title('{\phi} for RBPGS (41*41)grids');
subplot(3,2,6)
surfc(x,y,phi ADI);
```

```
xlabel('x');
ylabel('y');
title('{\phi} for ADI (41*41)grids');
%Plotting Errors
figure
subplot(3,2,1)
contour(x,y,phi Jacobi-phi analytic,'ShowText','on');
xlabel('x');
vlabel('v');
title('Error for Jacobi (41*41)grids');
subplot(3,2,2)
contour(x,y,phi Gauss-phi analytic,'ShowText','on');
xlabel('x');
ylabel('y');
title('Error for Gauss Seidel (41*41)grids');
subplot(3,2,3)
contour(x,y,phi SOR-phi analytic,'ShowText','on');
xlabel('x');
ylabel('y');
title('Error for SOR (\lambda=1.2) (41*41)grids');
subplot(3,2,4)
contour(x,y,phi RBPGS-phi analytic,'ShowText','on');
xlabel('x');
vlabel('v');
title('Error for RBPGS (41*41)grids');
subplot(3,2,5)
x=(9*deltaX/2 : deltaX : domainX-deltaX/2);
y=(9*deltaY/2 : deltaY : domainY-deltaY/2);
[x,y] = meshgrid(x,y);
contour(x,y,phi ADI(5:Nx,5:Ny)-phi analytic(5:Nx,5:Ny),'ShowText','on');
xlabel('x');
ylabel('y');
title('Error for ADI (41*41)grids');
%Ploting residuals for rate of convergence
figure
semilogy(iter Jacobi , residual Jacobi);
hold on
semilogy(iter Gauss , residual Gauss);
semilogy(iter SOR , residual SOR);
semilogy(iter RBPGS , residual RBPGS);
                    , residual ADI);
semilogy(iter ADI
yline(1e-6,'-.k','Tolerance = \overline{10}^{-6}');
legend('Jacobi','Gauss-Seidel','SOR for \lambda=1.2','RBPGS','ADI');
xlabel('Number of iterations');
ylabel('Residual on Log-scale');
```

```
title('Residual vs Number of iterations for 41*41 grids on semilog plot');
save('phi41.mat','phi analytic','phi Jacobi','phi Gauss','phi SOR','phi RBPGS','phi ADI');
save('residual41.mat', 'residual Jacobi', 'residual Gauss', 'residual SOR', 'residual RBPGS', 'residual ADI');
%Solution for 41*41 Grids Ends%
%For 81*81 GRIDS
%Initialization of domain, grids, coefficients of equations and Boundary Conditions
domainX=1;
domainY=1;
Nx=81;
Ny = 81;
nxp2=Nx+2;
nyp2=Ny+2;
deltaX=domainX/Nx;
deltaY=domainY/Nv;
pc = -2*(1/deltaX^2 + 1/deltaY^2);
px = 1/deltaX^2;
py = 1/deltaY^2;
x=(-deltaX/2 : deltaX : domainX+deltaX/2);
y=(-deltaY/2 : deltaY : domainY+deltaY/2);
bcx=zeros(2,nyp2); bcy=zeros(nxp2,2);
bcx(1,:)=0.25.*sinh(-5)+(y-0.5).^2.*sinh(10.*(y-0.5))+1;
bcx(2,:)=0.25.*sinh(5)+(y-0.5).^2.*sinh(10.*(y-0.5))+exp(2.*y);
bcv(:,1)=0.25.*sinh(-5)+(x-0.5).^2.*sinh(10.*(x-0.5))+1;
bcv(:,2)=0.25.*sinh(5)+(x-0.5).^2.*sinh(10.*(x-0.5))+exp(2.*x);
[x,y] = meshgrid(x,y);
phi=zeros(nxp2,nyp2); phi old=zeros(nxp2,nyp2);
Sphi=2.*sinh(10.*(x-0.5))+40.*(x-0.5).*cosh(10.*(x-0.5))+100.*((x-0.5).^2).*sinh(10.*(x-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5))+2.*sinh(10.*(y-0.5
(0.5) + 40.*(y-0.5).*\cosh(10.*(y-0.5)) + 100.*((y-0.5).^2).*\sinh(10.*(y-0.5)) + 4.*(x.^2+y.^2).*\exp(2.*x.*y);
%Solution for 41*41 Grids%
residual Jacobi=zeros(2,1);
iter Jacobi=zeros(2,1);
maxError = 1e-6;
error=1;
iter=0;
%Jacobi method%
while(error>maxError && iter<100000)</pre>
        iter=iter+1:
         phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
         phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
         phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
         phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
```

```
for i=2:Nx+1
        for j=2:Ny+1
            phi(i,j) = (Sphi(i,j) - (px*phi old(i-1,j) + px*phi old(i+1,j) + py*phi old(i,j-1) + py*phi old(i,j+1)))/pc;
        end
    end
응
      disp(iter);
    errorMatrix=(phi(2:Nx+1,2:Ny+1)-phi old(2:Nx+1,2:Ny+1));
    error = norm(errorMatrix,2);
    residual Jacobi(iter,1) = error;
    iter Jacobi(iter,1)=iter;
    phi old=phi;
      disp(error);
end
x=(deltaX/2 : deltaX : domainX-deltaX/2);
v=(deltaY/2 : deltaY : domainY-deltaY/2);
[x,y] = meshgrid(x,y);
phi analytic=(x-0.5).^2.*sinh(10.*(x-0.5))+(y-0.5).^2.*sinh(10.*(y-0.5))+exp(2.*x.*y);
phi Jacobi=phi(2:Nx+1,2:Ny+1);
phi=zeros(nxp2,nyp2); phi old=zeros(nxp2,nyp2);
residual Gauss=zeros(2,1);
iter Gauss=zeros(2,1);
error=1;
iter=0;
%Gauss Seidel method%
while(error>maxError && iter<100000)</pre>
    iter=iter+1;
    phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
    phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
    phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
    phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
    for i=2:Nx+1
        for j=2:Ny+1
            phi(i,j) = (Sphi(i,j) - (px*phi(i-1,j) + px*phi old(i+1,j) + py*phi(i,j-1) + py*phi old(i,j+1)))/pc;
        end
    end
      disp(iter);
    errorMatrix=(phi(2:Nx+1,2:Ny+1)-phi old(2:Nx+1,2:Ny+1));
    error = norm(errorMatrix,2);
    residual Gauss(iter,1)=error;
    iter Gauss(iter,1)=iter;
    phi old=phi;
```

```
disp(error);
end
phi Gauss=phi(2:Nx+1,2:Ny+1);
phi=zeros(nxp2,nyp2); phi old=zeros(nxp2,nyp2);
residual SOR=zeros(2,1);
iter SOR=zeros(2,1);
lambda=1.2;
error=1;
iter=0;
%Succesive Over Relaxation method%
while(error>maxError && iter<100000)</pre>
            iter=iter+1;
            phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
            phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
            phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
            phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
            for i=2:Nx+1
                        for j=2:Ny+1
                                   phi(i,j) = (1-lambda) * phi(i,j) + lambda * (Sphi(i,j) - (px*phi(i-1,j) + px*phi old(i+1,j) + py*phi(i,j-1,j) + px*phi old(i+1,j) + py*phi(i,j-1,j) + px*phi old(i+1,j) + px*phi old(i+1
1) +py*phi old(i,j+1)))/pc;
                        end
            end
응
                  disp(iter);
            errorMatrix=(phi(2:Nx+1,2:Ny+1)-phi old(2:Nx+1,2:Ny+1));
            error = norm(errorMatrix,2);
            residual SOR(iter,1) = error;
            iter SOR(iter,1)=iter;
            phi old=phi;
응
                  disp(error);
end
phi SOR=phi(2:Nx+1,2:Ny+1);
phi=zeros(nxp2,nyp2); phi old=zeros(nxp2,nyp2);
residual RBPGS=zeros(2,1);
iter RBPGS=zeros(2,1);
error=1;
iter=0;
%RBPGS method%
while(error>maxError && iter<100000)</pre>
            iter=iter+1;
```

```
phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
    phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
    phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
    phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
    for i=2:Nx+1
        m = 2 + mod(i, 2);
        for j=m:2:Ny+1
            phi(i,j) = (Sphi(i,j) - (px*phi old(i-1,j) + px*phi old(i+1,j) + py*phi old(i,j-1) + py*phi old(i,j+1)))/pc;
        end
    end
    for i=2:Nx+1
        m = 2 + mod(i+1, 2);
        for j=m:2:Ny+1
            phi(i,j) = (Sphi(i,j)-(px*phi(i-1,j)+px*phi(i+1,j)+py*phi(i,j-1)+py*phi(i,j+1)))/pc;
        end
    end
응
      disp(iter);
    errorMatrix=(phi(2:Nx+1,2:Ny+1)-phi old(2:Nx+1,2:Ny+1));
    error = norm(errorMatrix,2);
    residual RBPGS(iter,1) = error;
    iter RBPGS(iter,1)=iter;
    phi old=phi;
      disp(error);
end
phi RBPGS=phi(2:Nx+1,2:Ny+1);
%ADI method%
phi=zeros(nxp2,nyp2); phi old=zeros(nxp2,nyp2);
residual ADI=zeros(2,1);
iter ADI=zeros(2,1);
error=1;
iter=0;
phi old(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
phi old(nxp2, 2:Ny+1) = 2*bcx(2, 2:Ny+1) - phi(Nx+1, 2:Ny+1);
phi old(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
phi old(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
ar=zeros(Nx,1); ar(2:Nx,1)=-px;
br=zeros(Nx,1); br(:,1)=-pc;
cr=zeros(Nx,1); cr(2:Nx,1)=-px;
dr=zeros(Nx,1);
```

```
ac=zeros(Ny,1); ac(2:Ny,1)=-py;
bc=zeros(Ny,1); bc(:,1)=-pc;
cc=zeros(Ny,1); cc(2:Ny,1)=-py;
dc=zeros(Ny,1);
while(error>maxError && iter<100000)</pre>
    iter=iter+1;
    phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
    phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
    phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
    phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
    for j=2:Ny+1
        dr(1,1) = py.*phi(2,j-1) + py.*phi old(2,j+1) - Sphi(2,j) + px*phi(1,j);
        dr(2:Nx-1,1) = py.*phi(3:Nx,j-1) + py.*phi old(3:Nx,j+1) - Sphi(3:Nx,j);
        dr(Nx,1) = py.*phi(Nx+1,j-1) + py.*phi old(Nx+1,j+1) - Sphi(Nx+1,j) + px*phi(Nx+2,j);
        phi(2:Nx+1,j)=TDMA(ar,br,cr,dr);
    end
    phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
    phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
    phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
    phi(2:Nx+1,nvp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Nv+1);
    for i=2:Nx+1
        dc(1,1) = px.*phi(i-1,2) + px.*phi(i+1,2) - Sphi(i,2) + py*phi(i,1);
        dc(2:Ny-1,1)=px.*phi(i-1,3:Ny)+px.*phi(i+1,3:Ny)-Sphi(i,3:Ny);
        dc(Ny,1)=px.*phi(i-1,Ny+1)+px.*phi(i+1,Ny+1)-Sphi(i,Ny+1) + py*phi(i,Ny+2);
        phi(i, 2:Ny+1) = TDMA(ac, bc, cc, dc);
    end
    phi(1,2:Ny+1) = 2*bcx(1,2:Ny+1) - phi(2,2:Ny+1);
    phi(nxp2,2:Ny+1) = 2*bcx(2,2:Ny+1) - phi(Nx+1,2:Ny+1);
    phi(2:Nx+1,1) = 2*bcy(2:Nx+1,1) - phi(2:Nx+1,2);
    phi(2:Nx+1,nyp2) = 2*bcy(2:Nx+1,2) - phi(2:Nx+1,Ny+1);
응
      disp(iter);
    errorMatrix=(phi(2:Nx+1,2:Ny+1)-phi old(2:Nx+1,2:Ny+1));
    error = norm(errorMatrix,2);
    residual ADI(iter,1) = error;
    iter ADI(iter,1)=iter;
    phi old=phi;
      disp(error);
응
end
```

```
phi ADI=phi(2:Nx+1,2:Ny+1);
x=(deltaX/2 : deltaX : domainX-deltaX/2);
y=(deltaY/2 : deltaY : domainY-deltaY/2);
[x,y] = meshgrid(x,y);
%Plots for 81*81 Grids%
%Ploting phi contours
figure
subplot(3,2,1)
contour(x,y,phi analytic,25);
xlabel('x');
ylabel('y');
title('{\phi} {analytical}');
subplot(3,2,2)
contour(x,y,phi Jacobi,25);
xlabel('x');
ylabel('y');
title('{\phi} for Jacobi (81*81)grids');
subplot(3,2,3)
contour(x,y,phi Gauss,25);
xlabel('x');
ylabel('y');
title('{\phi} for Gauss Seidel (81*81)grids');
subplot(3,2,4)
contour(x,y,phi SOR,25);
xlabel('x');
ylabel('y');
title('{\phi} for SOR (81*81)grids');
subplot(3,2,5)
contour(x,y,phi RBPGS,25);
xlabel('x');
ylabel('y');
title('{\phi} for RBPGS (81*81)grids');
subplot(3,2,6)
contour(x,y,phi ADI,25);
xlabel('x');
ylabel('y');
title('{\phi} for ADI (81*81)grids');
%Ploting phi surfaces
figure
subplot(3,2,1)
surfc(x,y,phi analytic);
xlabel('x');
ylabel('y');
title('{\phi} {analytical}');
```

```
subplot(3,2,2)
surfc(x,y,phi Jacobi);
xlabel('x');
ylabel('y');
title('{\phi} for Jacobi (81*81)grids');
subplot(3,2,3)
surfc(x,y,phi Gauss);
xlabel('x');
ylabel('y');
title('{\phi} for Gauss Seidel (81*81)grids');
subplot(3,2,4)
surfc(x,y,phi SOR);
xlabel('x');
ylabel('y');
title('{\phi} for SOR (81*81)grids');
subplot(3,2,5)
surfc(x,y,phi RBPGS);
xlabel('x');
ylabel('y');
title('{\phi} for RBPGS (81*81)grids');
subplot(3,2,6)
surfc(x,y,phi ADI);
xlabel('x');
ylabel('y');
title('{\phi} for ADI (81*81)grids');
%Ploting Error contours
figure
subplot(3,2,1)
contour(x,y,phi Jacobi-phi analytic,'ShowText','on');
xlabel('x');
ylabel('y');
title('Error for Jacobi (81*81)grids');
subplot(3,2,2)
contour(x,y,phi Gauss-phi analytic,'ShowText','on');
xlabel('x');
ylabel('y');
title('Error for Gauss Seidel (81*81)grids');
subplot(3,2,3)
contour(x,y,phi SOR-phi analytic,'ShowText','on');
xlabel('x');
ylabel('y');
title('Error for SOR (81*81)grids');
subplot(3,2,4)
contour(x,y,phi RBPGS-phi analytic,'ShowText','on');
xlabel('x');
ylabel('y');
```

```
title('Error for RBPGS (81*81)grids');
subplot(3,2,5)
x=(13*deltaX/2 : deltaX : domainX-deltaX/2);
y=(13*deltaY/2 : deltaY : domainY-deltaY/2);
[x,y] = meshgrid(x,y);
contour(x,y,phi ADI(7:Nx,7:Ny)-phi analytic(7:Nx,7:Ny),'ShowText','on');
xlabel('x');
vlabel('v');
title('Error for ADI (81*81)grids');
%Ploting residuals
figure
semilogy(iter Jacobi ,residual Jacobi);
hold on
semilogy(iter Gauss ,residual Gauss);
semilogy(iter SOR , residual SOR);
semilogy(iter RBPGS , residual RBPGS);
semilogy(iter ADI , residual ADI);
yline (1e-6, '-.k', 'Tolerance = \overline{10^{-6}}');
legend('Jacobi','Gauss-Seidel','SOR for \lambda=1.2','RBPGS','ADI');
xlabel('Number of iterations');
ylabel('Residual on Log-scale');
title('Residual vs Number of iterations for (81*81 grids) on semilog plot');
save('phi81.mat','phi analytic','phi Jacobi','phi Gauss','phi SOR','phi RBPGS','phi ADI');
save('residual81.mat', 'residual Jacobi', 'residual Gauss', 'residual SOR', 'residual RBPGS', 'residual ADI');
%Solutions for (81*81) Grids ends%
```

#### Algorithm for TDMA

end

```
%Rikesh Sharma %180606 %ME630A (CFD) | < submission date: 22/09/2021 > function x = TDMA(a,b,c,d) %a, b, c are the column vectors for the compressed tridiagonal matrix, d is the right vector n = length(b); % n is the number of rows % Modify the first-row coefficients c(1) = c(1) / b(1); % Division by zero risk.
```