

Project 4

```
function [u,x,y,eps]=poisson(n,nIter,omega)
h=1/n;
x=linspace(0,1,n+1);
b=[];
for i=1:n
    for j=1:n
        b(i,j)=(h.^2)*fcn(x(i),x(j));
    end
end

u=zeros(n,n);
eps=[];
for k=1:nIter
    uold=u;
    for i=2:n-1
        for j=2:n-1
            u(i,j)=(1-omega)*u(i,j)+omega*((1/4)*(b(i,j)+u(i-1,j)+u(i+1,j)
+u(i,j-1)+u(i,j+1)));
            u(i,j)=u(i,j)*inGeom(i,j,n);
        end
    end
    eps(k)=max(max(abs(u-uold)));
end
[x,y]=meshgrid(0:h:1-h,0:h:1-h);

function f=fcn(x,y)
if (x>=.2) && (x<=.4) && (y>=.1) && (y<=.5)
    f=1;
elseif (x>=.1) && (x<=.9) && (y>=.8) && (y<=.9)
    f=1;
else
    f=0;
end
end

function m=inGeom(x,y,n)
    x=x/n;
    y=y/n;
if (x>=.4) && (x<=.8) && (y>=.3) && (y<=.6)
    m=0;
else
    m=1;
end
end

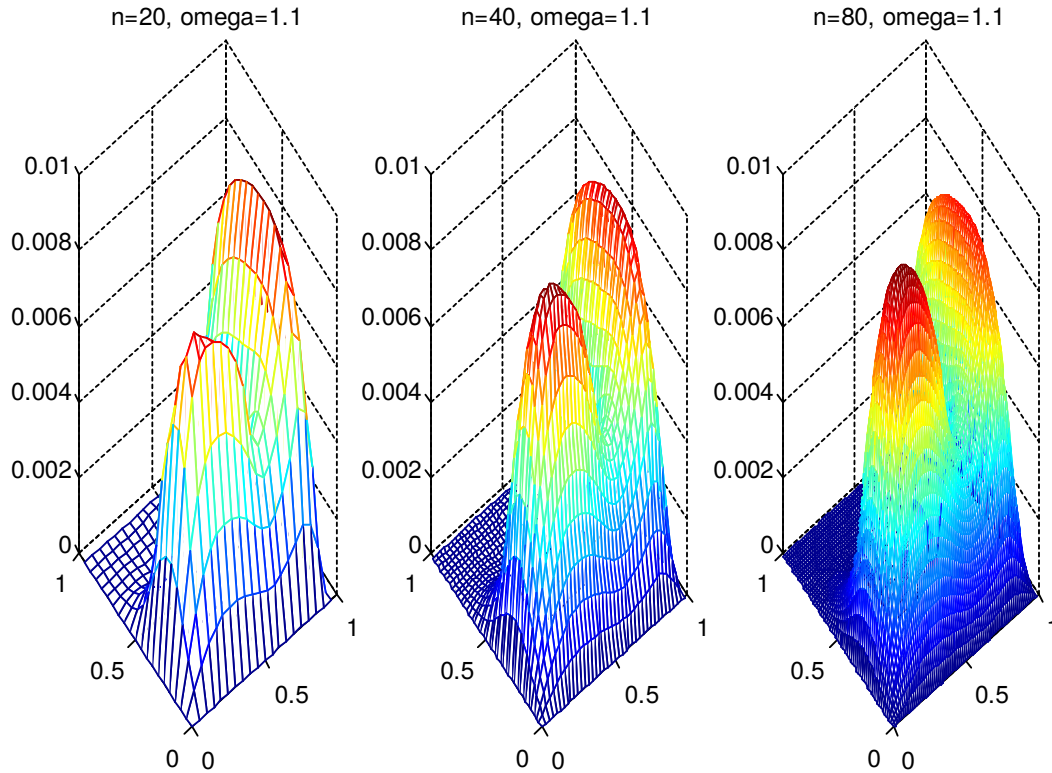
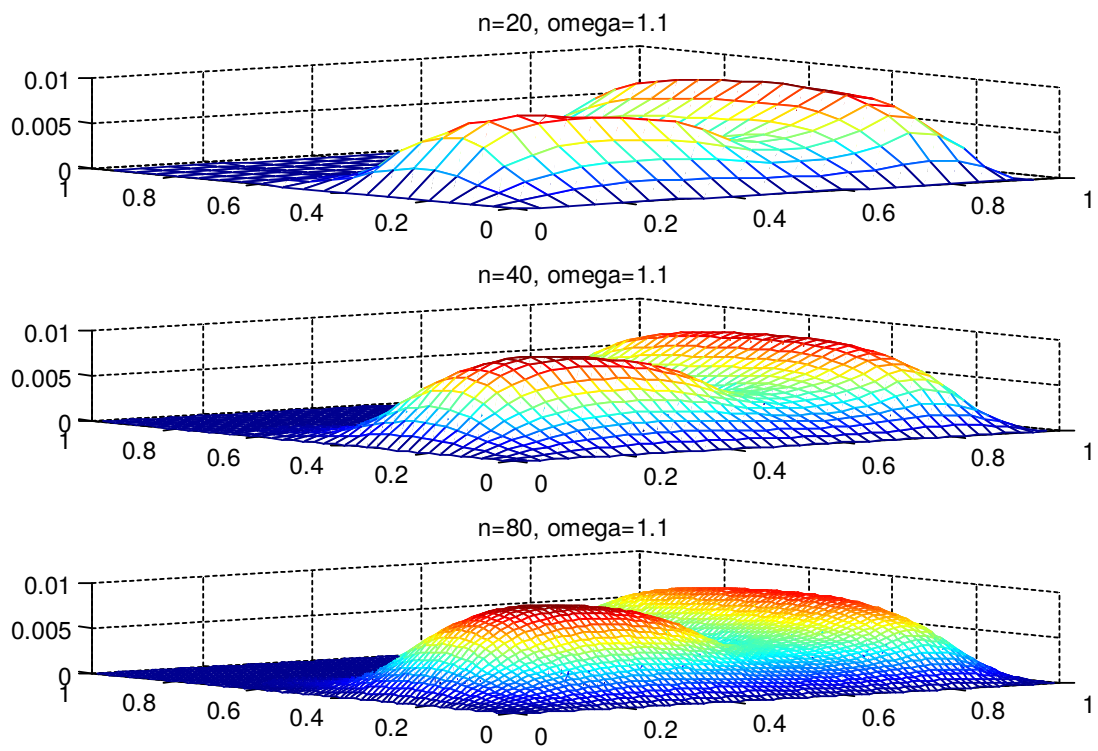
end
```

```

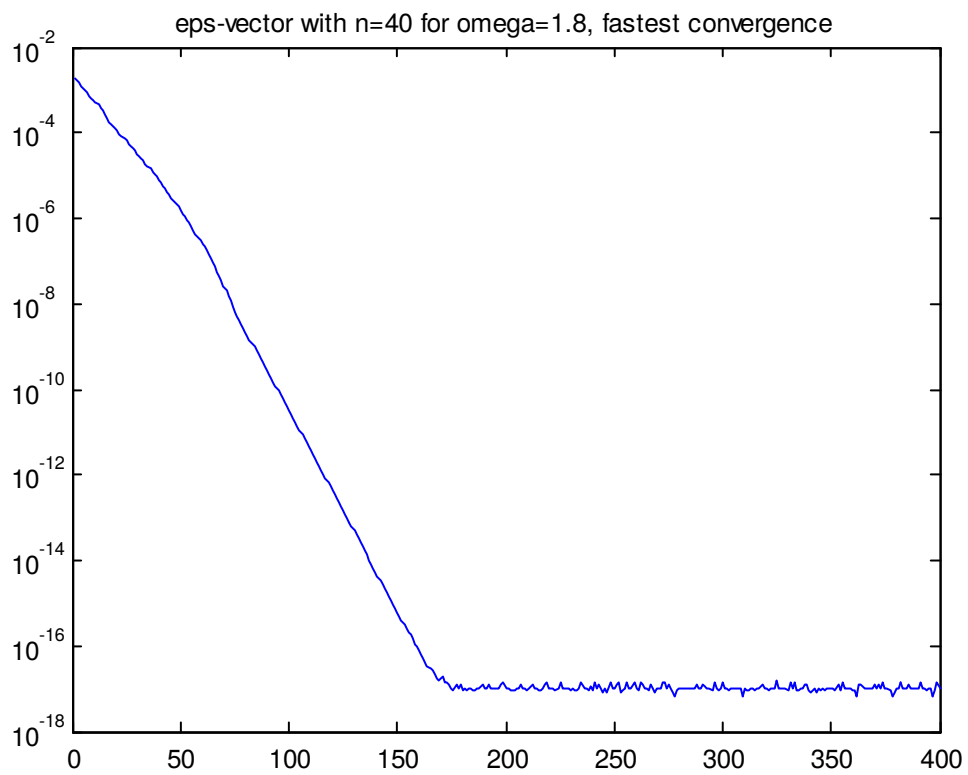
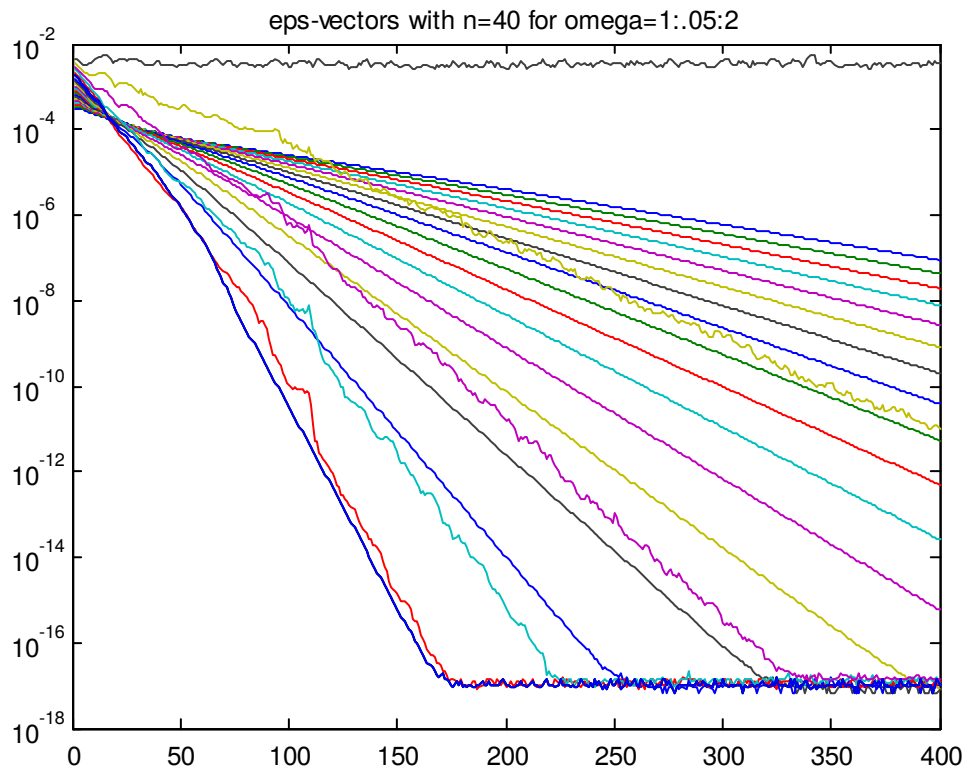
% Part 1.
% [u,x,y,eps]=poisson(20,200,1.1);
% subplot(3,1,1); mesh(x,y,u)
% hold on
% [u,x,y,eps]=poisson(40,400,1.1);
% subplot(3,1,2); mesh(x,y,u)
% hold on
% [u,x,y,eps]=poisson(80,800,1.1);
% subplot(3,1,3); mesh(x,y,u)
% axis([0 1 0 1 0 .01])
% Part 2.
% counter=0;
% eVec=[];
% for omega=1:.05:2
% counter=counter+1;
% [u,x,y,eps]=poisson(40,400,omega);
% eps=eps';
% eVec(:,counter)=eps;
% end
% semilogy(eVec)

```

Part I.



Part II.



Part I.

The solution mesh converged independently of n and for $n=20, 40$ and 80 the mesh graphs appear to be very similar and have the same max peaks, etc. But the graphs got much smoother as n increased due to having more nodes/data to graph/depict. For greater iterations and matrix size (when $n_{\text{iter}}=800$ for $n=80$) the function took longer to compute u, x and y , which was due to the number of flops (and thereby computation time) being dependent on the size of the matrix.

Part II.

Convergence, or the minimization of eps , the difference between the k th and $k-1$ th step of u , was fastest with an ω value of 1.8 . ω values greater than 1 represent over-relaxation, the goal of which is to make the convergence happen as fast as possible, and ω equal to 1 is the same as doing Gauss Seidel instead of SOR. By speeding up convergence we will achieve a minimal difference between the k th and $k-1$ th step of u in the fewest iterations possible, since once $u(k)$ and $u(k-1)$ are very close, our final answer cannot get much better. SOR converged with increasing speed from ω equal to 1 to ω equal to 1.8 , when convergence began to decrease in speed. If we wanted to really increase the speed of convergence, maybe we could use a variable ω method which alters ω depending on the difference between $u(k)$ and $u(k-1)$. We could also save computational time by exiting the SOR program once eps is below a certain error tolerance.