

# Assignment 5 : Laplace Equation

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EE15B025

## Introduction

The assignment is based on currents in a resistor. The currents depend on the shape of the resistor and we want to see if  $R = \rho \frac{L}{A}$  works or not.

A voltage  $V_{AB} = 1V$  is applied across the terminals of a resistor.



As a result, current flows. The current at each point can be described by a “current density”  $\vec{j}$ . This current density is related to the local Electric Field by the conductivity:

$$\vec{j} = \sigma \vec{E}$$

Now the Electric field is the gradient of the potential,

$$\vec{E} = -\nabla\phi$$

and continuity of charge yields

$$\nabla \cdot \vec{j} = -\frac{\partial \rho}{\partial t}$$

Combining these equations we obtain

$$\nabla \cdot (-\sigma \nabla \phi) = -\frac{\partial \rho}{\partial t}$$

Assuming that our resistor contains a material of constant conductivity, the equation becomes

$$\nabla^2 \phi = \frac{1}{\sigma} \frac{\partial \rho}{\partial t}$$

For DC currents, the right side is zero, and we obtain,

$$\nabla^2 \phi = 0$$

## Code

```
from pylab import *
import numpy as np
import mpl_toolkits.mplot3d.axes3d as p3
import sys

def update(phi):
    #Updating the potential
    phi[1:-1,1:-1]=0.25*(phi[:-2,1:-1]+\
        phi[1:-1,0:-2]+phi[2:,1:-1]+phi[1:-1,2:])
    return phi

def assert_boundaries(phi,Nbegin,Nend): #boundary conditions
    phi[1:-1,0]=phi[1:-1,1]
    phi[1:-1,-1]=phi[1:-1,-2]
    phi[0,1:Nbegin],phi[0,Nend+1:-1]=phi[1,1:Nbegin],phi[1,Nend+1:-1]
    phi[-1,1:Nbegin],phi[-1,Nend+1:-1]=phi[-2,1:Nbegin],phi[-2,Nend+1:-1]
    return phi

def fun(Nx=25,Ny=25,Nbegin=8,Nend=17,Niter=1500) :
    error=zeros(Niter)

    phi=zeros((Nx,Ny)) #create phi matrix
    phi[0,Nbegin:Nend+1]=1 #top potential 1
    phi[-1,Nbegin:Nend+1]=0 #lower potential 0

    for k in range(Niter):
        oldphi=phi.copy()
        phi=update(phi)
        phi=assert_boundaries(phi,Nbegin,Nend)
        error[k]=(abs(phi-oldphi)).max()

    logy=log(error)
    x=np.ones((len(error),2))
    x[:,1]=range(Niter)

    ans1=lstsq(x,logy)[0] #from index 0
    ans2=lstsq(x[500:,:],logy[500:])[0] #from index 500
```

```

A1,B1=exp(ans1[0]),ans1[1]
A2,B2=exp(ans2[0]),ans2[1]

q1 = B1*np.arange(Niter)
q2=B2*np.arange(Niter)

print "A and B values for error : "
print "Case 1 : {} and {}".format(A1,B1)
print "Case 2 : {} and {}".format(A2,B2)

title('Evaluation of Error with iteration')
xlabel('Iteration')
ylabel('log(Error)')
semilogy(range(Niter)[::50],error[::50], 'ro', label='error')
semilogy(range(Niter)[::50],A1*np.exp(q1)[::50], 'r', label='fit1')
semilogy(range(Niter)[500::50],A2*np.exp(q2)[500::50], 'g', label='fit2')
legend()
show()

fig1=figure(4)
ax=p3.Axes3D(fig1)
x=arange(1,Nx+1)
y=arange(1,Ny+1)
X,Y=meshgrid(x,y)
title('The 3D Surface plot of the potential')
surf=ax.plot_surface(Y,X,phi,rstride=1,cstride=1,cmap=cm.jet,linewidth=1)
show()

title('Contour plot of potential')
contour(Y,X,phi)
show()

Jx=np.zeros(phi.shape) #current densities
Jy=Jx.copy()

Jx[1:-1,1:-1]=(phi[1:-1,-2]-phi[1:-1,2])/2
Jy[1:-1,1:-1]=(phi[-2,1:-1]-phi[2:,1:-1])/2

title('Vector plot of the current flow')
quiver(y,x,Jy.transpose(),Jx.transpose())
show()
print "Electrode between indices : {} and {}".format(Nbegin,Nend)
print "The Iavg value : {}".format((sum(Jy[:,1]) + sum(Jy[:,-2]))/2)
#Iavg
print "The Idiff value : {}".format(abs(sum(Jy[:,1]) - sum(Jy[:,-2])))

```

```

        print "Resistance : {}".format(1/((sum(Jy[:,1]) + sum(Jy[:,-2]))/2))
    return phi,Nx,Ny

#Order : Nx,Ny,Nbegin,Nend,Niter
fun()
fun(int(sys.argv[1]),int(sys.argv[2]),int(sys.argv[3]),\
    int(sys.argv[4]),int(sys.argv[5])) #for fist case with small electrode
phi,Nx,Ny=fun(int(sys.argv[1]),int(sys.argv[2]),0,\
    int(sys.argv[2])-1,int(sys.argv[5])) #long electrode

array=np.zeros(phi.shape)
arr=np.linspace(Ny,Ny,Ny)-np.arange(0,Ny)
array=arr
epsilon2=sum((phi-array/Ny)**2)/(Nx*Ny)
print 'epsilon2 = {}'.format(epsilon2)

```

## Output

### For default arguments :

```

A and B values for error :
Case 1 : 0.00235717964439 and -0.00276031665063
Case 2 : 0.00141179338348 and -0.00226201175839
Electrode between indices : 8 and 17
The Iavg value : 0.432965780098
The Idiff value : 0.0540234032603
Resistance : 2.30965135345

```

### For specified arguments :

```

python Assg5.py 30 30 10 19 2000
A and B values for error :
Case 1 : 0.00149901445721 and -0.00180659748438
Case 2 : 0.000839141749588 and -0.00138543227524
Electrode between indices : 10 and 19
The Iavg value : 0.388145592997
The Idiff value : 1.66533453694e-16
Resistance : 2.57635283781

```

### Full length electrode :

```

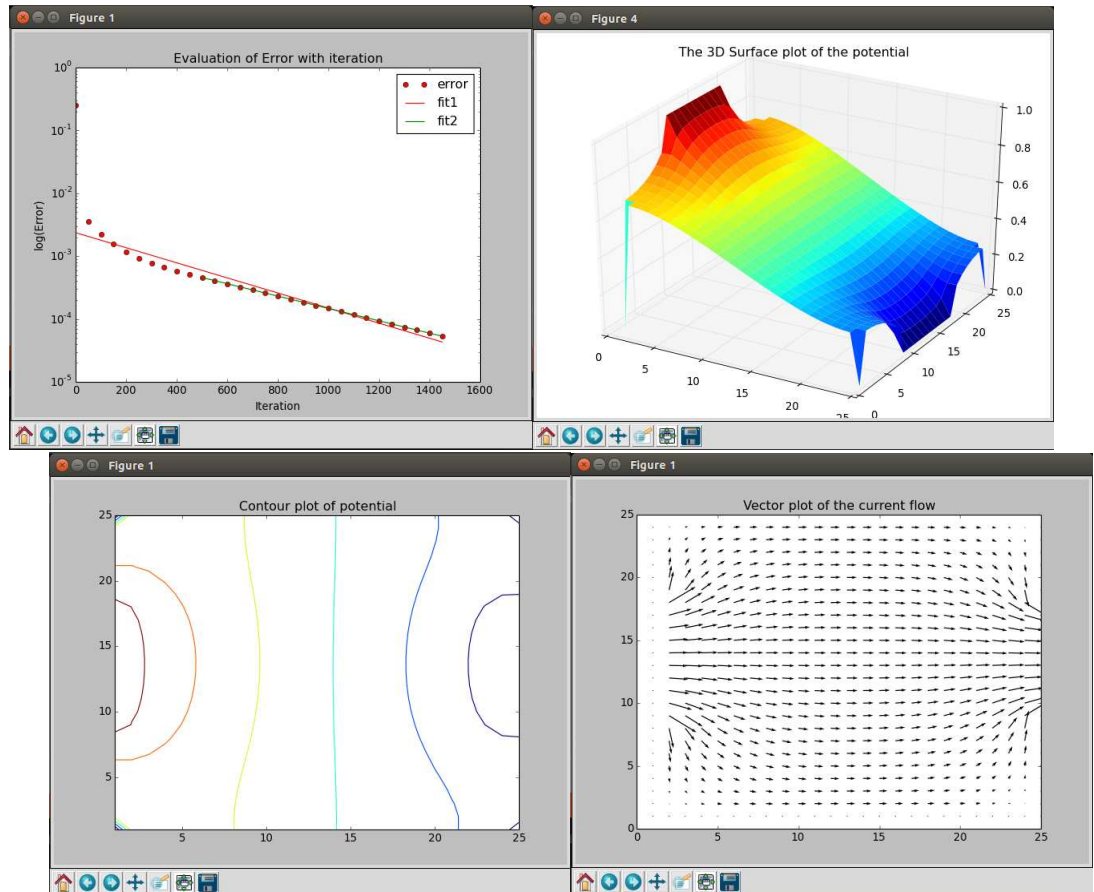
A and B values for error :
Case 1 : 0.00261456004012 and -0.00318576206597
Case 2 : 0.00186331495463 and -0.00293599566766
Electrode between indices : 0 and 29
The Iavg value : 0.965517241375

```

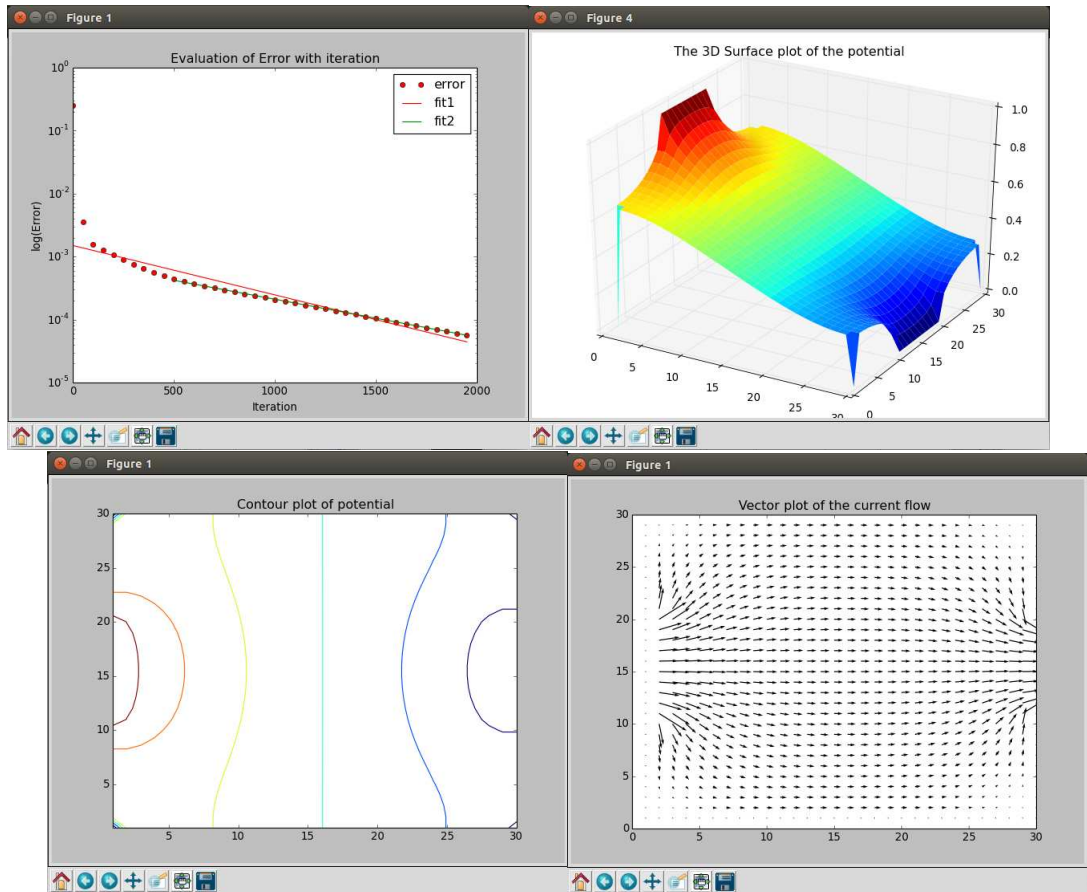
The Idiff value : 0.0  
Resistance : 1.03571428572  
epsilon2 = 0.172637302543

## Graphs :

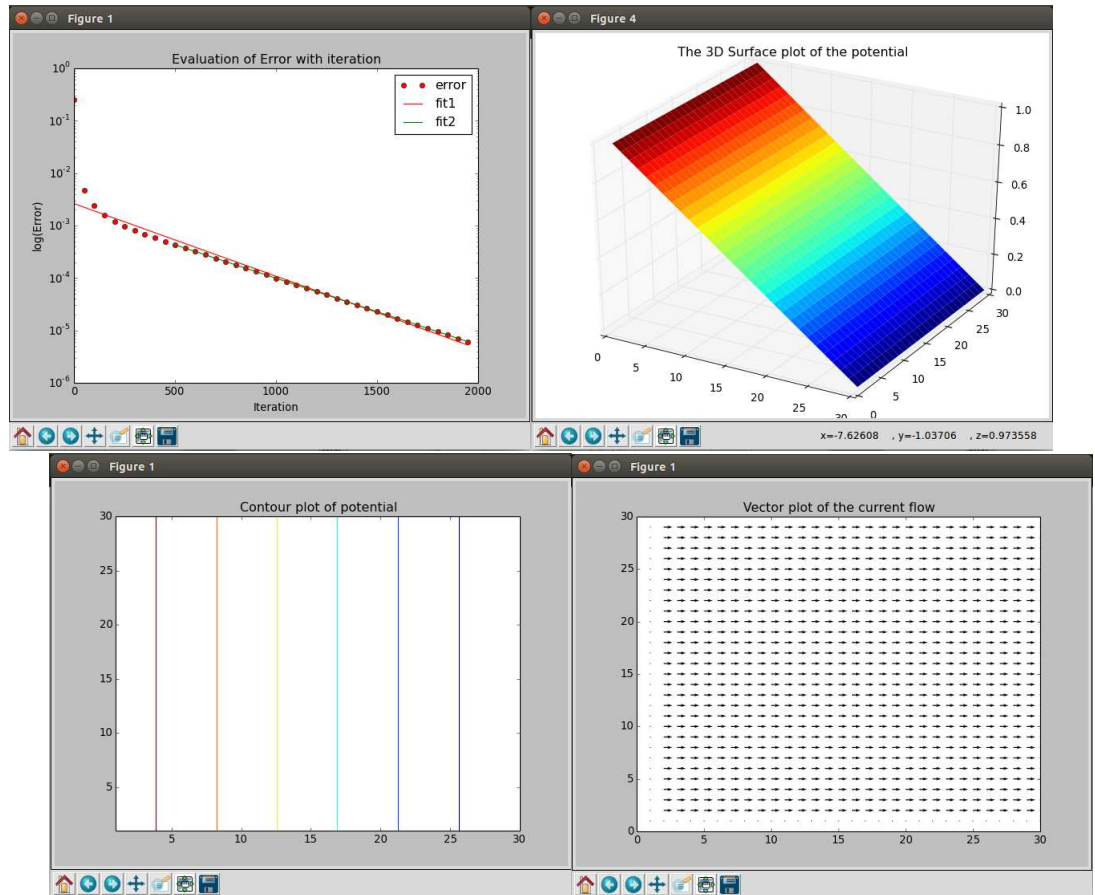
For default arguments :



For specified arguments :



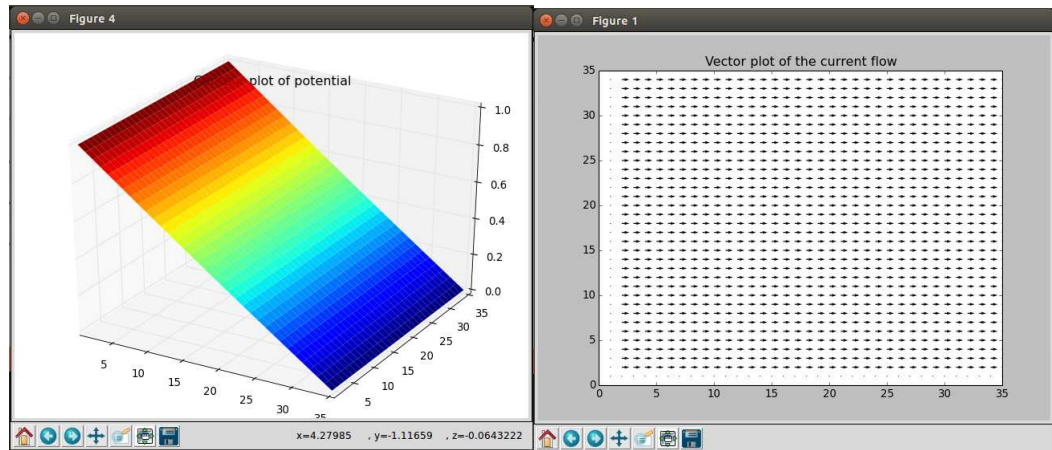
Full length electrode :



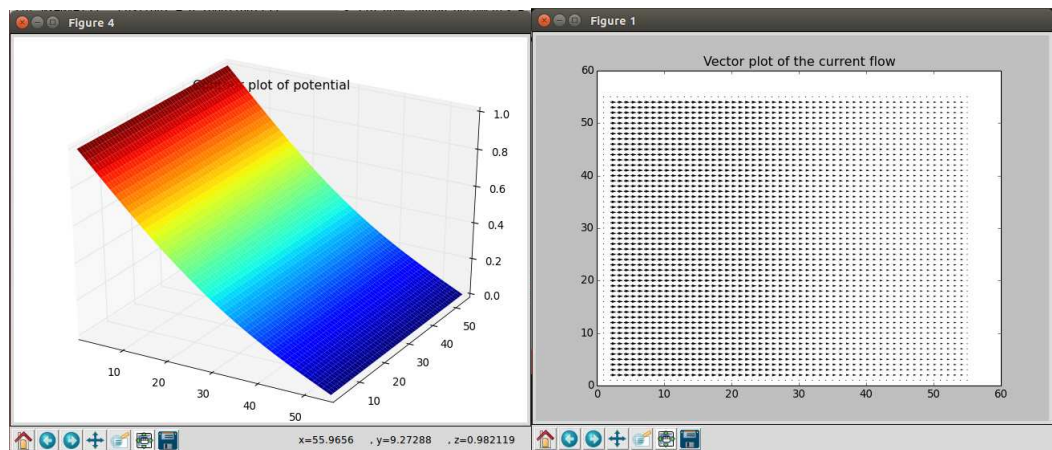
## Inference

The variation of epsilon with increase in  $N_x$  and  $N_y$  can be seen below :

For  $N_x=N_y=35$  ,  $\epsilon_2 = 0.171900070878$

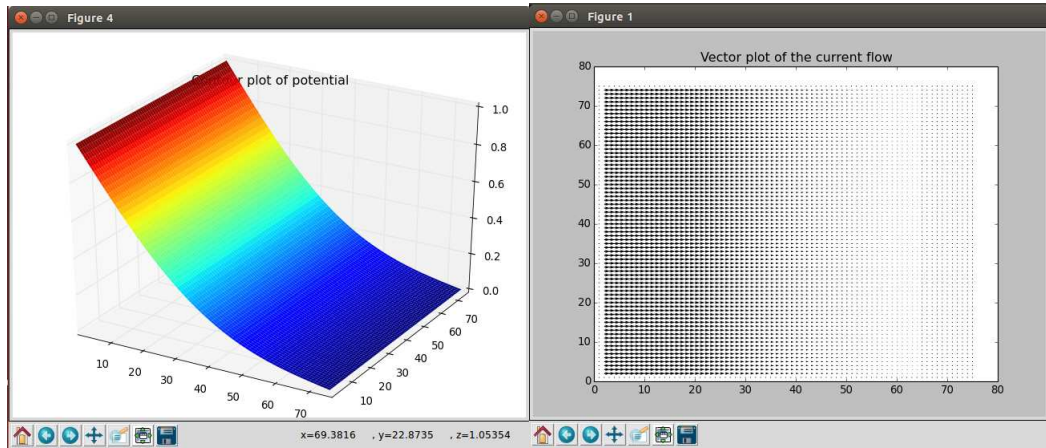


For  $N_x=N_y=55$  ,  $\epsilon_2 = 0.177761140725$

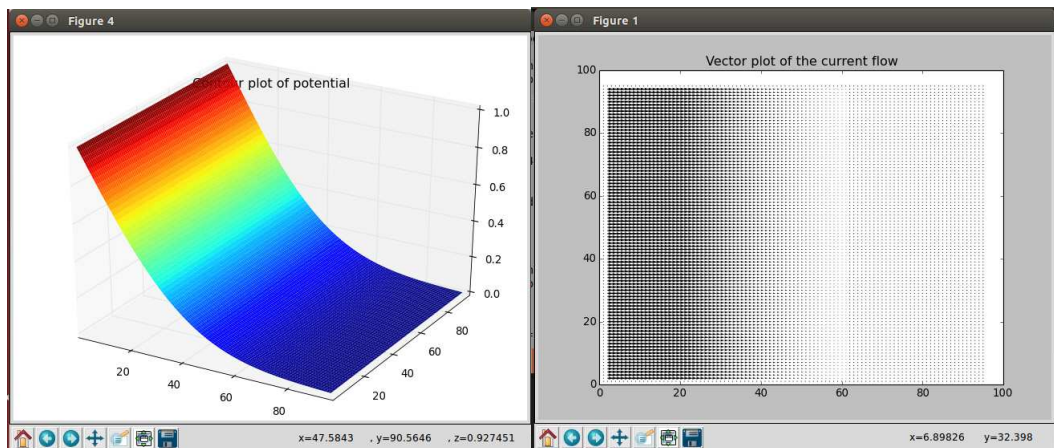




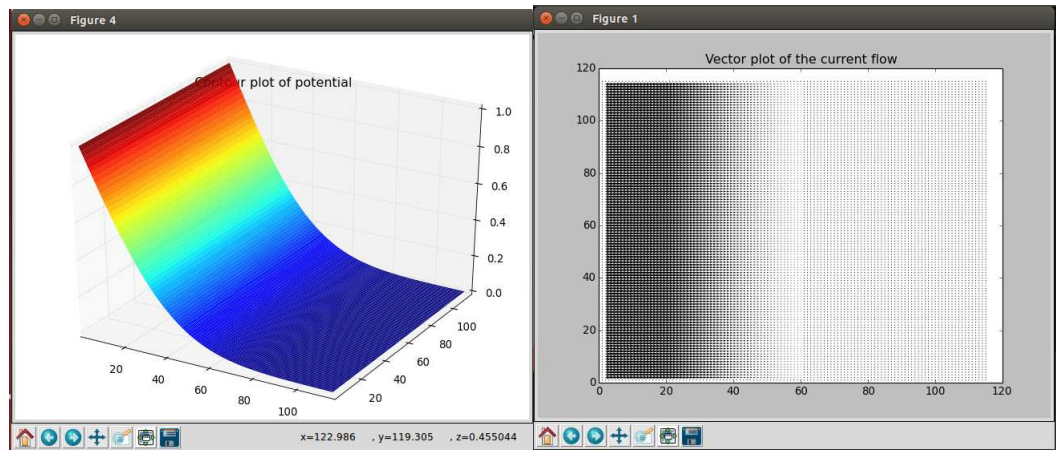
For  $N_x=N_y=75$  ,  $\epsilon_2 = 0.201395730006$



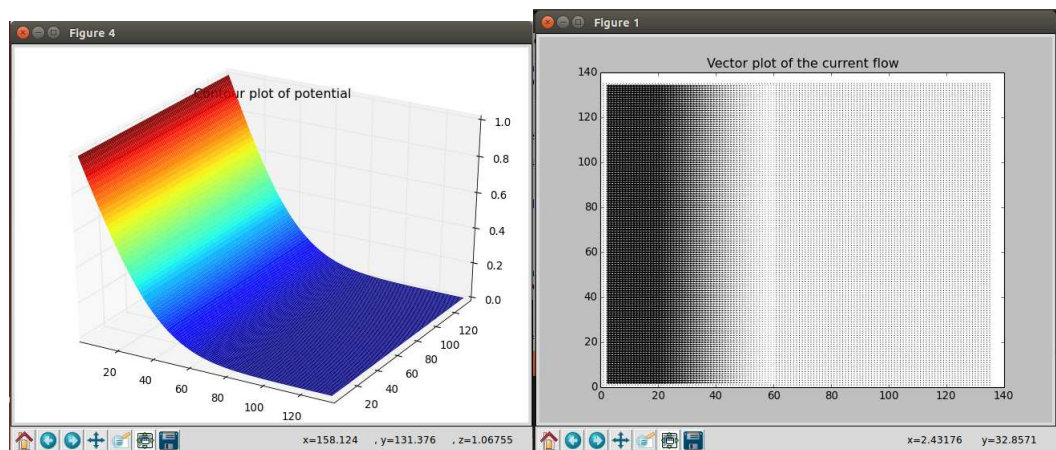
For  $N_x=N_y=95$  ,  $\epsilon_2 = 0.22632677789$



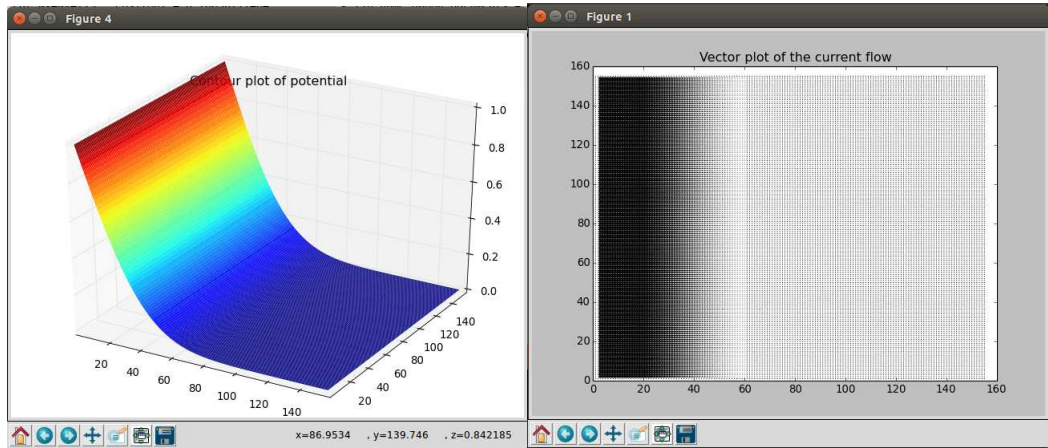
For  $N_x=N_y=115$  ,  $\epsilon_2 = 0.244922581726$



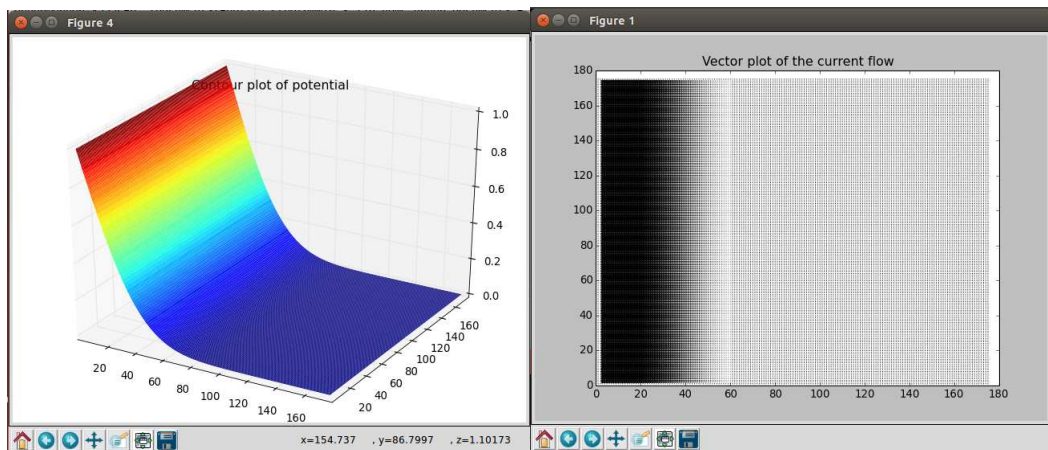
For  $N_x=N_y=135$  ,  $\epsilon_2 = 0.258230621492$



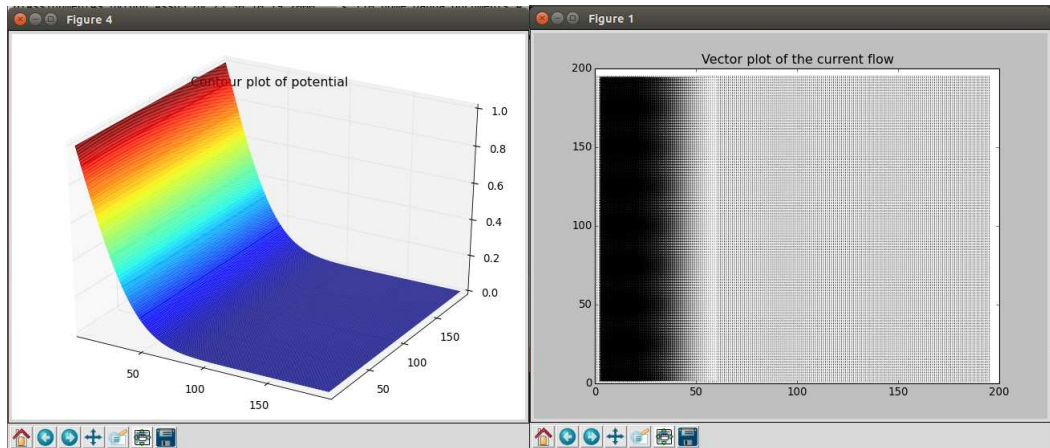
For  $N_x=N_y=155$  ,  $\epsilon_2 = 0.268077062159$



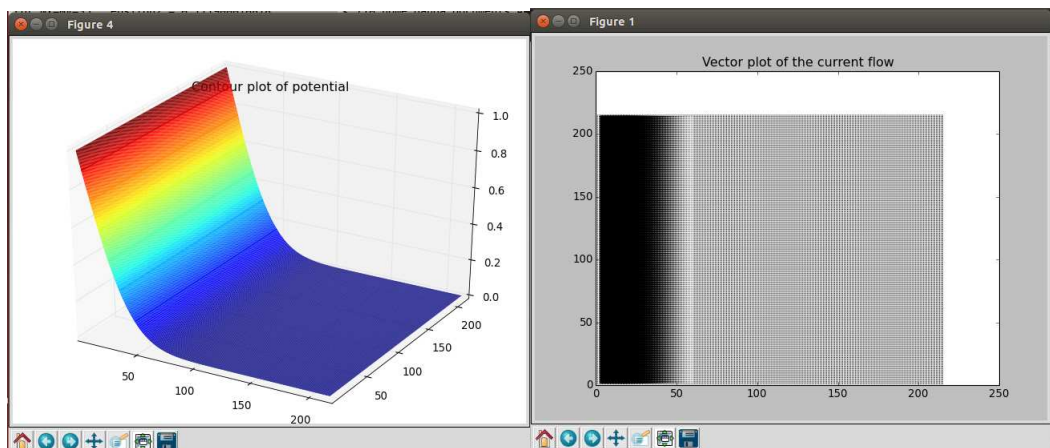
For  $N_x=N_y=175$  ,  $\epsilon_2 = 0.275642578605$



For  $N_x=N_y=195$  ,  $\epsilon_2 = 0.281636426928$



For  $N_x=N_y=215$  ,  $\epsilon_2 = 0.28650217484$



As we can see from the above figures, the epsilon(error) increases with  $N_x$  and  $N_y$ . This explains the change in behaviour of the figures as  $N_x$  and  $N_y$  increases. ie, the figures become more and more erroneous. This shows the ineffectiveness in using the above method for practical applications.

The code used for computing the above errors is as follows:

```
Nx=35
Ny=35
for i in range(10):
    phi ,Nx,Ny=fun(Nx,Ny,0 ,Ny-1,2000)
    array=np.zeros(phi.shape)
    arr=np.linspace(Ny,Ny,Ny)-np.arange(0,Ny)
    array=arr
    epsilon2=sum((phi-array/Ny)**2)/(Nx*Ny)
```

```
print 'For Nx=Ny={} , epsilon2 = {}'.format(Nx,epsilon2)
Nx+=20
Ny=Nx
```

Only requires figures were displayed and rest of the code commented during the process.