

2024 Physics 2 Homework 1

Firstly, we open the “aPart.py” to start the explanation.

```
import numpy as np
import matplotlib.pyplot as plt

# Create a 10x10 matrix
matrix = np.zeros((10, 10))

# Constants
k = 8.99e9 # Coulomb's constant in N m^2/C^2
q = 5e-9 # Charge in C

# Calculate potential for each cell
for i in range(10):
    for j in range(10):
        r = np.sqrt(i**2 + j**2) * 0.1 # Distance in meters
        matrix[i, j] = k * q / r

plt.imshow(matrix, cmap='viridis', origin='lower', extent=[0, 1, 0, 1])
plt.colorbar(label='Electric Potential (V)')
plt.xlabel('X')
plt.ylabel('Y')
plt.title('Electric Potential Distribution')
plt.show()

x_values = np.arange(0, 1, 0.1)
v_values = matrix[:, 0] # Potential values for j=0

plt.plot(x_values, v_values)
plt.xlabel('X')
plt.ylabel('Electric Potential (V)')
plt.title('Electric Potential vs. X')
plt.grid()
plt.show()

r_values = np.sqrt(np.arange(0, 10)**2 + np.arange(0, 10)**2) * 0.1

plt.plot(r_values, np.diag(matrix))
plt.xlabel('Diagonal Distance (m)')
plt.ylabel('Electric Potential (V)')
plt.title('Electric Potential vs. Diagonal Distance')
plt.grid()
plt.show()

# In this electric potential distribution, points that have the same electric potential
# value are those located on the same equipotential surface.
# An equipotential surface is a surface where the electric potential is constant.
# Electric Potential Distribution Plot:
# The first plot shows the electric potential distribution across the 10x10 matrix.
# Darker regions represent higher potential values.
# The potential decreases as you move away from the central charge (located at the
# origin).
# The circular contours indicate equipotential surfaces.

# The f part is here:
```

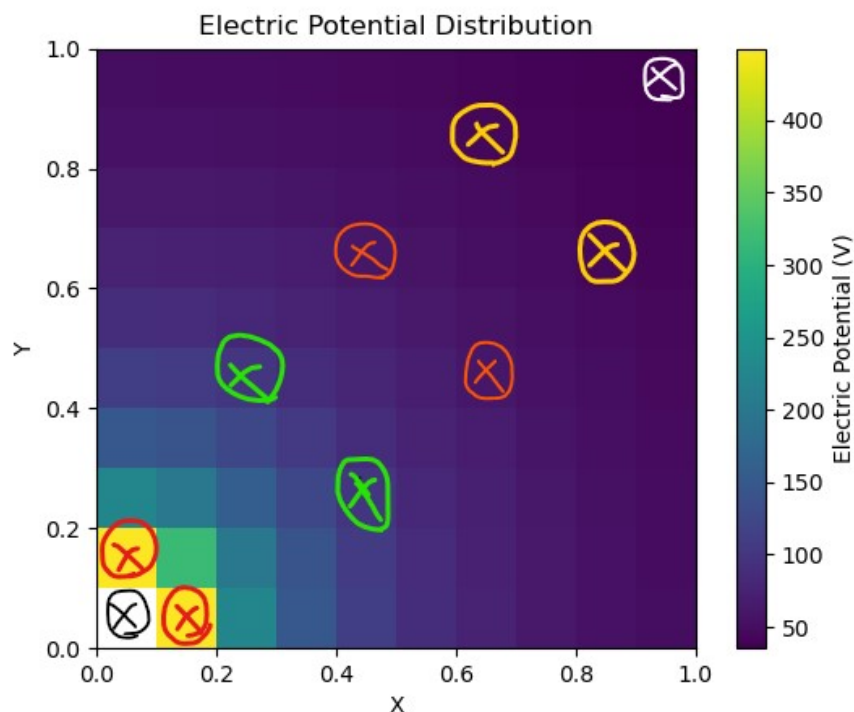
```
# By using a finer grid (100x100), the equipotential lines become smoother and more
detailed, providing a clearer representation of the potential distribution.

# Create a 100x100 matrix for a finer grid
matrix_fine = np.zeros((100, 100))

# Calculate potential for each cell in the finer grid
for i in range(100):
    for j in range(100):
        r = np.sqrt(i**2 + j**2) * 0.01 # Adjusted distance for finer grid
        if r != 0:
            matrix_fine[i, j] = k * q / r
        else:
            matrix_fine[i, j] = np.inf

# Plotting with finer grid
plt.imshow(matrix_fine, cmap='viridis', origin='lower', extent=[0, 1, 0, 1])
plt.colorbar(label='Electric Potential (V)')
plt.xlabel('X')
plt.ylabel('Y')
plt.title('Electric Potential Distribution with Finer Grid')
plt.contour(np.linspace(0, 1, 100), np.linspace(0, 1, 100), matrix_fine,
            colors='white', levels=np.linspace(0, np.max(matrix_fine[matrix_fine != np.inf]), 20))
plt.show()

# Below is the code that has been modified and edited so that we can draw equipotential
lines:
# aPartF.py
with open('aPartF.py') as file:
    exec(file.read())
```



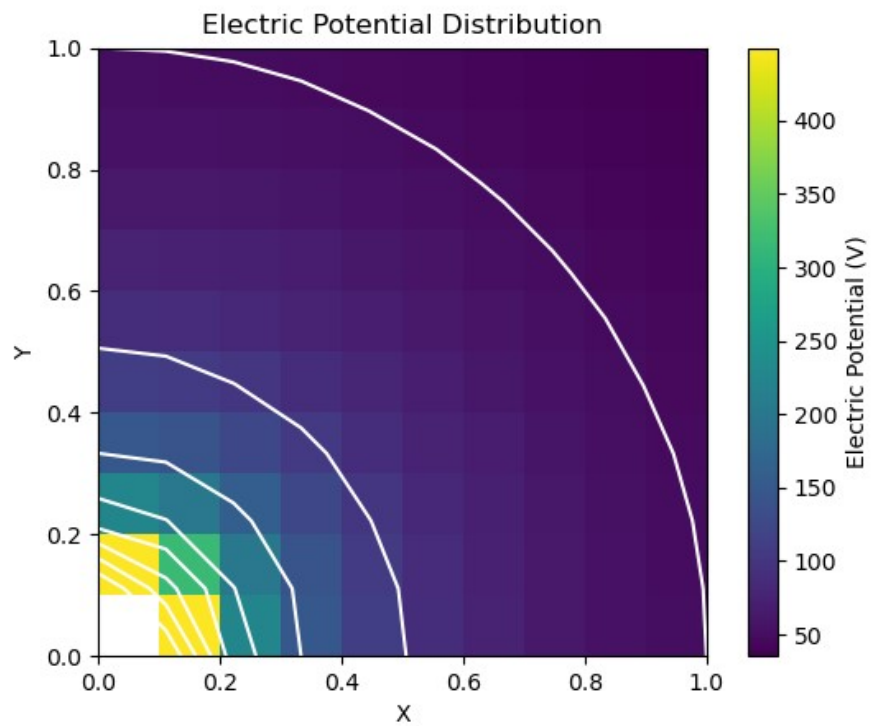
Points with equal electric potential are shown as the same colored markings on the matrix. For example, four identical points are marked: red, green, orange and yellow. Black and white markings symbolize maximum and minimum values.

```
C:\ProgramData\anaconda3\python.exe "C:\Users\enesi\OneDrive - Akdeniz Üniversitesi\Masaüstü\PhysicsHM\firstPart\savingMatrixData.py"
[[ 0. 449.375 224.6875 149.79166667 112.34375
 89.875 74.89583333 64.19642857 56.171875 49.93055556]
[449.375 317.7561098 200.96660948 142.10485235 108.9894465
 88.12968823 73.87679492 63.55122196 55.73810886 49.62516703]
[224.6875 200.96660948 158.8780549 124.63420034 100.48330474
 83.44684259 71.05242618 61.72640342 54.49472325 48.74156224]
[149.79166667 142.10485235 124.63420034 105.91870327 89.875
 77.0671767 66.98886983 59.00582827 52.59536552 47.36828412]
[112.34375 108.9894465 100.48330474 89.875 79.43902745
 70.18058425 62.31710017 55.73810886 50.24165237 45.6271183 ]
[ 89.875 88.12968823 83.44684259 77.0671767 70.18058425
 63.55122196 57.53657292 52.23878266 47.63365473 43.64715844]
[ 74.89583333 73.87679492 71.05242618 66.98886983 62.31710017
 57.53657292 52.95935163 48.74156224 44.9375 41.54473345]
[ 64.19642857 63.55122196 61.72640342 59.00582827 55.73810886
 52.23878266 48.74156224 45.39372997 42.27364402 39.41279474]
[ 56.171875 55.73810886 54.49472325 52.59536552 50.24165237
 47.63365473 44.9375 42.27364402 39.71951372 37.31856251]
[ 49.93055556 49.62516703 48.74156224 47.36828412 45.6271183
 43.64715844 41.54473345 39.41279474 37.31856251 35.30623442]]

Process finished with exit code 0
|
```

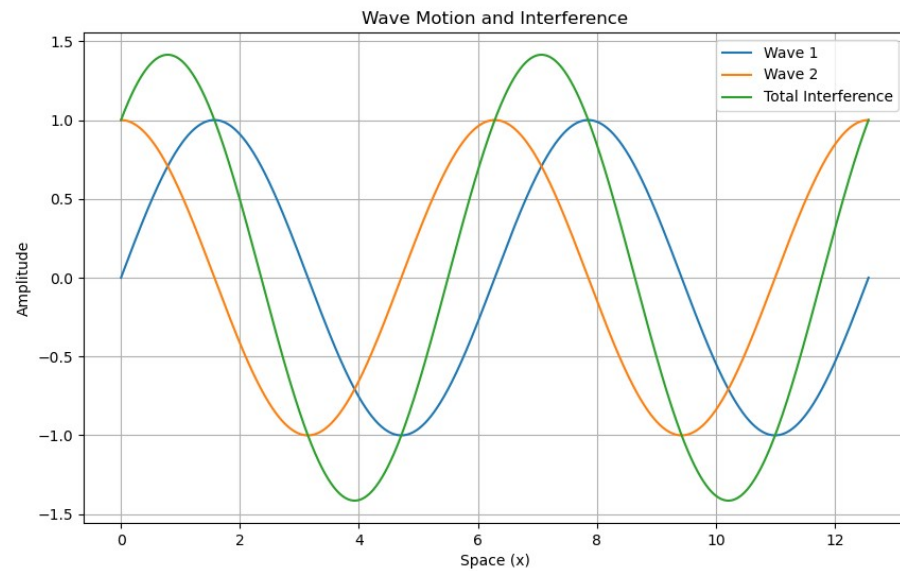
[illegible]

In the f part of the homework it runs “aPartF.py” as a compiler.

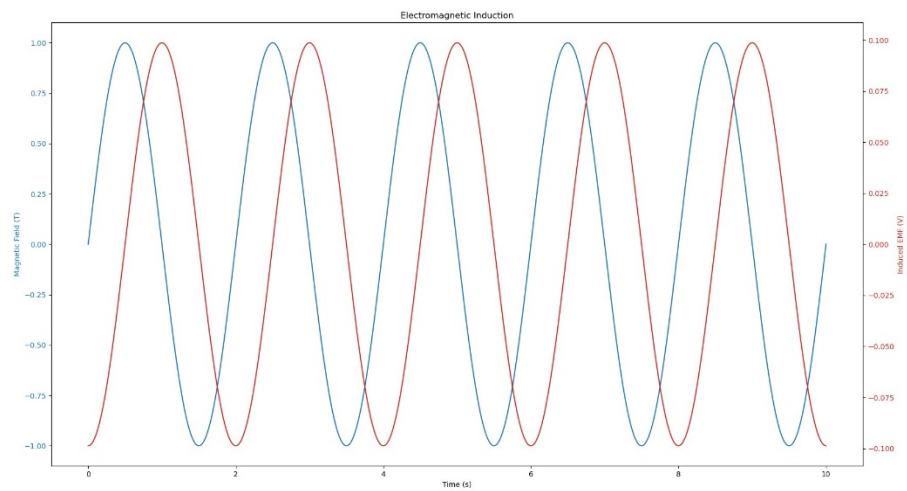


Then, when we begin the question 2. I want to do multiple simulation.

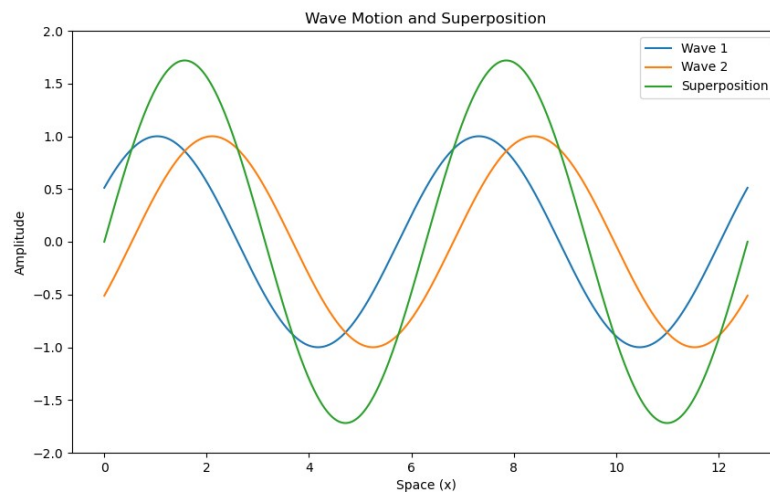
1. Wave Motion and Interference



2. Electromagnetic Induction

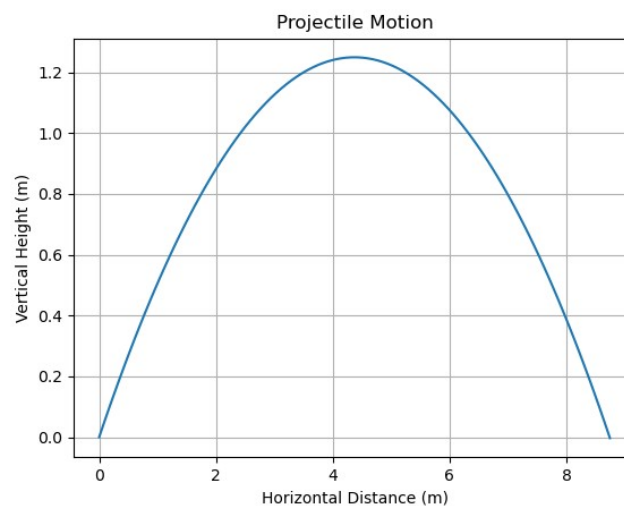


3. Wave Motion and Superposition



4. Shooting Movements

```
C:\ProgramData\anaconda3\python.exe "C:\Users\enesi\OneDrive - Akdeniz Üniversitesi\Masaüstü\PhysicsHM\bonusPart\1\bonusPart.py"  
Enter the force at launch (Newtons): 10  
Enter the launch angle (0 to 90 degrees): 30  
  
Process finished with exit code 0
```



You can reach the codes of this homework by the following link:

<https://github.com/dahaii/Physics2HM>