Chapter 31-5: Elliptic PDEs Using the Monte Carlo Method.

Monte Carlo Methods are a class of computational solutions that can be applied to various problems which rely on repeated random sampling to provide generally approximate solutions. Monte Carlo is a stochastic approach, in which a series of simulations (trials), representing the analyzed problem, with randomly selected input values, are performed. Among these trials, a specified number of properly defined successes is achieved. The ratio between the number of success trials to the number of all trials, scaled by dimensional quantity (e.g., area or function value) allows for the estimation of the unknown solution, providing the number of trials is large enough.

A capacitor is constructed out of two infinite metal plates spaced 10 cm apart from one another. One plate is placed at 5V potential with respect to the other and the space between the plates is free of charges. The number of random walks is taken to be 400 and the number of lattice points is taken to be 30.

(The description of the problem and the code solution was taken from the repository of s-ankur.)

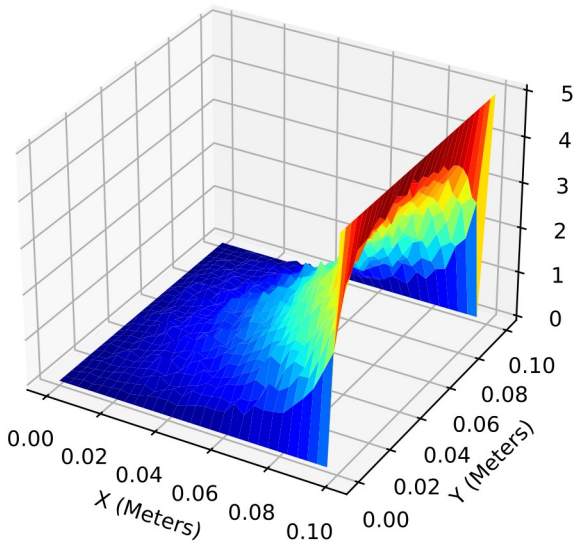
In [1]:

```
"""
All calculations in SI units
"""
import random
from mpl_toolkits.mplot3d import axes3d
import matplotlib.pyplot as plt
from matplotlib import cm
import numpy as np
%config InlineBackend.figure_formats = ['svg']

h = 10e-2  # Distance between plates = 10cm
lattice_points = 30  # Number of Points in lattice
d = h / lattice_points  # Lattice size
boundary_voltage_high = 5.0  # 5 Volts at Positive Plate
boundary_voltage_low = 0.0  # 0 Volts at Negative Plate
epsilon_naught = 8.854e-12  # Permittivity of Vaccum
charge_density = 6e-16  # Coulomb per meter cube
N = 400  # Number of Random Walks


def f(x):
    # The Function \nabla^2(phi)  = f
    # For Laplace f = 0
    return 0


def g(x):
    # Two Dimensional Boundary Conditions: two parallel metal plates at x=0,x=h
    # the plate at x=h is at high potential and x=0 is low potential
    # Assume that there are metal plates along y=0 and y=h (uncharged)
    # this is because I dont know how to simulate open boundry conditions
    if x[0] <= 0:
        return boundary_voltage_low
    if x[0] >= h:
        return boundary_voltage_high
    if x[1] <= 0 or x[1] >= h:
        return boundary_voltage_low

"""
def f_2(x):
    # Alternative charge distribution: A charged Sphere in the centre of metal box
```

```
    if (h / 2 - x[0]) ** 2 + (h / 2 - x[1]) ** 2 <= (h / 5) ** 2:
        return -charge_density * 5 / epsilon_naught
    else:
        return 0


def g_2(x):
    # Two Dimentional Alternative Boundary Conditions: uncharged metal box
    return 0
"""


def f_3(x):
    # Alternative charge distribution: TWO charged Sphere in the centre of metal box
    if (h / 3 - x[0]) ** 2 + (h / 2 - x[1]) ** 2 <= (h / 5) ** 2:
        return -charge_density * 5 / epsilon_naught
    if (2 * h / 3 - x[0]) ** 2 + (h / 2 - x[1]) ** 2 <= (h / 5) ** 2:
        return charge_density * 5 / epsilon_naught
    else:
        return 0


@np.vectorize
def poisson_approximation_fixed_step(*A):
    # Returns the Value of Potential Feild at a given point A with N random walks
    result = 0
    F = 0
    for i in range(N):
        x = list(A)
        while True:
            if x[0] <= 0 or x[0] >= h or x[1] <= 0 or x[1] >= h:
                break
            random_number = random.randint(0, 3)
            if random_number == 0:
                x[0] += d
            elif random_number == 1:
                x[0] -= d
            elif random_number == 2:
                x[1] += d
            elif random_number == 3:
                x[1] -= d
            F += f(x) * h ** 2
        result += g(x) / N
    result = result - F
    return result


def plot(x, y, z):
    # Function for plotting the potential
    fig = plt.figure()
    ax = fig.add_subplot(111, projection="3d")

    ax.plot_surface(x, y, np.array(z), cmap=cm.jet, linewidth=0.1)
    plt.xlabel("X (Meters)")
    plt.ylabel("Y (Meters)")
    ax.set_zlabel("Potential (Volts)")
    plt.show()


if __name__ == "__main__":
    # Experiment E: 2D Capacitor
    print(
        f"Calculating Monte Carlo with {lattice_points}x{lattice_points} lattice points an
    )
    lattice_x, lattice_y = np.mgrid[
        0 : h : lattice_points * 1j, 0 : h : lattice_points * 1j
    ]
    z = poisson_approximation_fixed_step(lattice_x.ravel(), lattice_y.ravel()).reshape(
        lattice_x.shape
    )
```

```
    plot(lattice_x, lattice_y, z)

    # Experiment F: Metal box with positively charged metal ball inside
#   f = f2
#   g = g2
    print(
        f"Calculating Monte Carlo with {lattice_points}x{lattice_points} lattice points an
    )
    lattice_x, lattice_y = np.mgrid[
        0 : h : lattice_points * 1j, 0 : h : lattice_points * 1j
    ]
    z = poisson_approximation_fixed_step(lattice_x.ravel(), lattice_y.ravel()).reshape(
        lattice_x.shape
    )
    plot(lattice_x, lattice_y, z)

    # Experiment G: Metal Box with two spheres (positive and negative)

    f = f_3
    g = g_2
    print(
        f"Calculating Monte Carlo with {lattice_points}x{lattice_points} lattice points an
    )
    lattice_x, lattice_y = np.mgrid[
        0 : h : lattice_points * 1j, 0 : h : lattice_points * 1j
    ]
    z = poisson_approximation_fixed_step(lattice_x.ravel(), lattice_y.ravel()).reshape(
        lattice_x.shape
    )
    plot(lattice_x, lattice_y, z)

```

```
Calculating Monte Carlo with 30x30 lattice points and 400 random walks
```




```
---------------------------------------------------------------------------
NameError                                 Traceback (most recent call last)
~\AppData\Local\Temp/ipykernel_11260/807967070.py in <module>
    116 #   g = g2
    117     print(
--> 118         f"Calculating Monte Carlo with {lattice_points}x{lattice_points} lattice p
oints and {N} random walks for {'Laplace' if laplace else 'Poisson'}"
    119     )
    120     lattice_x, lattice_y = np.mgrid[

NameError: name 'laplace' is not defined
```

In [ ]:

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