

Marine Hydrodynamics

Assignment 1

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Assignment

For specified (see below) two-dimensional geometries, assuming potential theory in unbounded fluid, and use of Green's second identity, calculate the velocity potential along the body and the added mass forces, for a circle, an ellipse, a square and a rectangle, moving laterally, and with rotation. Find also the cross coupling added mass coefficients. For the circle, the reference solution is: $\phi = -a^2/(x^2)$ where a denotes the cylinder radius, $r^2 = x^2 + y^2$.

Theory, numerics and program

In this assignment we use the method of panels. By splitting the geometry into N equal parts, and assuming that $\phi, \frac{\partial \phi}{\partial n}$ is constant every segment.

From Newman chapter 4 we have (79):

$$\int_C \phi \frac{\partial G}{\partial n} - G \frac{\partial \phi}{\partial n} = -\pi \phi(x, y)$$

where $G = \ln(r)$, a source in 2D, and C is the circumference of the geometry

$$-\pi \phi(x_0) + \int_C \phi \frac{\partial}{\partial n} \ln(r) dl = \int_C \ln(r) \frac{\partial \phi}{\partial n} dl$$

where x_0 states a point on our geometry.

We got a trick from the lectures, turning it into:

$$-\pi \phi(x_0) + \sum_{n=1}^N \phi(x_n)(\theta_a - \theta_b) = \int_C \ln(r) \frac{\partial \phi}{\partial n} dl$$

$$\begin{pmatrix} -\pi & (\theta_a - \theta_b)_0 & (\theta_a - \theta_b)_1 & \dots \\ (\theta_a - \theta_b)_N & -\pi & (\theta_a - \theta_b)_0 & \dots \\ (\theta_a - \theta_b)_{N-1} & (\theta_a - \theta_b)_N & -\pi & \dots \\ \dots & \dots & \dots & \dots \end{pmatrix} \begin{pmatrix} \phi_0 \\ \phi_1 \\ \phi_2 \\ \dots \\ \phi_N \end{pmatrix} = \begin{pmatrix} \int \ln(r_0) \frac{\partial \phi}{\partial n} dl \\ \int \ln(r_1) \frac{\partial \phi}{\partial n} dl \\ \int \ln(r_2) \frac{\partial \phi}{\partial n} dl \\ \dots \end{pmatrix}$$

To calculate the ϕ values we created a fictional point between the x_N points, where we stand in this ghostpoint and calculate the $\Delta\theta$ to the x_N values. This is done since we have assumed ϕ to be constant. To evaluate the integrals we used the trapezoidal method. Every line of the matrix on the right hand side is calculated by going around the geometry once.

The normal vector is defined as:

$$\bar{n} = \frac{\frac{x}{a^2} \bar{i} + \frac{y}{b^2} \bar{j}}{(\frac{x}{a^2})^2 + (\frac{y}{b^2})^2}$$

which is used for ellipse and for circle when $a = b$.

For a circle we have an exact solution for $\phi = -\frac{a^2x}{r^2}$, which is $\phi = -a\cos(\theta)$ and an exact solution for added mass $m_{11} = \rho r_a^2 \pi$

For an ellipse we only have the exact solution for added mass $m_{11} = \rho r_a^2 \pi$, $m_{22} = \rho r_b^2 \pi$, $m_{66} = \frac{\pi}{8} \rho (r_a^2 - r_b^2)^2$.

Next page is the program code in Python

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from matplotlib.pyplot import *
import numpy as np
import math as math

def make_points(num_points, r_a, r_b):
    dx = 2*pi/num_points
    x = np.zeros(num_points*3)
    y = np.zeros(num_points*3)
    for i in range(num_points*2+1):
        x[i] = r_a*np.cos(i*dx)
        y[i] = r_b*np.sin(i*dx)
    return x,y

def make_angle(N,r_a,r_b):
    angle = np.zeros(N-1)
    x,y = make_points(N,r_a,r_b)
    matrix_ = np.zeros((N,N))
    for i in range(N):
        x_first = (x[i]+x[i+1])/2.0
        y_first = (y[i]+y[i+1])/2.0
        for k in range(N):
            if i == k:
                matrix_[i][k] = -np.pi
            else:
                xa = x[k] - x_first
                xb = x[k+1] - x_first
                ya = y[k] - y_first
                yb = y[k+1] - y_first
                matrix_[i][k] = -np.arccos((xa*xb + ya*yb)/(np.sq

    return matrix_

def integral_1(N,r_a,r_b,direction1):
    x,y = make_points(N,r_a,r_b)
    traps1 = 0
    traps2 = 0
    Matrix_B = np.zeros(N)
    for i in range(N):
        integral_x = 0
        integral_66 = 0
        x0 = 0.5*(x[i]+x[i+1])
        y0 = 0.5*(y[i]+y[i+1])
        for j in range(N-1):
            rad_1 = np.sqrt((x0-x[j])**2 + (y0-y[j])**2)
            rad_2 = np.sqrt((x0-x[j+1])**2 + (y0-y[j+1])**2)

            if direction1 == 11:
                traps1x = -((x[j]+x[j+1])/(2*r_a**2)) / np.sqrt(
((x[j]+x[j+1])/(2*r_a**2))**2 + ((y[j]+y[j+1])/(2*r_b**2))**2 ) * np.log(rad_1)
                traps2x = -((x[j+1]+x[j+2])/(2*r_a**2)) / np.sqrt(
((x[j+1]+x[j+2])/(2*r_a**2))**2 + ((y[j+1]+y[j+2])/(2*r_b**2))**2 ) * np.log(rad_1)
                ds = np.sqrt((x[j+1]-x[j])**2 + (y[j+1] - y[j])**2)
                integral_x = integral_x + (traps1x + traps2x) * d

    * 0.5

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else :
    r1a = (x[j] + x[j+1])*0.5 ; r2a = 0.5*(y[j] + y[j+1]
    n1a = -((x[j]+x[j+1])/(2*r_a**2)) / np.sqrt(
((x[j]+x[j+1])/(2*r_a**2))**2 + ((y[j]+y[j+1])/(2*r_b**2))**2 )
    n2a = -((y[j]+y[j+1])/(2*r_b**2))/np.sqrt(
((x[j]+x[j+1])/(2*r_a**2))**2 + ((y[j]+y[j+1])/(2*r_b**2))**2 )
    crosses1 = r1a*n2a - r2a*n1a

    r1b = (x[j+1] + x[j+2])*0.5 ; r2b = 0.5*(y[j+1] +
    n1b = -((x[j+1]+x[j+2])/(2*r_a**2))/np.sqrt(
((x[j+1]+x[j+2])/(2*r_a**2))**2 + ((y[j+1]+y[j+2])/(2*r_b**2))**2 )
    n2b = -((y[j+1]+y[j+2])/(2*r_b**2))/np.sqrt(
((x[j+1]+x[j+2])/(2*r_a**2))**2 + ((y[j+1]+y[j+2])/(2*r_b**2))**2 )
    crosses2 = r1b*n2b - r2b*n1b

    traps3 = np.log(rad_1) * crosses1
    traps4 = np.log(rad_2) * crosses2
    ds = np.sqrt((x[j+1]-x[j])**2 + (y[j+1] - y[j])**2)
    integral_66 = integral_66 + (traps3 +traps4)* ds
* 0.5

    if direction1 == 11:
        Matrix_B[i] = integral_x
    else :
        Matrix_B[i] = integral_66
return Matrix_B

def solver1(N,r_a,r_b,direction1):
    return np.linalg.solve(make_angle(N,r_a,r_b),integral_1(N,r_a,r_b,direction1))

def added_mass(N,r_a,r_b):
    phix = solver1(N,r_a,r_b,direction1 =11)
    phi6 = solver1(N,r_a,r_b,direction1 =66)
    x,y = make_points(N,r_a,r_b)
    add_m11 = 0
    add_m22 = 0
    add_m66 = 0
    for i in range(N-2):
        rad_1 = np.sqrt( ((x[i]+x[i+1])/(2*r_a**2))**2 + ((y[i] +y[i+1]
        rad_2 = np.sqrt( ((x[i+1]+x[i+2])/(2*r_a**2))**2 +
((y[i+1]+y[i+2])/(2*r_b**2))**2)
        ds = np.sqrt((x[i+1]-x[i])**2 + (y[i+1] - y[i])**2)

        traps1 = phix[i] * -(x[i]+x[i+1])/(2*r_a**2)/rad_1
        traps2 = phix[i+1] * -(x[i+1]+x[i+2])/(2*r_a**2)/rad_2

        r1x = (x[i] + x[i+1])*0.5 ; r2x = 0.5*(y[i] + y[i+1])
        n1x = -(x[i] + x[i+1])/(2*r_a**2)/rad_1
        n2x = -(y[i] + y[i+1])/(2*r_b**2)/rad_1
        crosses1 = r1x*n2x - r2x*n1x

        r1y = (x[i+1] + x[i+2])*0.5 ; r2y = 0.5*(y[i+1] + y[i+2])
        n1y = -(x[i+1] + x[i+2])/(2*r_a**2)/rad_2
        n2y = -(y[i+1] + y[i+2])/(2*r_b**2)/rad_2
        crosses2 = r1y*n2y - r2y*n1y

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        traps3 = phi6[i] * crosses1
        traps4 = phi6[i+1] * crosses2

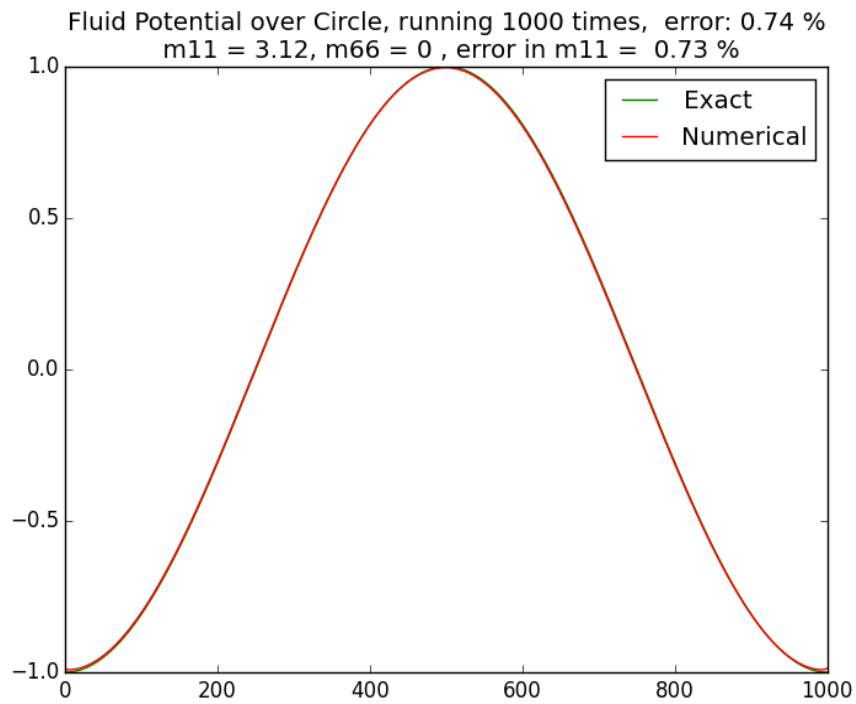
        add_m11 = add_m11 + (traps1 + traps2)*ds * 0.5
        add_m66 = add_m66 + (traps3 + traps4)*ds *0.5
    return add_m11, add_m66

N = 1000
r_a = 1
r_b = 3
equal = r_a - r_b

if equal == 0:
    a_m11,a_m66 = added_mass(N,r_a,r_b)
    error1 = (((r_a**2*np.pi)-float(a_m11))/(r_a**2*np.pi))*100
    print "m11= %.3f m66= %.f,m11 error in percent: %.2f %% " %(a_m11,a_m66,
    thet = np.zeros(N)
    om = 2.0*pi*r_a
    dx = 2*pi/N
    for i in range(N):
        thet[i] = i*dx
    solution = solver1(N,r_a,r_b,direction1 = 11)
    exact_1 = -np.cos(thet)
    diff = max(exact_1 - solution)
    plot(exact_1,"g")
    plot(solution, "r")
    plt.legend(['Exact ', 'Numerical'])
    title("Fluid Potential over Circle, running %.d times, error: %.2f %%
\n m11 = %.2f, m66 = %.f , error in m11 = %.2f %% " %(N,100*diff/max(exact_1), a
    show()

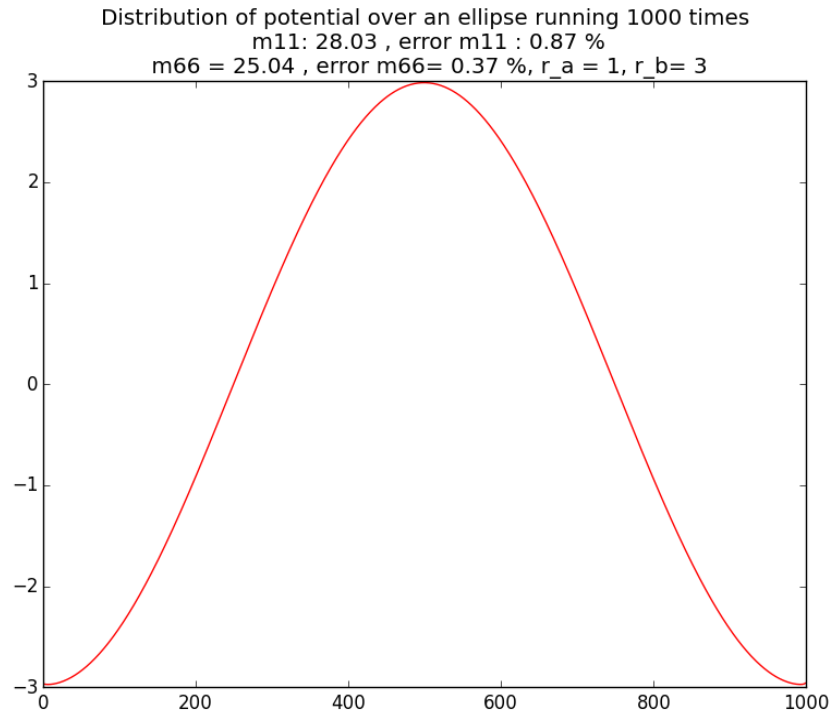
else :
    a_m11,a_m66 = added_mass(N,r_a,r_b)
    error1 = (((r_b**2*np.pi)-a_m11)/(r_b**2*np.pi))*100
    exact6 = (np.pi/8)*(r_a**2-r_b**2)**2
    error6 = (((exact6-a_m66 ) / (exact6)))*100
    print "m11: " ,a_m11 ,"error m11 : %.2f %% , error m66 %.2f %%
running %.d times" %(error1,error6, N)
    plot(solver1(N,r_a,r_b,direction1 = 11), "r")
    title("Distribution of potential over an ellipse running %.d times \n m11
" %(N,a_m11,error1,a_m66,error6,r_a,r_b))
    show()

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This shows a plot of ϕ over a circle, with the added mass and errors calculated.

We can see that the error of ϕ compared to the exact solution is 0.74% and the added mass m_{66} has an error of 0.73% . The added mass m_{66} is of course zero because we have a circle. With these small error i am convinced that the program works for a circle. Next up an ellipse:



This plot is over an ellipse. Again we can see that the errors are very small 0.87% and 0.37%. With both of our geometries having so small error compared to the analytics, I am convinced that the program is correct and the potential and added masses are correctly calculated. I have not calculated the added mass m_{22}, m_{12}, m_{21} as these will all be zero since the geometry only travels laterally.

Since i am not very good in Latex , the code does not look very good copied in. Please let me know if you want the Python code.