

# AST 5900: Problem Set 7

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May 4, 2024

## Problem 1a.

**Answer 1a.** Attached is the code used for implementing the first order upstream finite difference method. Using the code we create the plot in figure 1.

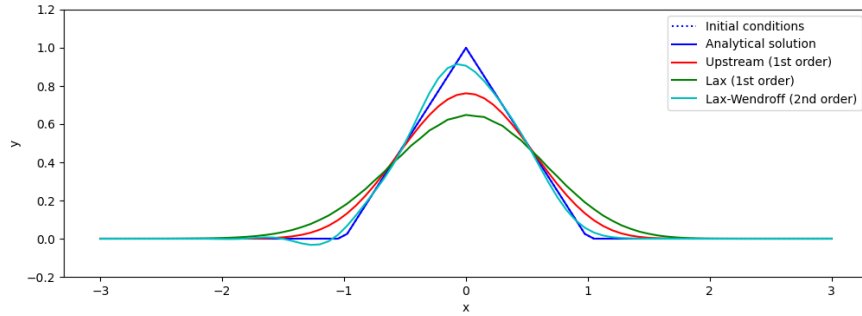


Figure 1: Solving the wave equation using different methods for a triangle wave.

## Problem 1b.

**Answer 1b.** We try different sets of  $nx$  and  $nt$  values until we arrive at the stable-most solution. That happens when  $C_{\max}=1$ . So, the range for stable solutions is  $0 < C < 1$ . Figure 2 shows the solution when  $C=1$ .

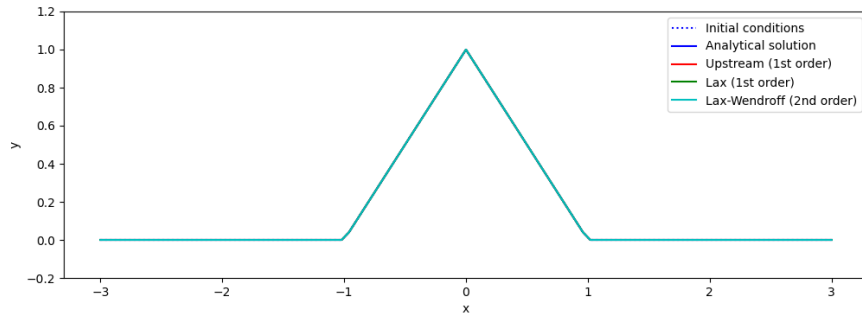


Figure 2: Numerical solutions with  $C = 1$ .

## Problem 1c.

**Answer 1c.** The solution for  $C=1$  is shown in figure 2. We also plot the solutions for  $C < 1$  in figure 3 and for  $C > 1$  in figure 4.

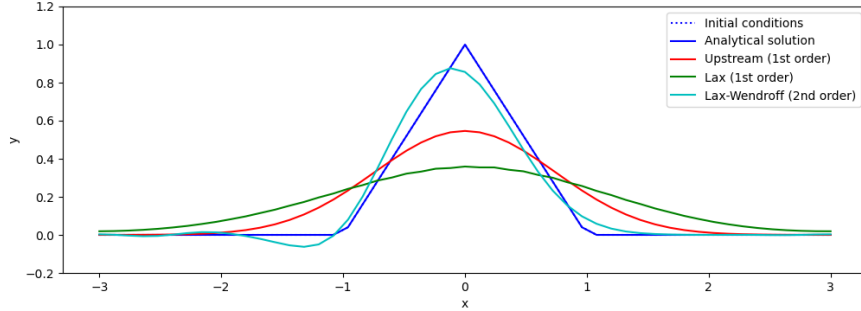


Figure 3: Numerical solutions with  $C = 0.5$ .

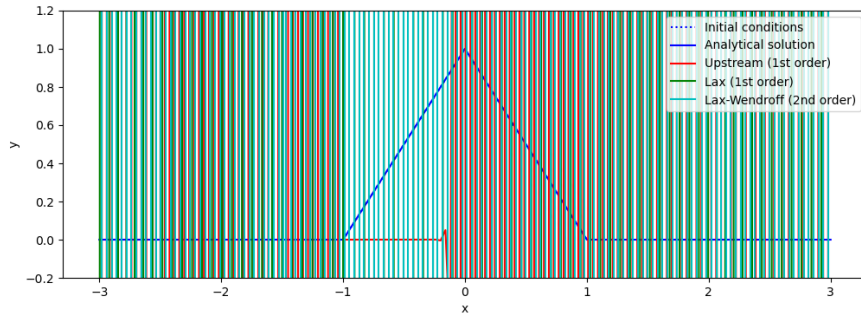


Figure 4: Numerical solutions with  $C = 1.5$ .

## Problem 2a.

**Answer 2a.** We examine the error of the numerical solution. We vary  $nt$  while keeping  $nx$  constant. We produce the absolute error plot in figure 5. We see that reducing the  $C$  number increases the error. In order to decrease  $C$ , we increase  $nt$  values, which, in turn, decreases  $\Delta t$ . So reducing  $\Delta t$  does not improve accuracy.

## Problem 2b.

**Answer 2b.** The modified equation that we got from lecture is

$$\frac{\partial u}{\partial t} + c \frac{\partial u}{\partial x} = (1-C) \frac{c \Delta x}{2} \frac{\partial^2 u}{\partial x^2} + \mathcal{O}(\Delta x^2) \quad (1)$$

The largest order error term is the first term on the left hand side. This term is what defines  $C_{\max}=1$  because  $C=1$  cancels out the term entirely. In this scenario, we arrive at the exact solution, as we saw in our answers above. For  $C > 1$ , we get a negative number which would be nonphysical, so our solution is unstable, as shown above. Thus, our range for  $C$  is between 0 and 1.

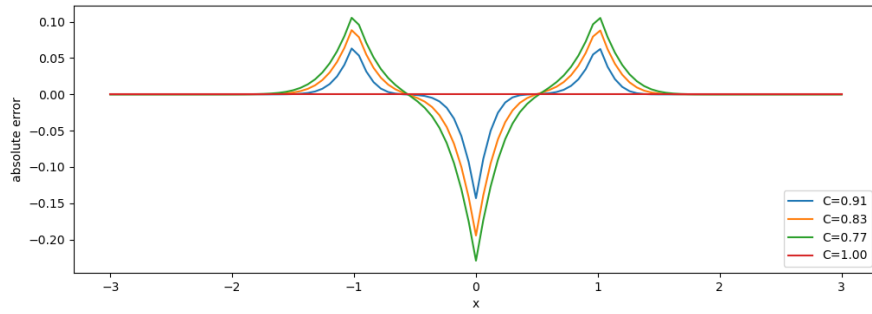


Figure 5: Error plot for various C values.

### Problem 2c.

**Answer 2c.** We keep  $C$  and  $nt$  constant at 0.9 and 110, respectively while varying  $nx$ . We examine how this changes the maximum error in figure 6. We see that the slope is roughly -1 so the error diminishes as we'd expect for a first order method.

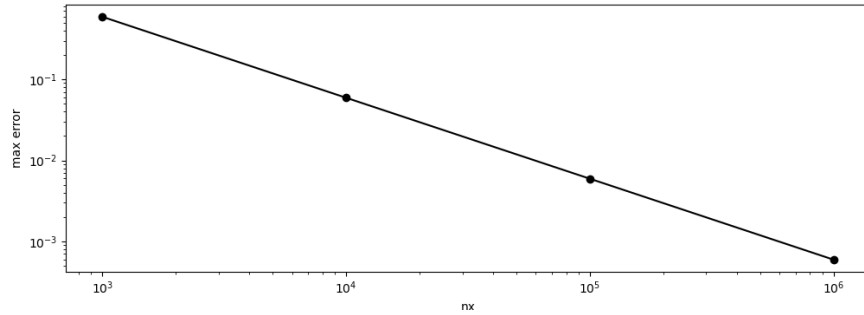


Figure 6: Error plot for various  $nx$  values, keeping  $nt$  and  $C$  constant.

### Problem 3.

**Answer 3.** We follow the code presented [here](#) and adjust our parameters so that our Reynold's number is 20. Reynold's number is given by

$$\text{Re} = \frac{uL}{\nu} \quad (2)$$

where  $u$  is flow speed,  $L$  is length, and  $\nu$  is kinematic viscosity. We choose  $u=1, L=2$ , and  $\nu=0.1$ . We produce the plot shown in figure 7. The code is attached.

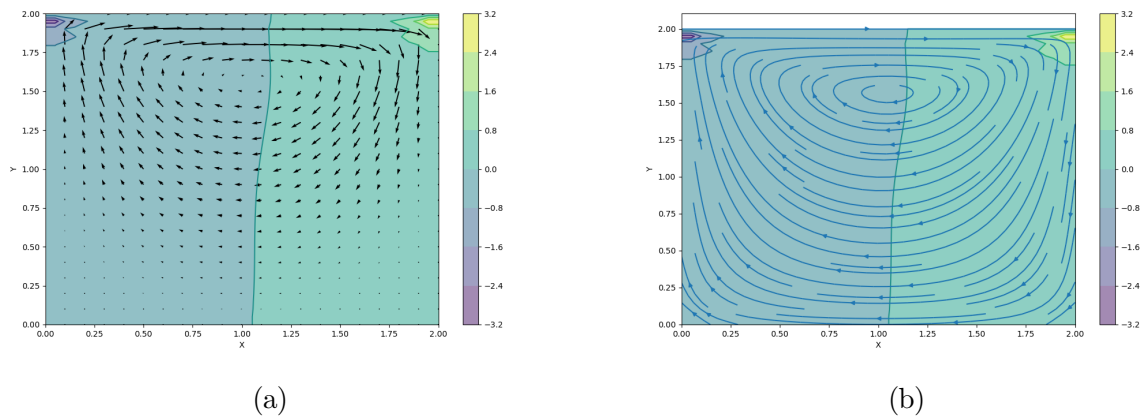


Figure 7: Cavity flow with  $Re = 20$  shown in a (a) vector plot and (b) quiver plot.

## Code

```

1  '''
2  Gilberto Garcia
3  ASTR5900
4  Homework #7
5  24 April 2024
6  '''
7
8  #import required libraries
9  import numpy as np
10 import matplotlib.pyplot as plt
11 from matplotlib import cm
12 from mpl_toolkits.mplot3d import Axes3D
13 plt.rcParams['figure.figsize'] = [12, 4]
14
15
16 #we update the code to match the conditions required for the hw
17 def wave_equation(nx, nt, c, plot=True):
18     # Set spatial parameters
19     xmin = -3.0
20     xmax = 3.0
21     dx = (xmax - xmin) / nx
22     print(f"{xmin} <= x <= {xmax} (Spatial domain)")
23     print(f"dx = {dx} (Spatial step size)")
24
25     x_points = np.linspace(xmin, xmax, nx + 1)
26
27     # Set temporal parameters
28     tmin = 0.
29     tmax = 3.
30     dt = (tmax - tmin) / nt
31     print(f"{tmin} <= t <= {tmax} (Temporal domain)")

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32     print(f"dt = {dt} (Temporal step size)")
33
34     t_points = np.linspace(tmin, tmax, nt + 1)
35
36     # Set speed for linear wave equation
37     print(f"c = {c} (Wave speed)")
38
39     # Calculate Courant Number
40     C = c * dt / dx
41     #C = 0.9 #uncomment for 2C
42     print(f"C = {C} (Courant Number)")
43
44     # Define numerical solver step functions
45     def upstream_step(u):
46         # u_old = u.copy()
47         # for x_index in range(1, nx + 1):
48         #     u[x] = u_old[x] - C * (u_old[x] - o_old[x - 1])
49         u[1:] = u[1:] - C * (u[1:] - u[:-1])
50
51     def upstream_periodic_step(u):
52         up = u.copy()
53         u[1:] = up[1:] - C * (up[1:] - up[:-1])
54         # Periodic boundary condition at x = xmin
55         u[0] = up[0] - C * (up[0] - up[-2])
56         # Periodic boundary condition at x = xmax
57         u[-1] = u[0]
58
59     def lax_step(u):
60         u[1:-1] = 0.5 * (u[2:] + u[:-2]) - 0.5 * C * (u[2:] - u[:-2])
61
62     def lax_periodic_step(u):
63         up = u.copy()
64         u[1:-1] = 0.5 * (up[2:] + up[:-2]) - 0.5 * C * (up[2:] - up[:-2])
65         # Periodic boundary condition at x = xmin
66         u[0] = 0.5 * (up[1] + up[-2]) - 0.5 * C * (up[1] - up[-2])
67         # Periodic boundary condition at x = xmax
68         u[-1] = u[0]
69
70     def lax_wendroff_step(u):
71         u[1:-1] = u[1:-1] -
72             0.5 * C * (u[2:] - u[:-2]) + 0.5 * C ** 2 * (u[2:] - 2 * u[1:-1] + u[:-2])
73
74     def lax_wendroff_periodic_step(u):
75         # Periodic boundary condition by wrapping at both ends
76         up = np.concatenate(( [u[-2]], u, [u[1]] ))
77         u[:] = up[1:-1] - 0.5 *
78             C * (up[2:] - up[:-2]) + 0.5 * C ** 2 * (up[2:] - 2 * up[1:-1] + up[:-2])
79         return 0
80
81     def triangle(x,t=0):

```

```

80     if -1.0 + c * t <= x <= 1.0 + c * t:
81         return 1.0 - abs(x - c*t)
82     else:
83         return 0
84
85 def periodic_triangle(x,t):
86     x_wrapped = x
87     while x_wrapped < xmin + c * t:
88         x_wrapped += (xmax - xmin)
89     while x_wrapped > xmax + c * t:
90         x_wrapped -= (xmax - xmin)
91     return triangle(x_wrapped, t)
92
93 # Save the initial and final analytical solutions
94 u0 = np.array([periodic_triangle(x, tmin) for x in x_points])
95 uf = np.array([periodic_triangle(x, tmax) for x in x_points])
96
97 # Initialize numerical solutions
98 u_upstream = u0.copy()
99 u_lax = u0.copy()
100 u_lax_wendroff = u0.copy()
101
102 # Advance numerical solution by nt timesteps
103 for t in t_points[1:]:
104     upstream_periodic_step(u_upstream)
105     lax_periodic_step(u_lax)
106     lax_wendroff_periodic_step(u_lax_wendroff)
107
108 if plot == True:
109     # Plot analytical and numerical solutions
110     plt.plot(x_points, u0, 'b:', label="Initial conditions")
111     plt.plot(x_points, uf, 'b-', label="Analytical solution")
112     plt.plot(x_points, u_upstream, 'r-', label="Upstream (1st order)")
113     plt.plot(x_points, u_lax, 'g-', label="Lax (1st order)")
114     plt.plot(x_points, u_lax_wendroff, 'c-', label="Lax-Wendroff (2nd order)")
115     plt.ylim(-0.2, 1.2)
116     plt.legend()
117     plt.xlabel('x')
118     plt.ylabel('y')
119     plt.show()
120
121 #calculate relative error:
122 abs_err = (u_upstream - uf)
123 #return x points, analytical, and 1st order upstream numerical soln
124 return x_points,abs_err,C
125
126 #1a - centering the propogation waves
127 wave_equation(nx = 80, nt = 100, c = 2)
128 #1b - we play around with nx and nt to find where Cmax is not stable
129 wave_equation(nx = 100, nt = 100, c = 2)

```

```

130 #we find that c_max of 1 is the upper limit for stability.
131
132 #1c - c_max for less than, equal to and greater than Cmax
133 #less than
134 wave_equation(nx = 50, nt = 100, c = 2)
135 #equal to
136 wave_equation(nx = 100, nt = 100, c = 2)
137 #greater than
138 wave_equation(nx = 150, nt = 100, c = 2)
139
140 #2a
141 print()
142 print('2a')
143 xpts1,abserr1,C1 = wave_equation(nx = 100, nt = 110, c = 2,plot=False)
144 xpts2,abserr2,C2 = wave_equation(nx = 100, nt = 120, c = 2,plot=False)
145 xpts3,abserr3,C3 = wave_equation(nx = 100, nt = 130, c = 2,plot=False)
146 xpts4,abserr4,C4 = wave_equation(nx = 100, nt = 100, c = 2,plot=False)
147
148
149 plt.plot(xpts1,abserr1,label='C={0:.2f}'.format(C1))
150 plt.plot(xpts2,abserr2,label='C={0:.2f}'.format(C2))
151 plt.plot(xpts3,abserr3,label='C={0:.2f}'.format(C3))
152 plt.plot(xpts4,abserr4,label='C={0:.2f}'.format(C4))
153 plt.ylabel('absolute error')
154 plt.xlabel('x')
155 plt.legend(loc='lower right')
156 plt.show()
157
158 #no, reducing delta t while keeping delta x constant does not improve accuracy.
159 # bigger nt means smaller delta t but also bigger error so less accuracy.
160
161 #2c - we will need to hard code C =0.9 in wave_equation()
162 #keep C at 0.9. we vary the nx values
163
164 nx_vals = [1_000,10_000,100_000,1_000_000]
165 xpts,err1,c = wave_equation(nx = nx_vals[0], nt = 110, c = 2,plot=False)
166 xpts,err2,c = wave_equation(nx = nx_vals[1], nt = 110, c = 2,plot=False)
167 xpts,err3,c = wave_equation(nx = nx_vals[2], nt = 110, c = 2,plot=False)
168 xpts,err4,c = wave_equation(nx = nx_vals[3], nt = 110, c = 2,plot=False)
169
170 #make a list of maximum errors:
171 max_err = [max(err1),max(err2),max(err3),max(err4)]
172
173 #plot nx vs max errors:
174
175 plt.loglog(nx_vals,max_err,'k')
176 plt.scatter(nx_vals,max_err,color='k')
177 plt.xlabel('nx')
178 plt.ylabel('max error')
179 plt.show()

```

```

180
181
182 #####
183 ### 3 #####
184 #####
185
186
187 #take the code from
188 #https
189     ://nbviewer.org/github/barbagroup/CFDPython/blob/master/lessons/14_Step_11.ipynb
190
191 nx = 41
192 ny = 41
193 nt = 500
194 nit = 50
195 c = 1
196 dx = 2 / (nx - 1)
197 dy = 2 / (ny - 1)
198 x = np.linspace(0, 2, nx)
199 y = np.linspace(0, 2, ny)
200 X, Y = np.meshgrid(x, y)
201
202 rho = 1
203 nu = .1
204 dt = .001
205
206 u = np.zeros((ny, nx))
207 v = np.zeros((ny, nx))
208 p = np.zeros((ny, nx))
209 b = np.zeros((ny, nx))
210
211 def build_up_b(b, rho, dt, u, v, dx, dy):
212
213     b[1:-1, 1:-1] = (rho * (1 / dt *
214         ((u[1:-1, 2:] - u[1:-1, 0:-2]) /
215         (2 * dx) + (v[2:, 1:-1] - v[0:-2, 1:-1]) / (2 * dy)) -
216         ((u[1:-1, 2:] - u[1:-1, 0:-2]) / (2 * dx))**2 -
217         2 * ((u[2:, 1:-1] - u[0:-2, 1:-1]) / (2 * dy) *
218         (v[1:-1, 2:] - v[1:-1, 0:-2]) / (2 * dx)) -
219         ((v[2:, 1:-1] - v[0:-2, 1:-1]) / (2 * dy))**2))
220
221     return b
222
223
224 def pressure_poisson(p, dx, dy, b):
225     pn = np.empty_like(p)
226     pn = p.copy()
227
228     for q in range(nit):

```



```

229     pn = p.copy()
230     p[1:-1, 1:-1] = (((pn[1:-1, 2:] + pn[1:-1, 0:-2]) * dy**2 +
231                      (pn[2:, 1:-1] + pn[0:-2, 1:-1]) * dx**2) /
232                      (2 * (dx**2 + dy**2))) -
233                      dx**2 * dy**2 / (2 * (dx**2 + dy**2)) *
234                      b[1:-1, 1:-1])
235
236     p[:, -1] = p[:, -2] # dp/dx = 0 at x = 2
237     p[0, :] = p[1, :] # dp/dy = 0 at y = 0
238     p[:, 0] = p[:, 1] # dp/dx = 0 at x = 0
239     p[-1, :] = 0 # p = 0 at y = 2
240
241     return p
242
243
244 def cavity_flow(nt, u, v, dt, dx, dy, p, rho, nu):
245     un = np.empty_like(u)
246     vn = np.empty_like(v)
247     b = np.zeros((ny, nx))
248
249     for n in range(nt):
250         un = u.copy()
251         vn = v.copy()
252
253         b = build_up_b(b, rho, dt, u, v, dx, dy)
254         p = pressure_poisson(p, dx, dy, b)
255
256         u[1:-1, 1:-1] = (un[1:-1, 1:-1] -
257                          un[1:-1, 1:-1] * dt / dx *
258                          (un[1:-1, 1:-1] - un[1:-1, 0:-2]) -
259                          vn[1:-1, 1:-1] * dt / dy *
260                          (un[1:-1, 1:-1] - un[0:-2, 1:-1]) -
261                          dt / (2 * rho * dx) * (p[1:-1, 2:] - p[1:-1, 0:-2]) +
262                          nu * (dt / dx**2 *
263                               (un[1:-1, 2:] - 2 * un[1:-1, 1:-1] + un[1:-1, 0:-2]) +
264                               dt / dy**2 *
265                               (un[2:, 1:-1] - 2 * un[1:-1, 1:-1] + un[0:-2, 1:-1])))
266
267         v[1:-1, 1:-1] = (vn[1:-1, 1:-1] -
268                          un[1:-1, 1:-1] * dt / dx *
269                          (vn[1:-1, 1:-1] - vn[1:-1, 0:-2]) -
270                          vn[1:-1, 1:-1] * dt / dy *
271                          (vn[1:-1, 1:-1] - vn[0:-2, 1:-1]) -
272                          dt / (2 * rho * dy) * (p[2:, 1:-1] - p[0:-2, 1:-1]) +
273                          nu * (dt / dx**2 *
274                               (vn[1:-1, 2:] - 2 * vn[1:-1, 1:-1] + vn[1:-1, 0:-2]) +
275                               dt / dy**2 *
276                               (vn[2:, 1:-1] - 2 * vn[1:-1, 1:-1] + vn[0:-2, 1:-1])))
277
278         u[0, :] = 0

```

```

279     u[:, 0] = 0
280     u[:, -1] = 0
281     u[-1, :] = 1    # set velocity on cavity lid equal to 1
282     v[0, :] = 0
283     v[-1, :] = 0
284     v[:, 0] = 0
285     v[:, -1] = 0
286
287
288     return u, v, p
289
290
291 u = np.zeros((ny, nx))
292 v = np.zeros((ny, nx))
293 p = np.zeros((ny, nx))
294 b = np.zeros((ny, nx))
295 nt = 700
296 u, v, p = cavity_flow(nt, u, v, dt, dx, dy, p, rho, nu)
297
298
299
300 fig = plt.figure(figsize=(11,7), dpi=100)
301 # plotting the pressure field as a contour
302 plt.contourf(X, Y, p, alpha=0.5, cmap=cm.viridis)
303 plt.colorbar()
304 # plotting the pressure field outlines
305 plt.contour(X, Y, p, cmap=cm.viridis)
306 # plotting velocity field
307 plt.quiver(X[::2, ::2], Y[::2, ::2], u[::2, ::2], v[::2, ::2])
308 plt.xlabel('X')
309 plt.ylabel('Y')
310 plt.show()
311
312
313 fig = plt.figure(figsize=(11, 7), dpi=100)
314 plt.contourf(X, Y, p, alpha=0.5, cmap=cm.viridis)
315 plt.colorbar()
316 plt.contour(X, Y, p, cmap=cm.viridis)
317 plt.streamplot(X, Y, u, v)
318 plt.xlabel('X')
319 plt.ylabel('Y')
320 plt.show()

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