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In [ ]: import numpy as np
        from scipy.integrate import trapz
        import matplotlib.pyplot as plt
        class CodeWrapper:
            def __init__(self, X1, X2, Y1, Y2, m, n, hx, hy, m_values, gamma_values, fV_values
                self.X1 = X1
                self.X2 = X2
                 self.Y1 = Y1
                 self.Y2 = Y2
                self.m = m
                 self.n = n
                 self.hx = hx
                 self.hy = hy
                 self.x = np.linspace(self.X1, self.X2, self.n)
                 self.y = np.linspace(self.Y1, self.Y2, self.m)
                 self.m_values = m_values
                 self.gamma_values = gamma_values
                 self.fV_values = fV_values
                self.B1 = B1
                 self.B2 = B2
                 self.XD = XD
            def f(self, x, M, G, F):
                return F**2 * ((M**4 + M**2 * G**2) / (24 * np.pi * ((x - M**2)**2 + M**2 * G*
            def define_functions(self):
                 self.f1 = self.f(self.x, self.m_values[0], self.gamma_values[0], 1)
                 self.f2 = self.f(self.x, self.m_values[1], self.gamma_values[1], 1)
                 self.f3 = self.f(self.x, self.m_values[2], self.gamma_values[2], 1)
                 self.f4 = self.f(self.x, self.m_values[3], self.gamma_values[3], 1)
                 self.F = self.f1 + self.f2 + self.f3 + self.f4
            def calculate_F1_F2(self):
                 self.F1 = self.f1
                 self.F2 = self.x - self.x
            def initialize_arrays(self):
                 self.G1 = np.zeros((self.m,))
                 self.G2 = np.zeros((self.m,))
                 self.G3 = np.zeros((self.m,))
                 self.g_exact = np.zeros((self.n,))
                 self.g_delta = np.zeros((self.n,))
                 self.U_exact = np.zeros((self.n,))
                 self.G_exact = np.zeros((self.n,))
                 self.G_delta = np.zeros((self.n,))
                 self.Aphi = np.zeros((self.m, self.n))
                 self.A = np.zeros((self.n, self.n))
                 self.ba = np.zeros(self.n)
                 self.M = np.zeros((self.n, self.n))
                 self.M1 = np.zeros((self.n, self.n))
                 self.H1 = np.zeros((self.n, self.n))
                 self.alpha = np.zeros(20)
                 self.ua = np.zeros((self.n - 2, 20))
                 self.Ua = np.zeros((self.n, 20))
                 self.fa = np.zeros((self.n, 20))
                 self.res = np.zeros(20)
                 self.err = np.zeros(20)
                 self.L = np.zeros(20)
                 self.LH = np.zeros(20)
                 self.V = np.zeros((20, 1))
                 self.Rho = np.zeros((20, 1))
            def calculate_G1_G2_G3(self):
                 for i in range(self.m):
                     self.G1[i] = trapz(1 / (self.y[i] - self.x) * self.F1, self.x)
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self.G2[i] = trapz(1 / (self.y[i] - self.x) * self.F2, self.x)
               self.G3[i] = trapz(1 / (self.y[i] - self.x) * self.V_exact, self.x)
def calculate_f_exact_f_delta(self):
       self.f_exact = self.B1 * self.F1 + self.B2 * self.F2
       self.f_delta = self.XD * self.B1 * self.F1 + self.XD * self.B2 * self.F2
def calculate_V_exact(self):
       MM = self.f_delta[0]
       NN = self.f_delta[-1]
       self.V_exact = (MM * (self.x - self.X2)) / (self.X1 - self.X2) + 
               (NN * (self.x - self.X1)) / (self.X2 - self.X1)
def calculate_q_exact_q_delta_U_exact_G_exact_G_delta(self):
       self.q_exact = self.B1 * self.G1 + self.B2 * self.G2
       self.g_delta = self.XD * self.B1 * self.G1 + self.XD * self.B2 * self.G2
       self.U_exact = self.f_exact - self.V_exact
       self.G_exact = self.q_exact - self.G3
       self.G_delta = self.g_delta - self.G3
def calculate_Aphi_A_ba(self):
       for j in range(1, self.n - 1):
               self.Aphi[:, j] = 1 / self.hx * ((self.y - self.x[j - 1]) * np.log(self.y)
                                                                            + (self.x[j - 1] - self.x[j]) + (self.y
                                                                            np.log(self.y - self.x[j + 1])
                                                                             - 2 * (self.y - self.x[j]) * np.log(self.
       self.Aphi[:, 0] = np.log(self.y - self.x[0]) + 1 / self.hx * ((self.y - self.x
                                                                                                                            - (self.y - self
       self.Aphi[:, -1] = -np.log(self.y - self.x[-1]) + 1 / self.hx * (-(self.y - self.x[-
                                                                                                                                 + (self.y - s
       for i in range(self.n):
               for j in range(self.n):
                      self.A[i, j] = trapz(self.Aphi[:, i] * self.Aphi[:, j], self.y)
               self.ba[i] = trapz(self.Aphi[:, i] * self.G_delta, self.y)
def calculate_M_M1_H1(self):
       self.M[0, 0] = 4
       self.M[0, 1] = 2
       for i in range(1, self.n - 1):
               self.M[i, i] = 8
               self.M[i, i - 1] = 2
               self.M[i, i + 1] = 2
       self.M[-1, -1] = 4
       self.M[-1, -2] = 2
       self.M = self.hx / 12 * self.M
       self.M1[0, 0] = 1
       self.M1[0, 1] = -1
       for i in range(1, self.n - 1):
               self.M1[i, i] = 2
               self.M1[i, i - 1] = -1
               self.M1[i, i + 1] = -1
       self.M1[-1, -1] = 1
       self.M1[-1, -2] = -1
       self.M1 = 1 / self.hx * self.M1
       self.H1 = self.M + self.M1
def remove_boundary_conditions(self):
       self.A = np.delete(self.A, [0, self.n - 2], axis=0)
       self.A = np.delete(self.A, [0, self.n - 2], axis=1)
       self.ba = np.delete(self.ba, [0, self.n - 2])
       self.H1 = np.delete(self.H1, [0, self.n - 2], axis=0)
       self.H1 = np.delete(self.H1, [0, self.n - 2], axis=1)
def iterate(self):
       for k in range(20):
               self.alpha[k] = 10**(-k-1)
               self.ua[:, k] = np.linalg.inv(
                      self.A + self.alpha[k] * self.H1) @ self.ba
               self.Ua[:, k] = np.concatenate(([0], self.ua[:, k], [0]))
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self.err[k] = np.sqrt(
                         trapz((self.fa[:, k] - self.f_exact)**2, self.x))
                    self.L[k] = np.sqrt(trapz(self.fa[:, k]**2, self.x)) * self.res[k]
                    self.LH[k] = np.sqrt(
                         trapz(self.fa[:, k]**2 + (np.gradient(self.fa[:, k]))**2, self.x)) * 
                     # Plot results (uncomment this block if you want to plot)
                    plt.figure(1)
                     # plt.figure(1, dpi=1000)
                    plt.subplot(4, 5, k + 1)
                    plt.plot(self.x, self.f_exact, self.x, self.fa[:, k], linewidth=3)
                    plt.title(str(self.alpha[k]))
            def find_indices(self):
                 kl = np.argmin(np.abs(self.L))
                 klh = np.argmin(np.abs(self.LH))
                for i in range(20):
                     self.V[i] = np.linalq.norm(self.Aphi @ np.linalq.inv(self.Aphi.T @
                                                                           self.Aphi + self.alpł
                         np.trace((np.eye(len(self.y)) - self.Aphi @ np.linalg.inv(self.Aphi.T
                                                                                    self.Aphi +
                for i in range(20):
                    self.Rho[i] = np.linalq.norm(self.fa[:, i] / (self.alpha[i] *
                                                                    (self.Aphi.T @ self.Aphi + s
                Vm = np.min(self.V)
                Vi = np.argmin(self.V)
                Rm = np.min(self.Rho)
                Ri = np.argmin(self.Rho)
                return [['L-curve+L^2', 'L-curve+H^1', 'GCV', 'Approx. Optimal'],
                         [kl + 1, klh + 1, Vi + 1, Ri + 1]] # Add 1 to indices to match MATLAL
            def run(self):
                 self.initialize_arrays()
                 self.define_functions()
                 self.calculate_F1_F2()
                 self.calculate_f_exact_f_delta()
                 self.calculate_V_exact()
                 self.calculate_G1_G2_G3()
                 self.calculate_g_exact_g_delta_U_exact_G_exact_G_delta()
                 self.calculate_Aphi_A_ba()
                 self.calculate M M1 H1()
                 self.remove_boundary_conditions()
                 self.iterate()
                return self.find_indices()
In [ ]: # if __name__ == "__main__":
        X1 = 0
        X2 = 2
        Y1 = X2 + 0.01
        Y2 = 30 + 0.01
        n = int((X2 - X1) * 1000 + 1)
        m = int((Y2 - Y1) * 100 + 1)
        hx = (X2 - X1) / (n - 1)
        hy = (Y2 - Y1) / (m - 1)
        m_{values} = [0.78, 1.46, 1.72, 1.90]
        gamma_values = [0.15, 0.4, 0.25, 0.1]
        fV_values = [0.22, 0.19, 0.14, 0.14]
        B1, B2 = 1, 2
        XD = 1.01
        code_wrapper = CodeWrapper(
            X1, X2, Y1, Y2, m, n, hx, hy, m_values, gamma_values, fV_values, B1, B2, XD)
In [ ]: | code_wrapper.initialize_arrays()
        code_wrapper.define_functions()
```

 $Selt.ta[:, K] = Selt.ua[:, K] + Selt.v_exact$ 

trapz((self.Aphi @ self.fa[:, k] - self.G\_delta)\*\*2, self.y))

self.res[k] = np.sqrt(

```
In [ ]:
           code_wrapper.calculate_F1_F2()
           code_wrapper.calculate_f_exact_f_delta()
           code_wrapper.calculate_V_exact()
In [ ]:
           code_wrapper.calculate_G1_G2_G3()
In [ ]:
           code_wrapper.calculate_q_exact_g_delta_U_exact_G_exact_G_delta()
           code_wrapper.calculate_Aphi_A_ba()
In [ ]:
In [ ]:
           code_wrapper.calculate_M_M1_H1()
           code_wrapper.remove_boundary_conditions()
In [ ]:
           code_wrapper.iterate()
                                             0.01
                                                              0.001
                                                                               0.0001
                           0.1
                                                                                                   1e-05
                                  0.25
                                                    0.25
                                                                                        0.25
               0.25
                                                                      0.25
                                            <del>1e-07<sup>0.0</sup>b</del>
                                                                                <del>1e-09<sup>0</sup> q</del>b
               0.00
                          <del>1e-00<sup>0.0</sup>0</del>
                                                               <del>1e-08<sup>0.00</sup></del>
                                                                                                   <del>1e-10</del>
               0.25
                                  0.25
                                                    0.25
                                                                      0.25
                                                                                        0.25
                                            <del>1e-12<sup>0.0</sup>b</del>
                                                                                <del>1e-14<sup>0.0</sup>b</del>
                                                              <del>1e-13<sup>0.0</sup>b</del>
               0.00
                          <del>1e-11<sup>0.0</sup>b</del>
                                                                                                   <del>1e-15</del>
```

2.b

o.b

2 0

1e-19

<del>2.</del>5

o.b

-2.5

2

0

.e8

1e-20

2

e10

1e-18<sub>-2.5</sub>

.e7



0.25

0.0þ

2.b

o.b

2 0

-2.5

0.25

0.00

5000

-5000

0

0.25

2 0

le6

<del>1e-16<sup>0.0</sup>b</del>