## ECSE 543 Assignment 2

Sean Stappas 260639512

November  $20^{th}$ , 2017

## Contents

1	Finite Element Triangles				
<b>2</b>	Finite Element Coaxial Cable	2			
	2.a Mesh	2			
	2.b Electrostatic Potential	2			
	2.c Capacitance	2			
3	Conjugate Gradient Coaxial Cable	3			
	3.a Positive Definite Test	3			
	3.b Matrix Solution	3			
	3.c Residual Norm	3			
	3.d Potential Comparison	4			
	3.e Capacitance Computation	4			
A	opendix A Code Listings	5			
A	opendix B Output Logs	14			
Α	opendix C Simple2D Data Files	16			

#### Introduction

### 1 Finite Element Triangles

The equation for the  $\alpha$  parameter for a general vertex i of a finite element triangle can be seen in Equation (1), where i+1 and i+2 implicitly wraps around when exceeding 3.

$$\alpha_{i}(x,y) = \frac{1}{2A} \left[ (x_{i+1}y_{i+2} - x_{i+2}y_{i+1}) + (y_{i+1} - y_{i+2})x + (x_{i+2} - x_{i+1})y \right]$$

$$(1)$$

Using Equation (1), we can solve for the entries of the local S matrix, as shown in Equation (2). This was used in the program to compute every entry for both example triangles.

$$S_{ij} = \int_{\Delta_e} \nabla \alpha_i \cdot \nabla \alpha_j dS$$

$$= \frac{1}{4A} \left[ (y_{i+1} - y_{i+2})(y_{j+1} - y_{j+2}) + (x_{i+2} - x_{i+1})(x_{j+2} - x_{j+1}) \right]$$
(2)

The local S matrix for the first triangle can be seen in Equation (3).

$$S_1 = \begin{bmatrix} 0.5 & -0.5 & 0.0 \\ -0.5 & 1.0 & -0.5 \\ 0.0 & -0.5 & 0.5 \end{bmatrix}$$
 (3)

The local S matrix for the second triangle can be seen in Equation (4).

$$S_2 = \begin{bmatrix} 1.0 & -0.5 & -0.5 \\ -0.5 & 0.5 & 0.0 \\ -0.5 & 0.0 & 0.5 \end{bmatrix} \tag{4}$$

The disjoint S matrix is then given by the following:

$$S_{dis} = \begin{bmatrix} 0.5 & -0.5 & 0.0 & 0 & 0 & 0 \\ -0.5 & 1.0 & -0.5 & 0 & 0 & 0 \\ 0.0 & -0.5 & 0.5 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1.0 & -0.5 & -0.5 \\ 0 & 0 & 0 & -0.5 & 0.5 & 0.0 \\ 0 & 0 & 0 & -0.5 & 0.0 & 0.5 \end{bmatrix}$$

The connectivity matrix C is given by Equation (5).

$$C = \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \\ 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 \end{bmatrix}$$
 (5)

The global matrix S is then given by Equation (6).

$$S = C^T S_{dis} C^T (6)$$

Using Equations (5) and (6), we can solve for the global S matrix, giving the value shown in Equation (7), which is computed by the finite\_element\_triangles.py script shown in Listing 3.

$$S = \begin{bmatrix} 1.0 & -0.5 & 0.0 & -0.5 \\ -0.5 & 1.0 & -0.5 & 0.0 \\ 0.0 & -0.5 & 1.0 & -0.5 \\ -0.5 & 0.0 & -0.5 & 1.0 \end{bmatrix}$$
 (7)

# 2 Finite Element Coaxial Cable

#### 2.a Mesh

The mesh to be used by the by SIMPLE2D program is generated the finite\_element\_mesh\_generator.py script shown in Listing 5. This input and output files of the SIMPLE2D program are shown in Listings 13 and 14 of Appendix C.

#### 2.b Electrostatic Potential

Based on the results from the SIMPLE2D program, the potential at  $(0.06,\,0.04)$  is  $5.5263\,\mathrm{V}$ . This corresponds to node 16 in the mesh arrangement we created.

#### 2.c Capacitance

The finite element functional equation for two conjoint finite element triangles forming a square i can be seen in Equation (8).

$$W_i = \frac{1}{2} U_{con_i}^T S U_{con_i} \tag{8}$$

where S is given in Equation (7) and  $U_{con_i}$  is the conjoint potential vector for square i, giving the potential at the four corners of the square defining the combination of two finite element triangles. This can be seen in Equation (9).

$$U_{con} = \begin{bmatrix} U_{i_1} \\ U_{i_2} \\ U_{i_3} \\ U_{i_4} \end{bmatrix}$$
 (9)

To find the total energy function W of the mesh, we must add the contributions from each square and multiply by 4, since our mesh is one quarter of the entire coaxial cable. This yields Equation (10).

$$W = 4\sum_{i}^{N} W_{i} = 2\sum_{i}^{N} U_{con_{i}}^{T} SU_{con_{i}}$$
 (10)

where N is the number of finite difference squares in the mesh.

Note that W is not equal to the energy. The relation between the energy per unit length E and W is shown in Equation (11).

$$E = \epsilon_0 W \tag{11}$$

We then know that the energy per unit length E is related to the capacitance per unit length C as shown in Equation (12).

$$E = \frac{1}{2}CV^2 \tag{12}$$

where V is the voltage across the coaxial cable.

Combining Equations (8) and (10) to (12), we obtain an expression for the capacitance per unit length which can be easily calculated, as shown in Equation 13.

$$C = \frac{2E}{V^2} = \frac{4\epsilon_0}{V^2} \sum_{i}^{N} U_{con_i}^T SU_{con_i}$$
 (13)

The capacitance per unit length is computed as  $5.2137 \times 10^{-11} \, \mathrm{F/m}$  by the finite\_element\_capacitance.py script shown in Listing 6 with output shown in Listing 11.

## 3 Conjugate Gradient Coaxial Cable

#### 3.a Positive Definite Test

To form the A matrix, we must consider all the free nodes in the mesh. The potential at the non-boundary free nodes is given by Equation (14).

$$-4\phi_{i,j} + \phi_{i+1,j} + \phi_{i-1,j} + \phi_{i,j+1} + \phi_{i,j-1} = 0$$
 (14)

The free nodes along a boundary must satisfy the Neumann boundary condition for symmetry. Since our quarter-mesh is the bottom left corner of the overall mesh, these boundary nodes defining planes of symmetry are along the top and the right. The Neumann boundary condition for the top nodes is given by Equation (15) and that for the right nodes is given by Equation (16).

$$\phi_{i,j+1} - \phi_{i,j-1} = 0 \tag{15}$$

$$\phi_{i+1,j} - \phi_{i-1,j} = 0 \tag{16}$$

Now, the simplified potential for boundary free nodes can be calculated, as seen in Equations (17) and (18).

$$-4\phi_{i,j} + \phi_{i+1,j} + \phi_{i-1,j} + 2\phi_{i,j-1} = 0$$
 (17)

$$-4\phi_{i,j} + 2\phi_{i-1,j} + \phi_{i,j+1} + \phi_{i,j-1} = 0$$
 (18)

The non-free nodes are fixed by the potentials of the conductors, i.e., 15 V and 0 V.

With Equations (14), (17) and (18), we can form the A matrix from every mesh node. This is done in finite\_difference\_mesh\_generator.py, as shown in Listing 8. The output A matrix can be seen in Listing 12.

If the matrix A is not positive definite, one can simply multiply both sides of the Ax = b equation by  $A^T$ , forming a new equation  $A^TAx = A^Tb$ . This is equivalent to A'x = b', where  $b' = A^Tb$  and  $A' = A^TA$ . Here, A' is now positive definite.

In our case, the matrix A is indeed not positive definite, and multiplying by  $A^T$  made it positive definite. The before and after positive definite test can be seen in Listing 12.

#### 3.b Matrix Solution

The matrix equation to be solved can be seen in Equation (19), where A is positive-definite matrix generated previously,  $\phi_c$  is the unknown potential vector and b contains the initial potential values along the boundaries.

$$A\phi_c = b \tag{19}$$

#### 3.c Residual Norm

Consider a vector  $\mathbf{v} = \{v_1, \dots, v_n\}$ . The infinity norm  $\|\mathbf{v}\|_{\infty}$  of  $\mathbf{v}$  is given by the maximum absolute element of  $\mathbf{v}$ , as shown in Equation (20).

$$\|\mathbf{v}\|_{\infty} = \max\{|v_1|, \dots, |v_n|\} \tag{20}$$

Similarly, the 2-norm  $\|\mathbf{v}\|_2$  of  $\mathbf{v}$  is given by Equation (21).

$$\|\mathbf{v}\|_2 = \sqrt{\sum_{i=1}^n v_i^2} \tag{21}$$

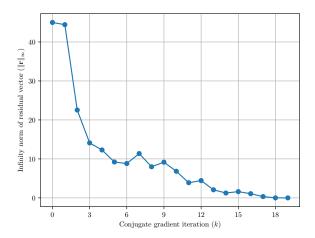


Figure 1: Value of the infinity norm of the residual vector versus iterations of the conjugate gradient algorithm.

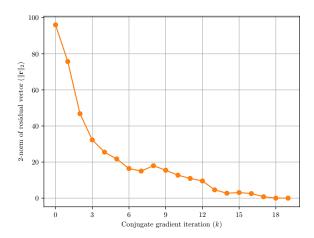


Figure 2: Value of the 2-norm of the residual vector versus iterations of the conjugate gradient algorithm.

#### 3.d Potential Comparison

#### 3.e Capacitance Computation

The capacitance can be calculated in the same way as in Question 2(c), i.e., with Equation (13). The node values must simply be mapped to same mesh used in the finite difference context.

## A Code Listings

```
Listing 1: Custom matrix package (matrices.py).
    from __future__ import division
2
    import copy
3
4
    import csv
    from ast import literal_eval
    import math
9
    class Matrix:
10
11
        def __init__(self, data):
             self.data = data
12
13
             self.num_rows = len(data)
             self.num_cols = len(data[0])
14
15
16
         def __str__(self):
             string = ''
17
18
             for row in self.data:
                 string += '\n
19
                 for val in row:
20
                     string += '{:6.2f} '.format(val)
21
             return string
22
23
         def integer_string(self):
             string = ''
25
             for row in self.data:
26
                 string += '\n'
27
                 for val in row:
28
                     string += '{:3.0f} '.format(val)
29
             return string
30
31
         def precision_string(self):
32
             string = ''
33
             for row in self.data:
34
                 string += '\n'
35
                 for val in row:
36
                     string += '{:6.4f} '.format(val)
37
38
             return string
39
         def __add__(self, other):
40
             if len(self) != len(other) or len(self[0]) != len(other[0]):
41
                 raise ValueError('Incompatible matrix sizes for addition. Matrix A is <math>{\}x\{}, but matrix B is
42
                  \hookrightarrow {}x{}.'
                                   .format(len(self), len(self[0]), len(other), len(other[0])))
43
44
             return Matrix([[self[row][col] + other[row][col] for col in range(self.num_cols)]
45
                            for row in range(self.num_rows)])
46
47
         def __sub__(self, other):
48
             if len(self) != len(other) or len(self[0]) != len(other[0]):
49
                 raise ValueError('Incompatible matrix sizes for subtraction. Matrix A is {}x{}, but matrix B
50
                  \hookrightarrow is \{\}x\{\}.
                                   .format(len(self), len(self[0]), len(other), len(other[0])))
51
52
             return Matrix([[self[row][col] - other[row][col] for col in range(self.num_cols)]
53
                            for row in range(self.num_rows)])
55
         def __mul__(self, other):
56
             if type(other) == float or type(other) == int:
57
                 return self.scalar_multiply(other)
58
59
             if self.num_cols != other.num_rows:
60
                 raise ValueError('Incompatible matrix sizes for multiplication. Matrix A is {}x{}, but matrix
61
                  \hookrightarrow B is \{\}x\{\}.
                                   .format(self.num_rows, self.num_cols, other.num_rows, other.num_cols))
62
```

```
63
              # Inspired from https://en.wikipedia.org/wiki/Matrix_multiplication
 64
             product = Matrix.empty(self.num_rows, other.num_cols)
65
             for i in range(self.num_rows):
66
                  for j in range(other.num_cols):
67
                      row_sum = 0
68
69
                      for k in range(self.num_cols):
70
                          row_sum += self[i][k] * other[k][j]
                      product[i][j] = row_sum
71
72
             return product
73
         def scalar_multiply(self, scalar):
74
             return Matrix([[self[row][col] * scalar for col in range(self.num_cols)] for row in
 75

    range(self.num_rows)])

 76
77
         def __div__(self, other):
78
79
             Element-wise division.
80
             if self.num_rows != other.num_rows or self.num_cols != other.num_cols:
81
82
                  raise ValueError('Incompatible matrix sizes.')
             return Matrix([[self[row][col] / other[row][col] for col in range(self.num_cols)]
83
84
                             for row in range(self.num_rows)])
85
         def __neg__(self):
86
87
             return Matrix([[-self[row][col] for col in range(self.num_cols)] for row in range(self.num_rows)])
88
         def __deepcopy__(self, memo):
89
             return Matrix(copy.deepcopy(self.data))
90
91
         def __getitem__(self, item):
92
             return self.data[item]
93
94
95
         def __len__(self):
             return len(self.data)
96
97
         def item(self):
98
99
              :return: the single element contained by this matrix, if it is 1x1.
100
101
             if not (self.num_rows == 1 and self.num_cols == 1):
102
103
                  raise ValueError('Matrix is not 1x1')
             return self.data[0][0]
104
105
         def is_positive_definite(self):
106
107
             :return: True if the matrix if positive-definite, False otherwise.
108
109
             A = copy.deepcopy(self.data)
110
111
             for j in range(self.num_rows):
                  if A[j][j] <= 0:
112
                      return False
113
114
                  A[j][j] = math.sqrt(A[j][j])
                  for i in range(j + 1, self.num_rows):
115
116
                      A[i][j] = A[i][j] / A[j][j]
                      for k in range(j + 1, i + 1):
117
                          A[i][k] = A[i][k] - A[i][j] * A[k][j]
118
             return True
119
120
         def transpose(self):
121
122
              :return: the transpose of the current matrix
123
124
             return Matrix([[self.data[row][col] for row in range(self.num_rows)] for col in
125

    range(self.num_cols)])

126
127
         def mirror_horizontal(self):
128
              :return: the horizontal mirror of the current matrix
129
130
```

```
return Matrix([[self.data[self.num_rows - row - 1][col] for col in range(self.num_cols)]
131
132
                              for row in range(self.num_rows)])
133
         def empty_copy(self):
134
135
              :return: an empty matrix of the same size as the current matrix.
136
137
138
              return Matrix.empty(self.num_rows, self.num_cols)
139
         def infinity_norm(self):
140
              if self.num_cols > 1:
141
                  raise ValueError('Not a column vector.')
142
              return max([abs(x) for x in self.transpose()[0]])
144
         def two_norm(self):
145
146
              if self.num_cols > 1:
                  raise ValueError('Not a column vector.')
147
148
              return math.sqrt(sum([x ** 2 for x in self.transpose()[0]]))
149
         def save_to_csv(self, filename):
150
151
              Saves the current matrix to a CSV file.
152
153
              :param filename: the name of the CSV file
154
155
              with open(filename, "wb") as f:
156
                  writer = csv.writer(f)
157
                  for row in self.data:
158
                      writer.writerow(row)
159
160
         def save_to_latex(self, filename):
161
162
             Saves the current matrix to a latex-readable matrix.
163
164
              :param filename: the name of the CSV file
165
166
              with open(filename, "wb") as f:
167
                  for row in range(self.num_rows):
168
169
                      for col in range(self.num_cols):
170
                          f.write('{}'.format(self.data[row][col]))
                          if col < self.num_cols - 1:</pre>
171
172
                              f.write('& ')
                      if row < self.num_rows - 1:</pre>
173
                          f.write('\\\\n')
174
         @staticmethod
176
         def multiply(*matrices):
177
178
             Computes the product of the given matrices.
179
180
              :param matrices: the matrix objects
181
              :return: the product of the given matrices
182
183
             n = matrices[0].rows
184
185
              product = Matrix.identity(n)
              for matrix in matrices:
186
                  product = product * matrix
187
188
             return product
189
         Ostaticmethod
190
         def empty(num_rows, num_cols):
191
192
              Returns an empty matrix (filled with zeroes) with the specified number of columns and rows.
193
194
              :param num_rows: number of rows
195
              : param\ num\_cols:\ number\ of\ columns
196
197
              :return: the empty matrix
198
              return Matrix([[0 for _ in range(num_cols)] for _ in range(num_rows)])
199
```

200

```
201
         Ostaticmethod
202
         def identity(n):
203
             Returns the identity matrix of the given size.
204
205
             :param n: the size of the identity matrix (number of rows or columns)
206
207
             :return: the identity matrix of size n
208
209
             return Matrix.diagonal_single_value(1, n)
210
         @staticmethod
211
         def diagonal(values):
212
             Returns a diagonal matrix with the given values along the main diagonal.
214
215
             :param values: the values along the main diagonal
216
             :return: a diagonal matrix with the given values along the main diagonal
217
218
             n = len(values)
219
             return Matrix([[values[row] if row == col else 0 for col in range(n)] for row in range(n)])
220
221
         Ostaticmethod
222
223
         def diagonal_single_value(value, n):
224
             Returns a diagonal matrix of the given size with the given value along the diagonal.
225
226
             :param value: the value of each element on the main diagonal
227
             :param n: the size of the matrix
228
             :return: a diagonal matrix of the given size with the given value along the diagonal.
230
             return Matrix([[value if row == col else 0 for col in range(n)] for row in range(n)])
231
232
         @staticmethod
233
234
         def column_vector(values):
235
             Transforms a row vector into a column vector.
236
237
             :param values: the values, one for each row of the column vector
238
239
             :return: the column vector
240
             return Matrix([[value] for value in values])
241
242
         Ostaticmethod
243
         def csv_to_matrix(filename):
244
245
             Reads a CSV file to a matrix.
246
247
             :param filename: the name of the CSV file
248
             :return: a matrix containing the values in the CSV file
249
250
             with open(filename, 'r') as csv_file:
251
                 reader = csv.reader(csv file)
252
253
                 data = []
                 for row_number, row in enumerate(reader):
254
255
                      data.append([literal_eval(val) for val in row])
                 return Matrix(data)
                                            Listing 2: Question 1 (q1.py).
    from finite_element_triangles import Triangle, find_local_s_matrix, find_global_s_matrix
     from matrices import Matrix
 2
 3
 5
     def q1():
         print('\n=== Question 1 ===')
 6
         S1 = build_triangle_and_find_local_S(
 8
             [0, 0, 0.02],
             [0.02, 0, 0])
 9
         S1.save_to_latex('report/matrices/S1.txt')
```

```
print('S1: {}'.format(S1))
11
12
        S2 = build_triangle_and_find_local_S(
13
            [0.02, 0, 0.02],
14
             [0.02, 0.02, 0])
15
        S2.save_to_latex('report/matrices/S2.txt')
16
        print('S2: {}'.format(S2))
17
18
19
        C = Matrix([
20
            [1, 0, 0, 0],
             [0, 1, 0, 0],
21
            [0, 0, 1, 0],
22
            [0, 0, 0, 1],
23
            [1, 0, 0, 0],
24
25
            [0, 0, 1, 0]])
        C.save_to_latex('report/matrices/C.txt')
26
        print('C: {}'.format(C))
27
28
        S = find_global_s_matrix(S1, S2, C)
29
        S.save_to_latex('report/matrices/S.txt')
30
31
        S.save_to_csv('report/csv/S.txt')
        print('S: {}'.format(S))
32
33
34
    def build_triangle_and_find_local_S(x, y):
35
36
        triangle = Triangle(x, y)
        S = find_local_s_matrix(triangle)
37
        return S
38
39
40
    if __name__ == '__main__':
41
        q1()
42
                      Listing 3: Finite element triangles (finite_element_triangles.py).
    from __future__ import division
    from matrices import Matrix
3
    class Triangle:
6
        def __init__(self, x, y):
            self.x = x
9
            self.y = y
            self.area = (x[1] * y[2] - x[2] * y[1] - x[0] * y[2] + x[2] * y[0] + x[0] * y[1] - x[1] * y[0]) /
10
11
12
    def find_local_s_matrix(triangle):
13
14
        x = triangle.x
        y = triangle.y
15
        S = Matrix.empty(3, 3)
16
17
        for i in range(3):
18
19
            for j in range(3):
                 S[i][j] = ((y[(i + 1) \% 3] - y[(i + 2) \% 3]) * (y[(j + 1) \% 3] - y[(j + 2) \% 3])
20
                            + (x[(i + 1) \% 3] - x[(i + 2) \% 3]) * <math>(x[(j + 1) \% 3] - x[(j + 2) \% 3])) / (4 * 
21
                             22
23
        return S
24
25
26
    def find_global_s_matrix(S1, S2, C):
        S_dis = find_disjoint_s_matrix(S1, S2)
27
        S_dis.save_to_latex('report/matrices/S_dis.txt')
28
29
        print('S_dis: {}'.format(S_dis))
        return C.transpose() * S_dis * C
30
31
```

```
33
    def find_disjoint_s_matrix(S1, S2):
        n = len(S1)
34
        S_{dis} = Matrix.empty(2 * n, 2 * n)
35
        for row in range(n):
36
             for col in range(n):
37
                 S_dis[row][col] = S1[row][col]
38
                S_dis[row + n][col + n] = S2[row][col]
39
40
        return S_dis
                                           Listing 4: Question 2 (q2.py).
    from finite_element_capacitance import find_capacitance
    from matrices import Matrix
    from \ finite\_element\_mesh\_generator \ import \ generate\_simple\_2d\_mesh
3
    INNER_CONDUCTOR_POINTS = [28, 29, 30, 34]
    OUTER_CONDUCTOR_POINTS = [1, 2, 3, 4, 5, 6, 7, 13, 19, 25, 31]
    MESH\_SIZE = 6
8
10
    def q2():
11
        print('\n=== Question 2 ===')
12
        q2a()
13
14
        q2c()
15
16
17
    def q2a():
        generate_simple_2d_mesh(MESH_SIZE, INNER_CONDUCTOR_POINTS, OUTER_CONDUCTOR_POINTS)
18
19
20
    def q2c():
21
        print('\n=== Question 2(c) ===')
22
23
        S = Matrix.csv_to_matrix('report/csv/S.txt')
        voltage = 15
24
25
        capacitance = find_capacitance(S, voltage, MESH_SIZE)
        print('Capacitance per unit length: {} F/m'.format(capacitance))
26
27
    if __name__ == '__main__':
29
30
        q2()
                Listing 5: Finite element mesh generator (finite_element_mesh_generator.py).
    def generate_simple_2d_mesh(mesh_size, inner_conductor_points, outer_conductor_points):
1
2
        with open('simple2d/mesh.dat', 'w') as f:
3
             generate_node_positions(f, mesh_size)
             generate_triangle_coordinates(f, mesh_size)
4
             generate_initial_potentials(f, inner_conductor_points, outer_conductor_points)
5
6
    def generate_node_positions(f, mesh_size):
        for row in range(mesh_size):
9
10
             y = row * 0.02
             for col in range(mesh_size):
11
12
                 x = col * 0.02
13
                 node = row * mesh_size + (col + 1)
                 if node <= 34: # Inner conductor</pre>
14
                     f.write('{} {} {}\n'.format(node, x, y))
15
        f.write('\n')
16
17
18
    def generate_triangle_coordinates(f, mesh_size):
19
        # Left triangles (left halves of squares)
20
21
        for row in range(mesh_size - 1):
            for col in range(mesh_size - 1):
22
                node = row * mesh\_size + (col + 1)
23
                 if node < 28:
```

```
25
                    f.write('{} {} {} {} 0\n'.format(node, node + 1, node + mesh_size))
26
        # Right triangles (right halves of squares)
27
        for row in range(mesh_size - 1):
28
            for col in range(1, mesh_size):
29
                node = row * mesh_size + (col + 1)
30
                if node <= 28:
31
32
                    33
        f.write('\n')
34
35
36
    def generate_initial_potentials(f, inner_conductor_points, outer_conductor_points):
37
        for point in outer_conductor_points:
38
            f.write('{} {}\n'.format(point, 0))
39
        for point in inner_conductor_points:
40
            f.write('{} {}\n'.format(point, 15))
41
                   Listing 6: Finite element capacitance (finite_element_capacitance.py).
    from matrices import Matrix
1
    E_0 = 8.854187817620E-12
4
    def extract_mesh():
        with open('simple2d/result.dat') as f:
7
8
            mesh = \{\}
            for line_number, line in enumerate(f):
9
10
                if line_number >= 2:
11
                    vals = line.split()
                    node = int(float(vals[0]))
12
13
                    voltage = float(vals[3])
                    mesh[node] = voltage
14
        return mesh
15
16
17
    def compute_half_energy(S, mesh, mesh_size):
18
19
        U_con = Matrix.empty(4, 1)
        half_energy = 0
20
        for row in range(mesh_size - 1):
21
            for col in range(mesh_size - 1):
22
                node = row * mesh_size + (col + 1) # 1-based
23
                if node < 28:
24
                    U_con[0][0] = mesh[node + mesh_size]
25
                    U_{con[1][0] = mesh[node]
26
27
                    U_{con[2][0] = mesh[node + 1]
                    U_{con[3][0]} = mesh[node + mesh_size + 1]
28
                    {\tt half\_energy\_contribution} \ = \ {\tt U\_con.transpose()} \ * \ {\tt S} \ * \ {\tt U\_con}
29
                    half_energy += half_energy_contribution[0][0]
30
        return half_energy
31
32
33
    def find_capacitance(S, voltage, mesh_size):
34
        mesh = extract_mesh()
35
        half_energy = compute_half_energy(S, mesh, mesh_size)
36
        capacitance = (4 * E_0 * half_energy) / voltage ** 2
37
        return capacitance
                                          Listing 7: Question 3 (q3.py).
    from copy import deepcopy
    import matplotlib.pyplot as plt
3
    from matplotlib import rc
5
    from matplotlib.ticker import MaxNLocator
```

```
from choleski import choleski_solve
    from conjugate_gradient import conjugate_gradient_solve
    from finite_difference_mesh_generator import generate_finite_diff_mesh
10
11
    MESH_SIZE = 6
12
    NUM_FREE_NODES = 19
13
    rc('font', **{'family': 'serif', 'serif': ['Computer Modern']})
14
    rc('text', usetex=True)
15
16
    def q3():
17
        print('\n=== Question 3 ===')
18
        A, b = q3a()
19
        choleski_potential, cg_potential, residual_vectors = q3b(A, b)
20
        q3c(residual_vectors)
21
22
        q3d(choleski_potential, cg_potential)
23
24
25
    def q3a():
        print('\n=== Question 3(a) ===')
26
        A, b = generate_finite_diff_mesh(MESH_SIZE, NUM_FREE_NODES)
27
28
        print('A: {}'.format(A.integer_string()))
        print('b: {}'.format(b.integer_string()))
29
30
        print('A is positive definite: {}'.format(A.is_positive_definite()))
31
        A_{prime} = A.transpose() * A
        b_prime = A.transpose() * b
32
        print("A' is positive definite: {}".format(A_prime.is_positive_definite()))
33
        return A_prime, b_prime
34
35
36
    def q3b(A, b):
37
        print('\n=== Question 3(b) ===')
38
39
        A = copy = deepcopy(A)
        b_{copy} = deepcopy(b)
40
41
        x_choleski = choleski_solve(A_copy, b_copy)
        print('Choleski x: {}'.format(x_choleski))
42
        residual_vectors = []
43
        x_cg = conjugate_gradient_solve(A, b, residual_vectors)
44
        print('Conjugate gradient x: {}'.format(x_cg))
45
46
        node_6_4 = 7
47
        return x_choleski[node_6_4][0], x_cg[node_6_4][0], residual_vectors
48
49
    def q3c(residual_vectors):
50
        print('\n=== Question 3(c) ===')
51
        plot_residual_norms(residual_vectors, infinity_norm=False)
52
        plot_residual_norms(residual_vectors, infinity_norm=True)
53
54
55
56
    def q3d(choleski_potential, cg_potential):
57
        print('\n=== Question 3(d) ===')
        print('Choleski potential at (0.06, 0.04): {} V'.format(choleski_potential))
58
        print('Conjugate gradient potential at (0.06, 0.04): {} V'.format(cg_potential))
59
60
61
62
    def plot_residual_norms(residual_vectors, infinity_norm=False):
63
        f = plt.figure()
        ax = f.gca()
64
65
        ax.xaxis.set_major_locator(MaxNLocator(integer=True))
66
        x_range = [i for i in range(len(residual_vectors))]
        y_range = [v.infinity_norm() if infinity_norm else v.two_norm() for v in residual_vectors]
67
        plt.plot(x_range, y_range, 'o-{}'.format('CO' if infinity_norm else 'C1'),
68
                  label=''.format('Infinity norm' if infinity_norm else '2-norm'))
69
        plt.xlabel('Conjugate gradient iteration ($k$)')
70
        plt.ylabel('Infinity norm of residual vector $(\\\\textbf{r}\\\|_\\infty)$' if infinity_norm
71
                    else '2-norm of residual vector ((\| \text{r})_{2})')
72
73
        plt.grid(True)
        f.savefig('report/plots/q3c_{}.pdf'.format('infinity' if infinity_norm else '2'), bbox_inches='tight')
74
75
76
    if __name__ == '__main__':
77
```

78 q3()

11

12

if residual\_vectors is not None:
 residual\_vectors.append(r)

for \_ in range(n):

Listing 8: Finite difference mesh generator (finite\_difference\_mesh\_generator.py). from matrices import Matrix def generate\_finite\_diff\_mesh(mesh\_size, num\_free\_nodes): 4 A = Matrix.empty(num\_free\_nodes, num\_free\_nodes) b = Matrix.empty(num\_free\_nodes, 1) 6 for row in range(mesh\_size - 3): 7 for col in range(mesh\_size - 1): node = row \* (mesh\_size - 1) + col 9 A[node][node] = -410 11 if row != 0: 12 13  $A[node][node - mesh\_size + 1] = 1$ if 12 <= node <= 14: 14 b[node][0] = -1515 16 else:  $A[node][node + mesh\_size - 1] = 1$ 17 18 # Right Neumann boundary 19 if col == mesh\_size - 2: 20 21 A[node][node - 1] = 222 if col != 0: 23 A[node][node - 1] = 124 A[node][node + 1] = 125 26 27 # Special nodes A[15][10] = 128 A[15][15] = -429 30 A[15][16] = 1A[15][17] = 131 32 A[16][11] = 133 A[16][15] = 134 A[16][16] = -4A[16][18] = 136 b[16][0] = -1537 38 A[17][15] = 239 A[17][17] = -440 A[17][18] = 141 42 A[18][16] = 243 A[18][17] = 144 45 A[18][18] = -446 b[18][0] = -1547 48 return A, b Listing 9: Conjugate gradient (conjugate\_gradient.py). 1 from copy import deepcopy from matrices import Matrix 3 4  ${\tt def\ conjugate\_gradient\_solve(A,\ b,\ residual\_vectors=None):}$ 6 n = len(A)x = Matrix.empty(n, 1)8 r = b - A \* x9 10 p = deepcopy(r)

```
14
            denom = p.transpose() * A * p
15
            alpha = (p.transpose() * r) / denom
            x = x + p * alpha.item()
16
            r = b - A * x
17
            beta = - (p.transpose() * A * r) / denom
18
            p = r + p * beta.item()
19
            if residual_vectors is not None:
20
21
                residual_vectors.append(r)
22
        return x
```

## B Output Logs

Listing 10: Output of Question 1 program (q1.txt).

```
=== Question 1 ===
2
     0.50 -0.50 0.00
3
           1.00 -0.50
     -0.50
     0.00 -0.50
                  0.50
5
    S2:
6
     1.00 -0.50 -0.50
     -0.50
            0.50
                  0.00
8
9
    -0.50
            0.00
                   0.50
   C:
10
     1.00
            0.00
                   0.00
                          0.00
11
12
     0.00
            1.00
                   0.00
                          0.00
            0.00
                         0.00
      0.00
                  1.00
13
14
     0.00
            0.00
                   0.00
                         1.00
15
      1.00
            0.00
                   0.00
                         0.00
     0.00
            0.00
                  1.00
                         0.00
16
17
   S_dis:
18
      0.50
           -0.50
                  0.00
                         0.00
                                0.00
           1.00 -0.50
                         0.00
                                0.00
                                      0.00
     -0.50
19
20
     0.00 - 0.50
                  0.50
                         0.00
                               0.00
                                      0.00
      0.00
            0.00
                   0.00
                         1.00
                               -0.50
                                      -0.50
21
      0.00
           0.00
                   0.00 -0.50
                               0.50
                                      0.00
22
      0.00
           0.00
                 0.00 -0.50
                                0.00
                                      0.50
    S:
24
     1.00 -0.50
                  0.00 -0.50
25
     -0.50
           1.00 -0.50 0.00
26
     0.00 -0.50
                  1.00 -0.50
27
                        1.00
           0.00 -0.50
28
     -0.50
```

Listing 11: Output of Question 2 program (q2.txt).

```
1 === Question 2 ===
2
3 === Question 2(c) ===
4 Capacitance per unit length: 5.21374340427e-11 F/m
```

Listing 12: Output of Question 3 program (q3. txt).

```
=== Question 3 ===
2
   === Question 3(a) ===
3
   A:
4
               0
                   0
5
    -4
        1
            0
                      1
                          0
                              0
                                 0
                                     0
                                         0
                                            0
                                                0
                                                   0
                                                       0
                                                          0
                                                              0
                                                                  0
                                                                     0
        -4
            1
                0
                   0
                       0
                          1
                              0
                                 0
                                     0
                                         0
                                            0
                                                0
                                                   0
                                                       0
                                                           0
                                                              0
                                                                  0
                                                                     0
6
                          0
     0
           -4
               1
                   0
                      0
                              1
                                 0
                                     0
                                         0
                                            0
                                                0
                                                   0
                                                       0
                                                           0
                                                              0
                                                                  0
     0
        0
           1
               -4
                   1
                       0
                          0
                              0
                                 1
                                     0
                                         0
                                            0
                                                0
                                                   0
                                                       0
                                                           0
                                                              0
                                                                  0
                                                                     0
        0
            0
               2
                   -4
                       0
                          0
                              0
9
                                     1
                  0 -4
        0 0 0
                              0
                                 0
                                     0
                                            0
                                                0
                                                   0
                                                       0
                                                          0
                                                              0
                                                                  0
                                                                     0
                          1
                                         1
10
     1
     0
        1 0 0 0
                      1 -4
                             1
                                 0
                                     0
                                         0
                                            1
                                                0
                                                   0
                                                       0 0
                                                             0
                                                                  0
                      0
                          1
                                                1
12
```

```
0
     0
         0
            0
                    0
                        0
                            0
                                1 -4
                                           0
                                               0
                                                  0
                                                      1
0
                                                              0
                                                                  0
                                                                      0
                1
                                       1
                                                                          0
13
                                       -4
14
      0
         0
            0
                0
                    1
                        0
                            0
                                0
                                   2
                                           0
                                               0
                                                   0
                                                          1
                                                              0
                                                                  0
                                                                      0
                                                                          0
         0
                            0
                                0
                                    0
                                       0
                                          -4
                                               1
                                                                  0
                                                                      0
                        1
15
             0
                 0
                    0
                        0
                                0
                                       0
                                               -4
                                                       0
                                                           0
                                                                      0
                                                                          0
      0
         0
                            1
                                    0
                                           1
                                                   1
                                                              0
                                                                  1
16
17
      0
         0
             0
                 0
                    0
                         0
                            0
                                1
                                    0
                                        0
                                           0
                                               1
                                                  -4
                                                       1
                                                           0
                                                              0
                                                                  0
                                                                      0
                                                                          0
             0
                     0
                        0
                            0
                                                  1
      0
         0
                 0
                                0
                                    1
                                        0
                                           0
                                               0
                                                      -4
                                                          1
                                                              0
                                                                  0
                                                                      0
                                                                          0
18
                 0
      0
         0
             0
                    0
                         0
                            0
                                0
                                    0
                                        1
                                           0
                                               0
                                                   0
                                                      2
                                                          -4
                                                              0
                                                                  0
                                                                      0
                                                                          0
19
20
      0
         0
             0
                 0
                     0
                         0
                            0
                                0
                                    0
                                        0
                                            1
                                               0
                                                   0
                                                       0
                                                          0
                                                              -4
                                                                          0
            0
                 0
                    0
                        0
                            0
                                0
                                    0
                                       0
                                           0
                                                   0
                                                      0
                                                          0
      0
         0
                                               1
                                                              1
                                                                 -4
                                                                     0
^{21}
                                                                          1
                0
                        0
                            0
                                0
                                       0
                                           0
                                                   0
                                                      0 0 2
22
      0
         0
            0
                    0
                                   0
                                               0
                                                                 0 -4
                                                                         1
     0
         0
             0
                 0
                    0
                        0
                            0
                                0
                                    0
                                        0
                                           0
                                               0
                                                   0
                                                                         -4
23
    b:
24
25
      0
      0
26
27
      0
      0
28
      0
29
30
      0
31
      0
32
33
      0
      0
34
35
     0
36
     0
    -15
37
38
    -15
39
    -15
     0
40
41
    -15
    0
42
   -15
43
44
   A is positive definite: False
   A' is positive definite: True
45
46
   === Question 3(b) ===
47
   Choleski x:
48
49
     0.96
     1.86
50
     2.61
51
52
      3.04
      3.17
53
54
     1.97
      3.88
55
      5.53
56
57
      6.37
      6.61
58
      3.03
59
60
      6.18
     9.25
61
     10.29
62
     10.55
63
     3.96
64
65
      8.56
     4.25
66
     9.09
67
68
    Conjugate gradient x:
     0.96
69
     1.86
70
71
      2.61
      3.04
72
      3.17
73
      1.97
74
      3.88
75
76
      5.53
      6.37
77
      6.61
78
79
      3.03
      6.18
80
81
     9.25
     10.29
82
```

```
10.55
83
      3.96
      8.56
85
      4.25
86
87
      9.09
88
    === Question 3(c) ===
89
90
    === Question 3(d) ===
91
   Choleski potential at (0.06, 0.04): 5.52634126517 V
    Conjugate gradient potential at (0.06, 0.04): 5.52634127414 V
93
```

## C Simple2D Data Files

Listing 13: Input mesh for the SIMPLE2D program.

```
1 1 0.0 0.0
2 2 0.02 0.0
   3 0.04 0.0
   4 0.06 0.0
5 5 0.08 0.0
   6 0.1 0.0
   7 0.0 0.02
   8 0.02 0.02
   9 0.04 0.02
   10 0.06 0.02
10
11 11 0.08 0.02
12 12 0.1 0.02
   13 0.0 0.04
14 14 0.02 0.04
16
   16 0.06 0.04
   17 0.08 0.04
17
   18 0.1 0.04
   19 0.0 0.06
19
   20 0.02 0.06
20
21 21 0.04 0.06
   22 0.06 0.06
22
   23 0.08 0.06
23
   24 0.1 0.06
   25 0.0 0.08
25
   26 0.02 0.08
26
27 27 0.04 0.08
   28 0.06 0.08
28
   29 0.08 0.08
   30 0.1 0.08
30
31
   31 0.0 0.1
   32 0.02 0.1
   33 0.04 0.1
33
   34 0.06 0.1
34
35
   1 2 7 0
36
   2 3 8 0
   3 4 9 0
38
   4 5 10 0
39
   5 6 11 0
   7 8 13 0
41
   8 9 14 0
42
   9 10 15 0
43
   10 11 16 0
44
   11 12 17 0
45
   13 14 19 0
46
47 14 15 20 0
   15 16 21 0
   16 17 22 0
49
50 17 18 23 0
```

19 20 25 0

```
20 21 26 0
52
    21 22 27 0
53
    22 23 28 0
54
    23 24 29 0
55
    25 26 31 0
    26 27 32 0
57
    27 28 33 0
59
    2 7 8 0
    3 8 9 0
60
    4 9 10 0
    5 10 11 0
62
    6 11 12 0
63
    8 13 14 0
    9 14 15 0
65
    10 15 16 0
66
    11 16 17 0
67
    12 17 18 0
68
69
    14 19 20 0
    15 20 21 0
70
    16 21 22 0
71
    17 22 23 0
    18 23 24 0
73
74
    20 25 26 0
    21 26 27 0
75
    22 27 28 0
76
    23 28 29 0
78
    24 29 30 0
    26 31 32 0
79
    27 32 33 0
    28 33 34 0
81
82
    1 0
83
    2 0
84
    3 0
85
86
    5 0
87
    6 0
    7 0
89
    13 0
90
91
    19 0
    25 0
92
93
    31 0
    28 15
94
    29 15
95
    30 15
    34 15
```

97

Listing 14: Resulting potentials generated by the SIMPLE2D program.

```
ans =
2
        1.0000
                        0
                                  0
3
        2.0000
                   0.0200
                                   0
                                             0
                   0.0400
                                  0
                                             0
        3.0000
5
        4.0000
                   0.0600
                                  0
                                             0
        5.0000
                   0.0800
                                  0
                                             0
        6.0000
                   0.1000
                             0.0200
9
        7.0000
                        0
        8.0000
                   0.0200
                             0.0200
                                       0.9571
10
        9.0000
                   0.0400
                             0.0200
                                       1.8616
11
12
       10.0000
                   0.0600
                             0.0200
                                        2.6060
       11.0000
                   0.0800
                             0.0200
                                        3.0360
13
       12.0000
                   0.1000
                             0.0200
                                        3.1714
       13.0000
                             0.0400
15
                   0.0200
       14.0000
                                        1.9667
                             0.0400
16
                   0.0400
                                       3.8834
17
       15.0000
                             0.0400
       16.0000
                   0.0600
                             0.0400
                                        5.5263
18
       17.0000
                                        6.3668
                   0.0800
                             0.0400
19
       18.0000
                   0.1000
                             0.0400
                                        6.6135
```

21	19.0000	0	0.0600	0
22	20.0000	0.0200	0.0600	3.0262
23	21.0000	0.0400	0.0600	6.1791
24	22.0000	0.0600	0.0600	9.2492
25	23.0000	0.0800	0.0600	10.2912
26	24.0000	0.1000	0.0600	10.5490
27	25.0000	0	0.0800	0
28	26.0000	0.0200	0.0800	3.9590
29	27.0000	0.0400	0.0800	8.5575
30	28.0000	0.0600	0.0800	15.0000
31	29.0000	0.0800	0.0800	15.0000
32	30.0000	0.1000	0.0800	15.0000
33	31.0000	0	0.1000	0
34	32.0000	0.0200	0.1000	4.2525
35	33.0000	0.0400	0.1000	9.0919
36	34.0000	0.0600	0.1000	15.0000