ECSE 543 Assignment 2

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Introduction

1 Finite Element Triangles

The equation for the α 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, ϕ_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).

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

The infinity norm of the residual vector versus conjugate gradient iterations can be seen in Figure 1. The 2-norm versus iterations can be seen in Figure 2. Both norms converge to 0 after n=19 iterations, as expected.

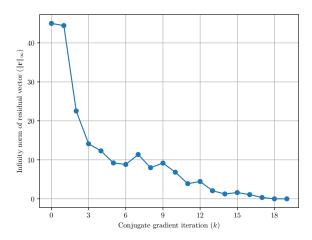


Figure 1: Value of the infinity norm $\|\mathbf{r}\|_{\infty}$ of the residual vector versus iterations of the conjugate gradient algorithm.

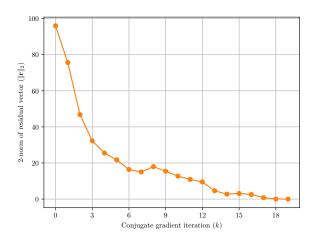


Figure 2: Value of the 2-norm $\|\mathbf{r}\|_2$ of the residual vector versus iterations of the conjugate gradient algorithm.

3.d Potential Comparison

A comparison of the potential at (0.06, 0.04) to three decimal places for various numerical methods can be seen in Table 1. It can be seen that the

potential found with all the methods is the same to three decimal places.

Table 1: Comparison of potential at (0.06, 0.04) to three decimal places for various numerical methods.

Method	Potential (V)	
Choleski	5.526	
Conjugate gradient	5.526	
Finite element	5.526	
Finite difference (SOR)	5.526	

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 __add__(self, other):
32
                                        if len(self) != len(other) or len(self[0]) != len(other[0]):
33
                                                      \textbf{raise ValueError('Incompatible matrix sizes for addition. Matrix A is $\{\}x\{\}$, but matrix B is $\{\}x\{\}$, but matrix B
34
                                                        \hookrightarrow {}x{}.'
                                                                                                             .format(len(self), len(self[0]), len(other), len(other[0])))
35
36
                                        return Matrix([[self[row][col] + other[row][col] for col in range(self.num_cols)]
37
38
                                                                                        for row in range(self.num_rows)])
                           def __sub__(self, other):
40
                                         if len(self) != len(other) or len(self[0]) != len(other[0]):
 41
                                                    raise ValueError('Incompatible matrix sizes for subtraction. Matrix A is {}x{}, but matrix B
 42
                                                        \hookrightarrow is \{\}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
 46
                                                                                        for row in range(self.num_rows)])
47
48
                           def __mul__(self, other):
                                         if type(other) == float or type(other) == int:
 49
                                                    return self.scalar_multiply(other)
50
51
                                         if self.num_cols != other.num_rows:
52
                                                     \textbf{raise ValueError('Incompatible matrix sizes for multiplication. Matrix A is \{}x\{\}, \ but \ matrix \ and \ an arrive in the property of th
53
                                                                                                            .format(self.num_rows, self.num_cols, other.num_rows, other.num_cols))
54
55
                                        # Inspired from https://en.wikipedia.org/wiki/Matrix_multiplication
56
                                        product = Matrix.empty(self.num_rows, other.num_cols)
57
58
                                        for i in range(self.num_rows):
                                                     for j in range(other.num_cols):
59
60
                                                                 row_sum = 0
                                                                  for k in range(self.num_cols):
                                                                              row_sum += self[i][k] * other[k][j]
62
```

```
product[i][j] = row_sum
63
             return product
64
65
         def scalar_multiply(self, scalar):
66
              return Matrix([[self[row][col] * scalar for col in range(self.num_cols)] for row in
67

    range(self.num_rows)])

 68
69
         def __div__(self, other):
70
71
             Element-wise division.
72
             if self.num_rows != other.num_rows or self.num_cols != other.num_cols:
73
                 raise ValueError('Incompatible matrix sizes.')
             return Matrix([[self[row][col] / other[row][col] for col in range(self.num_cols)]
75
76
                             for row in range(self.num_rows)])
77
         def __neg__(self):
78
79
              return Matrix([[-self[row][col] for col in range(self.num_cols)] for row in range(self.num_rows)])
80
         def __deepcopy__(self, memo):
81
82
             return Matrix(copy.deepcopy(self.data))
83
84
         def __getitem__(self, item):
             return self.data[item]
85
86
87
         def __len__(self):
             return len(self.data)
88
89
         def item(self):
90
91
             :return: the single element contained by this matrix, if it is 1x1.
92
93
             if not (self.num_rows == 1 and self.num_cols == 1):
94
                 raise ValueError('Matrix is not 1x1')
95
             return self.data[0][0]
96
97
         def is_positive_definite(self):
98
99
              : return: \ \textit{True if the matrix if positive-definite, False otherwise}.
100
101
             A = copy.deepcopy(self.data)
102
103
             for j in range(self.num_rows):
                  if A[j][j] <= 0:
104
                      return False
105
                  A[j][j] = math.sqrt(A[j][j])
                  for i in range(j + 1, self.num_rows):
107
                      A[i][j] = A[i][j] / A[j][j]
108
                      for k in range(j + 1, i + 1):
109
                          A[i][k] = A[i][k] - A[i][j] * A[k][j]
110
111
             return True
112
         def transpose(self):
113
114
              :return: the transpose of the current matrix
115
116
             return Matrix([[self.data[row][col] for row in range(self.num_rows)] for col in
117

    range(self.num_cols)])

118
119
         def mirror_horizontal(self):
120
              :return: the horizontal mirror of the current matrix
121
122
             return Matrix([[self.data[self.num_rows - row - 1][col] for col in range(self.num_cols)]
123
                             for row in range(self.num_rows)])
124
125
126
         def empty_copy(self):
127
              :return: an empty matrix of the same size as the current matrix.
128
129
             return Matrix.empty(self.num_rows, self.num_cols)
130
```

```
131
132
          def infinity_norm(self):
              if self.num_cols > 1:
133
                  raise ValueError('Not a column vector.')
134
              return max([abs(x) for x in self.transpose()[0]])
135
136
137
         def two_norm(self):
138
              if self.num_cols > 1:
                 raise ValueError('Not a column vector.')
139
140
              return math.sqrt(sum([x ** 2 for x in self.transpose()[0]]))
141
          def save_to_csv(self, filename):
142
              Saves the current matrix to a CSV file.
144
145
              :param filename: the name of the CSV file
146
147
              with open(filename, "wb") as f:
148
                  writer = csv.writer(f)
149
                  for row in self.data:
150
151
                      writer.writerow(row)
152
153
          def save_to_latex(self, filename):
154
              Saves the current matrix to a latex-readable matrix.
155
156
              :param filename: the name of the CSV file
157
158
              with open(filename, "wb") as f:
                  for row in range(self.num_rows):
160
161
                      for col in range(self.num_cols):
                           f.write('{}'.format(self.data[row][col]))
162
                           if col < self.num_cols - 1:</pre>
163
164
                               f.write('&')
                      if row < self.num_rows - 1:</pre>
165
                          f.write('\\\\n')
166
167
          @staticmethod
168
         def multiply(*matrices):
169
170
              Computes the product of the given matrices.
171
172
              :param matrices: the matrix objects
173
              :return: the product of the given matrices
174
175
              n = matrices[0].rows
176
              product = Matrix.identity(n)
177
              for matrix in matrices:
178
                  product = product * matrix
179
180
              return product
181
          Ostaticmethod
182
183
          def empty(num_rows, num_cols):
184
              Returns an empty matrix (filled with zeroes) with the specified number of columns and rows.
185
186
              :param num_rows: number of rows
187
188
              :param num_cols: number of columns
              :return: the empty matrix
189
190
              return Matrix([[0 for _ in range(num_cols)] for _ in range(num_rows)])
191
192
          Ostaticmethod
193
          def identity(n):
194
195
              Returns the identity matrix of the given size.
196
197
              :param n: the size of the identity matrix (number of rows or columns)
198
              : return: \ the \ identity \ matrix \ of \ size \ n
199
200
```

```
201
             return Matrix.diagonal_single_value(1, n)
202
         @staticmethod
203
         def diagonal(values):
204
205
             Returns a diagonal matrix with the given values along the main diagonal.
206
207
208
             :param values: the values along the main diagonal
             :return: a diagonal matrix with the given values along the main diagonal
209
210
             n = len(values)
211
             return Matrix([[values[row] if row == col else 0 for col in range(n)] for row in range(n)])
212
213
         Ostaticmethod
214
215
         def diagonal_single_value(value, n):
216
             Returns a diagonal matrix of the given size with the given value along the diagonal.
217
218
             :param value: the value of each element on the main diagonal
219
             :param n: the size of the matrix
220
221
             :return: a diagonal matrix of the given size with the given value along the diagonal.
222
223
             return Matrix([[value if row == col else 0 for col in range(n)] for row in range(n)])
224
         @staticmethod
225
226
         def column_vector(values):
227
             Transforms a row vector into a column vector.
228
             :param values: the values, one for each row of the column vector
230
231
             :return: the column vector
232
             return Matrix([[value] for value in values])
233
234
         @staticmethod
235
         def csv_to_matrix(filename):
236
237
             Reads a CSV file to a matrix.
238
239
240
             :param filename: the name of the CSV file
             :return: a matrix containing the values in the CSV file
241
242
             with open(filename, 'r') as csv_file:
243
                 reader = csv.reader(csv_file)
244
                 data = []
                 for row_number, row in enumerate(reader):
246
                     data.append([literal_eval(val) for val in row])
247
                 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
 3
     def q1():
 5
         print('\n=== Question 1 ===')
 6
 7
         S1 = build_triangle_and_find_local_S(
             [0, 0, 0.02],
             [0.02, 0, 0])
 9
         S1.save_to_latex('report/matrices/S1.txt')
 10
         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
```

```
C = Matrix([
19
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
                             [0, 0, 1, 0]])
25
26
                   C.save_to_latex('report/matrices/C.txt')
                   print('C: {}'.format(C))
27
28
                   S = find_global_s_matrix(S1, S2, C)
29
                   S.save_to_latex('report/matrices/S.txt')
30
                   S.save_to_csv('report/csv/S.txt')
31
                   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
41
          if __name__ == '__main__':
                   q1()
42
                                                   Listing 3: Finite element triangles (finite_element_triangles.py).
          from __future__ import division
         from matrices import Matrix
 3
 5
          class Triangle:
  6
                    """Represents a finite-difference triangle."""
                   def __init__(self, x, y):
 8
 9
                             self.x = x
                            self.y = y
10
                            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]) /
11
12
13
14
          def find_local_s_matrix(triangle):
15
16
                   Finds the local S matrix for a finite-difference triangle.
17
                    :param triangle: the finite-difference triangle
18
                    : return \colon \ the \ local \ S \ matrix
19
20
                   x = triangle.x
21
22
                   y = triangle.y
                   S = Matrix.empty(3, 3)
23
24
                   for i in range(3):
                            for j in range(3):
26
27
                                      S[i][j] = ((y[(i + 1) \% 3] - y[(i + 2) \% 3]) * (y[(j + 1) \% 3] - y[(j + 2) \% 3])
                                                               + (x[(i + 1) \% 3] - x[(i + 2) \% 3]) * <math>(x[(j + 1) \% 3] - x[(j + 2) \% 3])) / (4 * (x[(i + 1) \% 3]) + (x[(i + 1) \% 3])) / (4 * (x[(i + 1) \% 3])) / (4 * (x[(i + 1) \% 3])) / (4 * (x[(i + 1) \% 3]))) / (4 * (x[(i + 1) \% 3])) / (4 * (x[(i + 1) \% 3]))) / (4 * (x[(i + 1) \% 3])))) / (4 * (x[(i + 1) \% 3])))))
28
                                                                 \hookrightarrow triangle.area)
29
                   return S
30
31
32
          def find_global_s_matrix(S1, S2, C):
33
34
                   Finds the global S matrix given by two local S matrices and the the connectivity matrix.
35
36
37
                    :param S1: the first local S matrix
                    :param S2: the second local S matrix
38
                    :param C: the connectivity matrix
39
                   :return: the global S matrix
```

```
41
42
        S_dis = find_disjoint_s_matrix(S1, S2)
        S_dis.save_to_latex('report/matrices/S_dis.txt')
43
        print('S_dis: {}'.format(S_dis))
44
        return C.transpose() * S_dis * C
45
46
47
48
    def find_disjoint_s_matrix(S1, S2):
49
50
        Finds the disjoint S matrix given by the two provided local S matrices.
51
         :param S1: the first local S matrix
52
        :param S2: the second local S matrix
53
         :return: the disjoint S matrix
54
55
        n = len(S1)
56
        S_{dis} = Matrix.empty(2 * n, 2 * n)
57
58
        for row in range(n):
            for col in range(n):
59
                S_dis[row][col] = S1[row][col]
60
61
                S_dis[row + n][col + n] = S2[row][col]
        return S_dis
62
                                           Listing 4: Question 2 (q2.py).
    from finite_element_capacitance import find_capacitance
    from matrices import Matrix
2
    from finite_element_mesh_generator import generate_simple_2d_mesh
    INNER_CONDUCTOR_POINTS = [28, 29, 30, 34]
5
    OUTER_CONDUCTOR_POINTS = [1, 2, 3, 4, 5, 6, 7, 13, 19, 25, 31]
6
    MESH_SIZE = 6
9
10
11
    def q2():
        print('\n=== Question 2 ===')
12
13
        q2a()
         q2c()
15
16
17
        generate_simple_2d_mesh(MESH_SIZE, INNER_CONDUCTOR_POINTS, OUTER_CONDUCTOR_POINTS)
18
19
20
    def q2c():
21
22
        print('\n=== Question 2(c) ===')
        S = Matrix.csv_to_matrix('report/csv/S.txt')
23
24
        voltage = 15
         capacitance = find_capacitance(S, voltage, MESH_SIZE)
25
        print('Capacitance per unit length: {} F/m'.format(capacitance))
26
27
28
    if __name__ == '__main__':
29
        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
         Generates the input mesh needed for the SIMPLE2D program.
3
4
5
         :param mesh_size: the mesh size
         : param\ inner\_conductor\_points \colon\ the\ inner\ conductor\ points
6
         : param\ outer\_conductor\_points:\ the\ outer\ conductor\ points
8
        with open('simple2d/mesh.dat', 'w') as f:
9
             generate_node_positions(f, mesh_size)
```

```
11
             generate_triangle_coordinates(f, mesh_size)
12
             generate_initial_potentials(f, inner_conductor_points, outer_conductor_points)
13
14
    def generate_node_positions(f, mesh_size):
15
16
         {\it Generates the node positions for the SIMPLE2D program.}
17
18
19
         :param f: the mesh file
20
         :param\ \textit{mesh\_size}\colon\ the\ \textit{mesh}\ \textit{size}
21
         for row in range(mesh_size):
22
             y = row * 0.02
23
             for col in range(mesh_size):
24
25
                 x = col * 0.02
                 node = row * mesh_size + (col + 1)
26
                  if node <= 34: # Inner conductor</pre>
27
28
                      f.write('{} {} {}) /n'.format(node, x, y))
         f.write('\n')
29
30
31
    def generate_triangle_coordinates(f, mesh_size):
32
33
         # Left triangles (left halves of squares)
34
         Generates the triangle coordinates for the SIMPLE2D program.
35
36
         :param f: the mesh file
37
         :param\ \textit{mesh\_size}\colon\ \textit{the mesh\ size}
38
39
         for row in range(mesh_size - 1):
40
41
             for col in range(mesh_size - 1):
                 node = row * mesh_size + (col + 1)
42
                  if node < 28:
43
44
                      f.write('{} {} {} {} 0\n'.format(node, node + 1, node + mesh_size))
45
         # Right triangles (right halves of squares)
46
47
         for row in range(mesh_size - 1):
             for col in range(1, mesh_size):
48
49
                 node = row * mesh_size + (col + 1)
50
                  if node <= 28:
                      f.write('{} {} {} {} 0\n'.format(node, node + mesh_size - 1, node + mesh_size))
51
52
         f.write('\n')
53
54
55
    def generate_initial_potentials(f, inner_conductor_points, outer_conductor_points):
56
57
         Generates the initial potentials for the SIMPLE2D program.
58
59
60
         :param f: the mesh file
         :param inner_conductor_points: the inner conductor points
61
         : param\ outer\_conductor\_points:\ the\ outer\ conductor\ points
62
63
         for point in outer_conductor_points:
64
65
             f.write('{} {}\n'.format(point, 0))
         for point in inner_conductor_points:
66
             f.write('{} {}\n'.format(point, 15))
67
                     Listing 6: Finite element capacitance (finite_element_capacitance.py).
    from matrices import Matrix
    E_0 = 8.854187817620E-12
    def extract_mesh():
7
         {\it Extracts mesh information from the SIMPLE2D file.}
```

```
10
         :return: the extracted mesh
11
         with open('simple2d/result.dat') as f:
12
             mesh = {}
13
             for line_number, line in enumerate(f):
14
                 if line_number >= 2:
15
                      vals = line.split()
16
17
                      node = int(float(vals[0]))
                      voltage = float(vals[3])
18
19
                      mesh[node] = voltage
         return mesh
20
21
22
    def compute_half_energy(S, mesh, mesh_size):
23
24
         Computes the half-energy needed to compute the capacitance of the mesh.
25
26
27
         :param S: the S matrix
         :param mesh: the mesh
28
         :param\ \textit{mesh\_size}\colon\ \textit{the mesh\ size}
29
30
         :return: the half-energy
31
32
         U_con = Matrix.empty(4, 1)
33
         half_energy = 0
         for row in range(mesh_size - 1):
34
35
             for col in range(mesh_size - 1):
                 node = row * mesh_size + (col + 1) # 1-based
36
                 if node < 28:
37
                      U_con[0][0] = mesh[node + mesh_size]
38
                      U_con[1][0] = mesh[node]
U_con[2][0] = mesh[node + 1]
39
40
                      U_con[3][0] = mesh[node + mesh_size + 1]
41
                      {\tt half\_energy\_contribution} \ = \ {\tt U\_con.transpose()} \ * \ {\tt S} \ * \ {\tt U\_con}
42
43
                      half_energy += half_energy_contribution[0][0]
         return half_energy
44
45
46
    def find_capacitance(S, voltage, mesh_size):
47
48
49
         Finds the capacitance per unit length of the mesh.
50
51
         :param S: the S matrix
         :param voltage: the voltage difference
52
         :param mesh_size: the mesh size
53
         :return: the capacitance per unit length
55
         mesh = extract_mesh()
56
         half_energy = compute_half_energy(S, mesh, mesh_size)
57
         capacitance = (4 * E_0 * half_energy) / voltage ** 2
58
59
         return capacitance
                                             Listing 7: Question 3 (q3.py).
    from copy import deepcopy
1
    import matplotlib.pyplot as plt
4
    from matplotlib import rc
    from matplotlib.ticker import MaxNLocator
    from choleski import choleski_solve
    from conjugate_gradient import conjugate_gradient_solve
10
    from finite_difference_mesh_generator import generate_finite_diff_mesh
11
    MESH_SIZE = 6
12
    NUM_FREE_NODES = 19
    rc('font', **{'family': 'serif', 'serif': ['Computer Modern']})
14
    rc('text', usetex=True)
15
```

```
17
        def q3():
               print('\n=== Question 3 ===')
18
               A, b = q3a()
19
               choleski_potential, cg_potential, residual_vectors = q3b(A, b)
20
               q3c(residual_vectors)
21
               q3d(choleski_potential, cg_potential)
22
23
24
25
        def q3a():
               print('\n=== Question 3(a) ===')
26
               A, b = generate_finite_diff_mesh(MESH_SIZE, NUM_FREE_NODES)
27
               print('A: {}'.format(A.integer_string()))
28
               print('b: {}'.format(b.integer_string()))
               print('A is positive definite: {}'.format(A.is_positive_definite()))
30
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
37
        def q3b(A, b):
               print('\n=== Question 3(b) ===')
38
39
               A_{copy} = deepcopy(A)
40
               b_copy = deepcopy(b)
               x_choleski = choleski_solve(A_copy, b_copy)
41
               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
               \tt return \ x\_choleski[node\_6\_4][0], \ x\_cg[node\_6\_4][0], \ residual\_vectors
47
48
49
50
        def q3c(residual_vectors):
               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):
               print('\n=== Question 3(d) ===')
57
58
               print('Choleski potential at (0.06, 0.04): {} V'.format(choleski_potential))
               print('Conjugate gradient potential at (0.06, 0.04): {} V'.format(cg_potential))
59
60
61
        def plot_residual_norms(residual_vectors, infinity_norm=False):
62
               f = plt.figure()
63
               ax = f.gca()
64
               ax.xaxis.set_major_locator(MaxNLocator(integer=True))
65
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
69
                                label=''.format('Infinity norm' if infinity_norm else '2-norm'))
               plt.xlabel('Conjugate gradient iteration ($k$)')
70
               plt.ylabel('Infinity norm of residual vector $((\|\textbf{r}\\|_\))$' if infinity_norm of residual vector $((\|\textbf{r},\textbf{r},\textbf{r},\textbf{r})$' if infinity_norm of residual vector $((\|\textbf{r},\textbf{r},\textbf{r},\textbf{r},\textbf{r},\textbf{r},\textbf{r})$' if infinity_norm of residual vector $((\|\textbf{r},\textbf{r},\textbf{r},\textbf{r},\textbf{r})$' if infinity_norm of residual vector $((\|\textbf{r},\textbf{r},\textbf{r},\textbf{r})$' if infinity_norm of residual vector $((\|\textbf{r},\textbf{r},\textbf{r})$')$' if infinity_norm of residual vector $((\|\textbf{r},\textbf{r})$')$' if infinity_norm of residual vector $(\|\textbf{r},\textbf{r})$' if infinity_norm of residual vector $(\|\textbf{r})$' if infinity_no
71
                                   else '2-norm of residual vector ((\| \text{r})_2)')
72
               plt.grid(True)
73
               f.savefig('report/plots/q3c_{}.pdf'.format('infinity' if infinity_norm else '2'), bbox_inches='tight')
74
75
76
        if __name__ == '__main__':
77
               q3()
78
                       Listing 8: Finite difference mesh generator (finite_difference_mesh_generator.py).
        from matrices import Matrix
 2
        def generate_finite_diff_mesh(mesh_size, num_free_nodes):
```

```
5
6
        Generates a finite-difference mesh with the given size and number of free nodes.
7
        :param mesh_size: the mesh size
8
9
         :param num_free_nodes: the number of free nodes
         :return: the A and b matrices defining the mesh equation (Ax = b)
10
11
12
        A = Matrix.empty(num_free_nodes, num_free_nodes)
13
        b = Matrix.empty(num_free_nodes, 1)
14
        for row in range(mesh_size - 3):
            for col in range(mesh_size - 1):
15
                node = row * (mesh_size - 1) + col
16
                A[node][node] = -4
17
18
                if row != 0:
19
                     A[node][node - mesh\_size + 1] = 1
20
                 if 12 <= node <= 14:
21
22
                     b[node][0] = -15
23
                     A[node][node + mesh\_size - 1] = 1
24
25
                 # Right Neumann boundary
26
27
                if col == mesh_size - 2:
                     A[node][node - 1] = 2
28
                 else:
29
                     if col != 0:
30
                         A[node][node - 1] = 1
31
                     A[node][node + 1] = 1
32
33
        # Special nodes
34
        A[15][10] = 1
35
        A[15][15] = -4
36
        A[15][16] = 1

A[15][17] = 1
37
38
39
        A[16][11] = 1
40
41
        A[16][15] = 1
        A[16][16] = -4
42
        A[16][18] = 1
43
44
        b[16][0] = -15
45
46
        A[17][15] = 2
        A[17][17] = -4
47
        A[17][18] = 1
48
        A[18][16] = 2
50
        A[18][17] = 1
51
        A[18][18] = -4
52
        b[18][0] = -15
53
54
        return A, b
55
                             Listing 9: Conjugate gradient (conjugate_gradient.py).
    from copy import deepcopy
3
    from matrices import Matrix
5
    def conjugate_gradient_solve(A, b, residual_vectors=None):
6
        Solves the Ax = b matrix equation given by the given A and b matrices
8
9
10
         :param A: the A matrix
         :param b: the b matrix
11
        :param residual_vectors: the list to store the residual vectors in
12
13
         :return: the solved x vector
14
        n = len(A)
```

```
x = Matrix.empty(n, 1)
16
17
        r = b - A * x
        p = deepcopy(r)
18
        if residual_vectors is not None:
19
            residual_vectors.append(r)
20
        for _ in range(n):
21
            \tt denom = p.transpose() * A * p
22
23
            alpha = (p.transpose() * r) / denom
            x = x + p * alpha.item()
24
            r = b - A * x
            beta = - (p.transpose() * A * r) / denom
26
            p = r + p * beta.item()
27
             if residual_vectors is not None:
                residual_vectors.append(r)
29
30
        return x
```

B Output Logs

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

```
=== Question 1 ===
   S1:
2
     0.50 -0.50 0.00
     -0.50 1.00 -0.50
     0.00 -0.50 0.50
   S2:
    1.00 -0.50 -0.50
    -0.50
           0.50
                  0.00
9
    -0.50
            0.00
                  0.50
10
11
     1.00
            0.00
                  0.00
                         0.00
12
     0.00
            1.00
                  0.00
            0.00
                  1.00
                         0.00
     0.00
13
14
     0.00
            0.00
                  0.00
                         1.00
     1.00
            0.00
                  0.00
                         0.00
15
     0.00
            0.00
16
                  1.00
                        0.00
   S_dis:
     0.50 -0.50
                  0.00
                         0.00
                               0.00
                                      0.00
18
           1.00
19
     -0.50
                 -0.50
                         0.00
                               0.00
                                      0.00
     0.00 -0.50
                  0.50
                        0.00
                               0.00
20
     0.00
            0.00
                  0.00
                        1.00
                              -0.50 -0.50
21
                  0.00 -0.50
22
     0.00
           0.00
                               0.50
                                     0.00
                 0.00 -0.50
     0.00
           0.00
                               0.00
                                     0.50
23
   S:
24
25
     1.00 -0.50
                  0.00 -0.50
     -0.50 1.00 -0.50 0.00
26
27
     0.00 -0.50 1.00 -0.50
           0.00 -0.50
     -0.50
                        1.00
```

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).

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

```
0
         1 -4
                    0
                        0
                                   0
                                       0
                                           0
                                               0
                                                  0
                                                      0
                                                          0
                                                              0
                                                                 0
     0
                1
                                1
                                                                      0
                                                                         0
7
8
      0
         0
            1
                -4
                    1
                        0
                            0
                                0
                                   1
                                       0
                                           0
                                               0
                                                  0
                                                      0
                                                          0
                                                              0
                                                                 0
                                                                      0
                                                                         0
                2
                   -4
                                           0
                                                                      0
                                       1
9
                0
                    0
                       -4
                                0
                                               0
                                                   0
                                                       0
                                                          0
                                                              0
                                                                      0
                                                                         0
      1
         0
             0
                            1
                                    0
                                       0
                                           1
                                                                  0
10
11
      0
         1
             0
                 0
                    0
                        1
                           -4
                                1
                                    0
                                       0
                                           0
                                               1
                                                   0
                                                       0
                                                          0
                                                              0
                                                                  0
                                                                      0
                                                                          0
      0
         0
            1
                 0
                    0
                        0
                            1
                               -4
                                   1
                                        0
                                           0
                                               0
                                                   1
                                                       0
                                                          0
                                                              0
                                                                  0
                                                                      0
                                                                          0
12
            0
                1
                    0
                        0
                            0
                                   -4
                                           0
                                               0
                                                   0
                                                      1
                                                          0
                                                              0
                                                                  0
                                                                      0
                                                                         0
13
      0
         0
                               1
                                       1
14
      0
         0
             0
                 0
                     1
                        0
                            0
                                0
                                   2
                                       -4
                                           0
                                               0
                                                   0
                                                       0
                                                          1
                                                              0
                                                                  0
                                                                      0
                                                                          0
            0
                0
                    0
                            0
                                0
                                       0
                                           -4
                                                       0
      0
         0
                        1
                                   0
                                               1
                                                   0
                                                          0
                                                                  0
                                                                      0
                                                                         0
15
                                                              1
16
      0
         0
            0
                0
                    0
                        0
                            1
                                0
                                   0
                                       0
                                           1
                                              -4
                                                  1
                                                       0
                                                          0
                                                              0
                                                                 1
                                                                      0
                                                                         0
      0
         0
             0
                 0
                     0
                        0
                            0
                                1
                                    0
                                       0
                                           0
                                               1
                                                  -4
                                                      1
                                                          0
                                                              0
                                                                  0
                                                                      0
                                                                          0
17
      0
             0
                 0
                     0
                        0
                            0
                                0
                                       0
                                           0
                                               0
                                                  1
                                                      -4
                                                              0
                                                                 0
                                                                      0
                                                                         0
         0
                                    1
                                                          1
18
19
      0
         0
             0
                 0
                     0
                        0
                            0
                                0
                                    0
                                       1
                                           0
                                               0
                                                  0
                                                      2
                                                         -4
                                                             0
                                                                 0
                                                                      0
                                                                          0
      0
         0
             0
                 0
                     0
                        0
                            0
                                0
                                    0
                                       0
                                           1
                                               0
                                                   0
                                                       0
                                                          0
                                                             -4
                                                                 1
                                                                      1
                                                                         0
20
         0
             0
                 0
                    0
                        0
                            0
                                0
                                    0
                                       0
                                           0
                                               1
                                                   0
                                                      0
                                                          0
                                                             1
                                                                 -4
                                                                      0
21
      0
                                                                         1
      0
         0
             0
                 0
                     0
                        0
                            0
                                0
                                   0
                                       0
                                           0
                                               0
                                                   0
                                                       0
                                                          0 2
                                                                 0 -4
22
                                0
                                               0
     0
         0
             0
                 0
                     0
                        0
                            0
                                    0
                                       0
                                           0
                                                   0
                                                       0
                                                          0
                                                              0
                                                                         -4
23
24
    b:
      0
25
      0
26
27
      0
      0
28
29
      0
30
      0
      0
31
32
      0
33
      0
34
35
      0
     0
36
    -15
37
    -15
38
    -15
39
40
     0
    -15
41
    0
42
43
    -15
   A is positive definite: False
44
   A' is positive definite: True
45
46
   === Question 3(b) ===
47
48
   Choleski x:
49
     0.96
     1.86
50
51
     2.61
      3.04
52
      3.17
53
54
      1.97
      3.88
55
      5.53
56
      6.37
57
      6.61
58
59
      3.03
      6.18
60
     9.25
61
62
     10.29
     10.55
63
     3.96
64
65
      8.56
     4.25
66
67
     9.09
    Conjugate gradient x:
68
     0.96
69
70
      1.86
71
      2.61
      3.04
72
73
      3.17
      1.97
74
75
      3.88
      5.53
```

```
6.37
77
78
      6.61
      3.03
79
      6.18
80
81
      9.25
     10.29
82
     10.55
83
84
      3.96
      8.56
85
      4.25
87
88
89
    === Question 3(c) ===
90
    === Question 3(d) ===
91
    Choleski potential at (0.06, 0.04): 5.52634126517 V
92
    Conjugate gradient potential at (0.06, 0.04): 5.52634127414 \mbox{\em V}
93
```

C Simple2D Data Files

Listing 13: Input mesh for the SIMPLE2D program.

```
1 1 0.0 0.0
   2 0.02 0.0
   3 0.04 0.0
   4 0.06 0.0
   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
   10 0.06 0.02
   11 0.08 0.02
11
12
   12 0.1 0.02
   13 0.0 0.04
13
   14 0.02 0.04
14
   15 0.04 0.04
   16 0.06 0.04
16
   17 0.08 0.04
17
   18 0.1 0.04
18
   19 0.0 0.06
19
   20 0.02 0.06
20
   21 0.04 0.06
21
   22 0.06 0.06
22
   23 0.08 0.06
   24 0.1 0.06
24
25
   25 0.0 0.08
   26 0.02 0.08
   27 0.04 0.08
27
   28 0.06 0.08
   29 0.08 0.08
29
   30 0.1 0.08
30
   31 0.0 0.1
   32 0.02 0.1
32
   33 0.04 0.1
33
   34 0.06 0.1
35
   1 2 7 0
36
   2 3 8 0
37
   3 4 9 0
38
   4 5 10 0
   5 6 11 0
40
   7 8 13 0
41
   8 9 14 0
   9 10 15 0
43
44 10 11 16 0
   11 12 17 0
```

```
13 14 19 0
46
47
    14 15 20 0
    15 16 21 0
48
    16 17 22 0
49
50
    17 18 23 0
    19 20 25 0
51
    20 21 26 0
52
53
    21 22 27 0
    22 23 28 0
54
    23 24 29 0
    25 26 31 0
56
    26 27 32 0
57
    27 28 33 0
58
    2 7 8 0
59
    3 8 9 0
60
    4 9 10 0
61
    5 10 11 0
62
63
    6 11 12 0
    8 13 14 0
64
    9 14 15 0
65
66
    10 15 16 0
    11 16 17 0
67
68
    12 17 18 0
69
    14 19 20 0
    15 20 21 0
70
71
    16 21 22 0
72
    17 22 23 0
    18 23 24 0
73
    20 25 26 0
    21 26 27 0
75
    22 27 28 0
76
    23 28 29 0
77
    24 29 30 0
78
    26 31 32 0
79
    27 32 33 0
80
    28 33 34 0
81
82
    1 0
83
    2 0
84
85
    3 0
    4 0
86
87
    5 0
    6 0
88
    7 0
89
    13 0
    19 0
91
    25 0
92
93
    31 0
    28 15
94
    29 15
95
    30 15
96
    34 15
97
```

Listing 14: Resulting potentials generated by the SIMPLE2D program.

```
ans =
1
2
        1.0000
                        0
                                  0
                                             0
3
        2.0000
                   0.0200
                                  0
                                             0
4
                                  0
        3.0000
                   0.0400
                                             0
5
6
        4.0000
                   0.0600
                                  0
                                             0
        5.0000
                   0.0800
                                  0
                                             0
7
        6.0000
                   0.1000
                                  0
                                             0
9
        7.0000
                             0.0200
                   0.0200
        8.0000
                                       0.9571
                             0.0200
10
                   0.0400
                                       1.8616
11
        9.0000
                             0.0200
       10.0000
                   0.0600
                             0.0200
                                       2.6060
12
       11.0000
                                       3.0360
                   0.0800
                             0.0200
13
       12.0000
                   0.1000
                             0.0200
                                       3.1714
```

15	13.0000	0	0.0400	0
16	14.0000	0.0200	0.0400	1.9667
17	15.0000	0.0400	0.0400	3.8834
18	16.0000	0.0600	0.0400	5.5263
19	17.0000	0.0800	0.0400	6.3668
20	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