

Bioinformatics III

Sixth Assignment

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June 2, 2018

Exercise 5.1: Boolean Networks

- (a) Listing 1 shows the source code of our propagation matrix class, which behaves like such a matrix. Internally it uses a adjacency matrix to efficiently calculate next states. Therefore it depends on the class AdjacencyMatrix shown in Listing 2. Further, the networks states are encoded in the class State, see 3.

Listing 1: Listing of source code

```
0 from os.path import exists
  from AdjacencyMatrix import AdjacencyMatrix
  from State import State

  class PropagationMatrix:
5      """
      Implements a propagation matrix. The propagated states are not
      preprocessed and are computed on demand to reduce the required
      amount of memory. Thereby increased runtime-changes are negligible
      """
10     def __init__(self, filename):
        """
        Create a class object behaving like a propagation matrix
        for a gene network obtained by textfile
        Textfile specification:
15     A > B indicates that gene A turns gene B activate
        A | B indicates that gene A turns gene B inactive
        (space separated)
        """

        # temporary storage for gene labels (to avoid redundancy)
20     nodes = set()
        # temporary storage for linkages (to avoid double file reading)
        links = list()
        if exists(filename):
            with open(filename) as openfile:
25                 for line in openfile:
                    content = line[0:(len(line)-1)].split(" ")
                    nodes.add(content[0])
                    nodes.add(content[2])
                    if (content[1] == ">"):
30                        links.append((content[0], content[2], 1))
                    elif (content[1] == "|"):
                        links.append((content[0], content[2], -3))

        # sort genes in alphabetiv order by
35     # by making use of list functionalities
        nodes = list(nodes)
        nodes.sort()

        # initialize a adjacency matrix representing the obtained
40     # gene network
```

```
        self.matrix = AdjacencyMatrix(0, nodes)
        for triple in links:
            self.matrix.setByLable(triple[0], triple[1], triple[2])
    else:
        print(filename, "does_not_exist")

45
def propagate(self, state):
    """
    Derive the propagated state form a given state
    Each gene is set inactive if the scalar product
    50 of the given state and its column in the adjacency matrix
    is smaller or equal to zero. Otherwise, the gene is set active
    """
    prop_state = State(state.getLables())
    for lab_y in state.getLables():
        # scalar product
        temp = 0
        for lab_x in state.getLables():
            temp += state.getByLable(lab_x) * self.matrix.getByLable(lab_x, lab_y)
        60 # threshold task specific
        prop_state.setByLable(lab_y, 0 if temp <= 0 else 1)
    return prop_state

def size(self):
    65 """
    Returns the number of states
    """
    return self.matrix.size()

70 def basinsAndAttractor(self, state):
    """
    Track the states obtained by simulating the regulatory
    network until the states rePEAT. Return the results divided in
    orbit states and pure basins
    75 """
    order = list()
    order.append(state.getInt())
    temp = 0
    # simulate network until repetition starts
    80 while True:
        state = self.propagate(state)
        temp = state.getInt()
        if not (temp in order):
            order.append(temp)
        85 else:
            break
    # divide states in upper described distinct subsets
    result = [[], []]
    switch = 0
    90 for i in range(0, len(order)):
        if order[i] == temp:
            result[0] = order[0:i]
            result[1] = order[i:len(order)]
    return result

95
def simplify(self, basatt):
    """
    Simplify redundand and incomplete basins of attractions and
    orbits
    100 """
    # Since every state can have only one next propagated state,
    # a minimum state-value is unique for every orbit and can be
    # used as key for our simplification
    attractors = dict()
    105 basins = dict()
    for ba in basatt:
        temp = min(ba[1])
```

```

    if not (temp in attractors.keys()):
        attractors[temp] = ba[1]
110     basins[temp] = ba[0]
    else:
        basins[temp].extend(ba[0])
        basins[temp].extend(ba[1])
    simple = list()
115     for i in attractors.keys():
        simple.append([set(basins[i]), attractors[i]])
    return simple

def orbit(self):
120     """
        Find basins of attraction and according orbits
        """
    # Storage for for attractors and basins
    periodes = list()
125     # simulate the network starting from every possible state
    candidates = list(range(0, 2**self.matrix.size()))
    while len(candidates) > 0:
        state = State(self.matrix.getLables())
        state.setInt(candidates[0])
130         # simulate regulatory network
        temp = self.basinsAndAttractor(state)
        # store obtained results in the according categorie,
        # further remove observed states from the candidate list
        # to minimize runtime
135         for basin in temp[0]:
            if basin in candidates:
                candidates.remove(basin)
        for attractor in temp[1]:
            if attractor in candidates:
140                 candidates.remove(attractor)
        periodes.append(temp)
    # results might contain reduncdancies, therefore a last
    # simlification step is possible
    return self.simplify(periodes)

```

Listing 2: Listing of source code

```

0 class AdjacencyMatrix:
    """
    Adjancency matrix, encoding a squared matrix with edge weight
    information between nodes.
    Initilialization and data access is only possible over the
5     row and column lables, not by positions
    """

    def __init__(self, initial, lables):
        """
10         Initialization of the adjacency matrix, required are an
            initial default weight and row lables, which are also used as
            column lables -> leading to a squared matrix
            """
        self.lables = lables
15         # default setup
        self.matrix = [0] * len(lables)
        for i in range(0, len(self.matrix)):
            self.matrix[i] = [initial] * len(lables)

20         # Similar to the state class, data access is only possible
            # with row and column lables, dicts enable an acces in linear
            # time
        self.access = dict()
        counter = 0
25         for i in self.lables:
            self.access[i] = counter
            counter += 1

```

```
def setByLable(self, a, b, value):
    """
    Set the weight for the edge from a to b
    """
    self.matrix[self.access[a]][self.access[b]] = value

def getByLable(self, a, b):
    """
    Get the weight for the edge from a to b
    """
    return self.matrix[self.access[a]][self.access[b]]

def size(self):
    """
    Returns the size of the matrix
    expressed by the number of lables
    """
    return len(self.lables)

def getLables(self):
    return self.lables

def show(self):
    """
    Output the matrix as collection of lists
    """
    print("Square_Matrix:")
    for i in range(0, len(self.lables)):
        print(self.matrix[i])
```

Listing 3: Listing of source code

```
class State:
    """
    Encode if genes are active (1) or inactive (0) or inactive
    Covers a simple list and extends it with functionality required
    by the PropagationMatrix and SquareMatrix class
    Data acces can only be done by using the gene lables.
    """

    def __init__(self, lables):
        """
        Initialize a state with the set of all gene lables
        """
        self.lables = lables
        # dict allows to access the data by lable in linear time
        self.access = dict()
        counter = 0
        for l in lables:
            self.access[l] = counter
            counter += 1
        # encondes if genes are active or inactive
        self.state = [0] * len(lables)

    def setByLable(self, lable, value):
        """
        Set the gene with a given lable avtive or inactive
        Every input will be translated to active (1) or inactive (0)
        """
        self.state[self.access[lable]] = 0 if value <= 0 else 1

    def getByLable(self, lable):
        """
        Returns 1 if the gene with a certain lable is active,
        0 otherwise
        """
        return self.state[self.access[lable]]
```

```
35     def getLables(self):
        """
        Returns the lables of genes
        """
40     return self.lables

    def size(self):
        """
        Returns the lenght of the state,
        i.g. the number of encoded genes
45     """
        return len(self.lables)

    def getInt(self):
        """
50     Returns the unique integer obtained by the binary encoded
        genes (active or inactive)
        """
        value = 0
55     for n in range(0, len(self.lables)):
        value += self.state[n] * (2**n)
        return value

    def setInt(self, value):
60     """
        Initializes a state whose binary representation equals the
        provided integer value
        """
        if value == 0:
65         return
        binaries = 2**(len(self.lables)-1)
        pos = len(self.lables) - 1
        while pos >= 0:
            if value >= binaries:
70                 self.state[pos] = 1
                value -= binaries
                binaries /= 2
                pos -= 1
        return

75     def show(self):
        """
        Output the state as binary list
        """
80     print("State:")
        print(self.state)
```

- (b) Listing 4 shows source code applying the the functionality of the code shown in Listing 1, which includes the network simulation.

1) It makes sense to stop the propagation when a state is observed a second time, from then on we will only observe orbiting behavior of the network states. The results of 2) are shown in this way. e.g. the first repeting state is the last shown.

2) Programs output for the required initial states:

Initial state 1:

1 -> 3 -> 7 -> 23 -> 55 -> 63 -> 13 -> 1

Initial state 4:

4 -> 18 -> 36 -> 26 -> 4

Initial state 21:

21 -> 51 -> 47 -> 13 -> 1 -> 3 -> 7 -> 23 -> 55 -> 63 -> 13

Initial state 33:

33 -> 11 -> 5 -> 19 -> 39 -> 31 -> 5

Listing 4: Listing of source code

```

0 from PropagationMatrix import PropagationMatrix
  from State import State

  def trackPropagation(state, repeats):
      """
      5   Visualize the propagations by a sequence of integers
          """
          track = str(state.getInt())
          for i in range(1, repeats):
              state = prop.propagate(state)
              track += " -> "
              track += str(state.getInt())
          print(track)

      # Initialize propagation network with text file
      15 # File contains structural informations of the given
          # gene regulatory network
          prop = PropagationMatrix("net.txt")

          # initialize with the state integer 13
      20 state_a = State(['A', 'B', 'C', 'D', 'E', 'F'])
          print("-----Exercise 6.1 b)-----")
          print("-----")
          print("\nInitial state 1:")
          state_a.setInt(1)
      25 trackPropagation(state_a, 8)

          # initialize with the state integer 13
          print("\nInitial state 4:")
          state_b = State(['A', 'B', 'C', 'D', 'E', 'F'])
      30 state_b.setInt(4)
          trackPropagation(state_b, 5)

          # initialize with the state integer 13
          print("\nInitial state 21:")
      35 state_c = State(['A', 'B', 'C', 'D', 'E', 'F'])
          state_c.setInt(21)
          trackPropagation(state_c, 11)

          # initialize with the state integer 13

```

```
40 print("\nInitial_state_33:")
    state_d = State(['A', 'B', 'C', 'D', 'E', 'F'])
    state_d.setInt(33)
    trackPropagation(state_d, 7)

45 print()
    print("-----Exercise_6.1_c)-----")
    print("-----")
    orbits = prop.orbit()
    for i in range(0, len(orbits)):
50     print()
        length = len(orbits[i][1])
        print("Orbit_" + str(i + 1) + "_with_length_" + str(length) + ":")
        print(orbits[i][1])
        print("Set_of_basins:")
55     print(orbits[i][0])
        coverage = float(len(orbits[i][0]))
        coverage /= float(2**prop.size())
        coverage *= 100.0
        print("Relative_coverage:_" + str(coverage) + "%")
```

(c) Output of the program listing the orbits:

Orbit 1 with length 1:

0

Set of basins:

0, 6, 8, 12, 16, 20, 22, 24, 28, 32, 34, 40, 42, 44, 46, 48, 50, 52, 54, 56, 58,
60, 62

Relative coverage: 35.9375%

Orbit 2 with length 7:

1, 3, 7, 23, 55, 63, 13

Set of basins:

1, 3, 7, 9, 13, 21, 23, 25, 29, 41, 43, 45, 47, 49, 51, 53, 55, 57, 59, 61, 63

Relative coverage: 32.8125%

Orbit 3 with length 4:

4, 18, 36, 26

Set of basins:

2, 4, 36, 38, 10, 14, 18, 26, 30

Relative coverage: 14.0625%

Orbit 4 with length 4:

5, 19, 39, 31

Set of basins:

33, 35, 5, 37, 39, 11, 15, 17, 19, 27, 31

Relative coverage: 17.1875%

(d) Listing ?? shows source code.

Exercise 5.2: Differential Expression Analysis