nb-2021-04-15

April 29, 2021

1 Exercises for 2021-04-29

Note: You will need to do import Pkg; Pkg.add("GraphRecipes") if you do not already have it installed.

```
[]: # import Pkg; Pkg.add("GraphRecipes")

[]: using LinearAlgebra
using Plots
using GraphRecipes
using SparseArrays
using MatrixNetworks
```

1.1 Data setup

The Zachary karate club is a standard example in network analysis; for this and some other examples, see Mark Newman's web site, the Network Repository, and the KONECT repository. Other example repositories include SNAP and UCI.

```
[ ]: m = 78
     n = 34
     adj = [1 \ 2]
               1 3
               2 3
               1 4
               2 4
               3 4
               1 5
               1 6
               1 7
               5 7
               6 7
               18
               2 8
               3 8
               4 8
               1 9
               3 9
```

```
3 10
1 11
5 11
6 11
1 12
1 13
4 13
1 14
2 14
3 14
4 14
6 17
7 17
1 18
2 18
1 20
2 20
1 22
2 22
24 26
25 26
3 28
24 28
25 28
3 29
24 30
27 30
2 31
9 31
1 32
25 32
26 32
29 32
3 33
9 33
15 33
16 33
19 33
21 33
23 33
24 33
30 33
31 33
32 33
9 34
10 34
14 34
```

```
15 34
      16 34
      19 34
      20 34
      21 34
      23 34
      24 34
      27 34
      28 34
      29 34
      30 34
      31 34
      32 34
      33 34]
A = sparse(adj[:,1], adj[:,2], ones(m), n, n)
A = A+A'
spy(A)
```

```
[]: graphplot(A, markersize=0.4, curves=false, names=labels)
```

1.2 Semi-supervised labeling

Suppose we label node 1 as 0 and node 34 as 1.

(4 points): Compute the vector of "soft" labels using the Laplacian-based approach from the 4/22 lecture. Plot the soft labels vs the "ground truth" labels. Does this approach work well?

[]:

1.3 Spectral partitioning

This is a small enough matrix that we can just use the Julia eigen function. For larger matrices, we would want to use eigs (which is now in the Arpack.jl package).

We start by plotting the labeling vs the first eigenvector of the combinatorial Laplacian. What we notice is that this ordering does a pretty good job of separating the two "ground truth" clusters in the graph.

```
[]: d = [sum(A[i,:]) for i = 1:n]
D = spdiagm(0 => d)
L = D-A
Ldense = Matrix(L)
s, V = eigen(Ldense)
spectral1d = V[:,2]
scatter(spectral1d, labels')
```

(3 points): In your own words, explain what was meant by "does a pretty good job of separating the two 'ground truth' clusters in the graph."

The spectral approximation to maximizing the modularity (defined in terms of $B = A - \frac{dd^T}{2m}$ involves similarly looking at the eigenvector associated with the largest eigenvalue of B.

(3 points): Plot the "ground truth" labels against the components of this eigenvector – you should see it again does a pretty good job of separating out the two pieces.

[]:

1.4 Students choice

Suppose you had to ask a question to gently probe students knowledge of the graph learning methods discussed in the past two weeks.

(2 points): What question do you think you would ask?

(2 points): What would your answer to that question be?