Module IV - Basic Analysis

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Loading data from multiple sources

- Local network data files
- Building directly from the Internet

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Brief review of Python dictionaries

- Why is the di ct so useful?
- ► How NetworkX utilizes it?

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- Calculating degree distribution
- Plotting statistics using matpl otlib
- Calculating cliques, clustering and transitivity

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Basic visualization

- Review of NetworkX's plotting algorithms
- Adding analysis to visualization

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► For those that are unfamiliar with working at the command-line, however, the process can be confusing



Let's try!

- We will load the edge list of Hartford drug users network
- Specify that the network be a directed graph, and the nodes be integers
- Use nx. i nfo() to check that data has been loaded correctly

Starting NetworkX and loading data

- $1 >>> hartford=nx.read_edgelist(".../.../data/hartford_drug.txt",create_using=nx.DiGraph(),nodetype=int)$
- 2 >>> nx.info(hartford)
- 3 Name:

4 Type: DiGraph
5 Number of nodes: 212
6 Number of edges: 337
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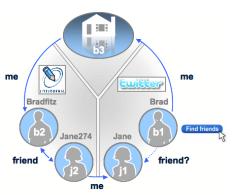
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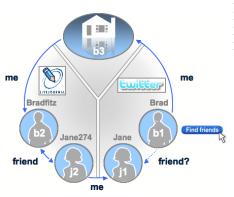
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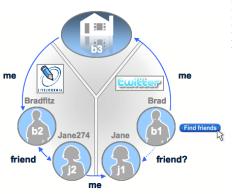
Some formats may have more or less options, always check the documentations!



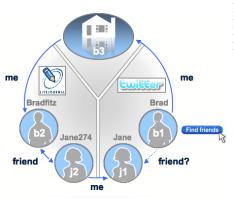


Perhaps the most powerful aspect of NetworkX is its ability to work in Python to generate networks from live-streaming data

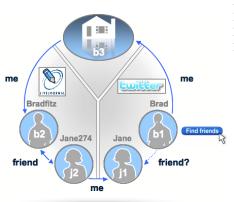
 In Python, use NetworkX, cj son and a other standard scientific libraries to parse Google's SocialGraph data



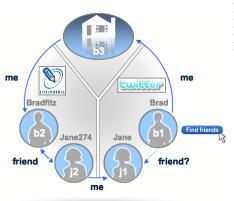
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- Through a process called "k-snowball searching" $seed \rightarrow friend \rightarrow \cdots \rightarrow friend_k$
 - Seed: imichaeldotorg.livejournal.com
 - ▶ k = 3
- Note the low value of k

The code, part 1

Loading the libraries

```
from cjson import *
  from urllib import *
   from time import *
   from scipy import array, unique
   if __name__ == "__main__":
       seed="imichaeldotorg"
 7
       seed_url="http://"+seed+".livejournal.com"
 8
 9
       sg=get sg(seed url)
10
       net,newnodes=create_egonet(sg)
11
       nx.write paiek(net."../../data/"+seed+" ego.net")
12
       nx. info (net)
13
```

Scraping egonet relationships from seed

```
1 def get_sg(seed_url):
    sgapl_url="http://socialgraph.apis.google.com/lookup?q="+seed_url+"&edo=1&edi=1&fme=1&pretty=0"
    try:
        furl=urlopen(sgapi_url)
        f=furl.read()
        furl.close()
        return fr
        except IOError:
        print "Could not connect to website"
        print sgapi_url
        return {}
```

Creating the egonet

```
def create egonet(s):
 3
            raw=decode(s)
            G=nx.DiGraph()
            pendants=[]
            n=raw['nodes']
            nk=n.kevs()
 7
            G.name=str(nk)
 9
            pendants=[]
10
            for a in range(0.len(nk)):
                for b in range(0,len(nk)):
11
12
                    if al=b:
                        G.add edge(nk[a].nk[b])
13
            for k in nk:
14
15
                ego=n[k]
                ego_out=ego['nodes_referenced']
16
                for o in ego_out:
18
                    G. add edge(k.o)
                    pendants.append(o)
19
                ego in=ego['nodes referenced by']
20
21
                for i in ego_in:
22
                    G.add_edge(i,k)
                    pendants.append(i)
23
24
            pendants=array (pendants, dtype=str)
25
            pendants. flatten()
26
            pendants=unique (pendants)
27
            return G, pendants
        except DecodeError:
28
29
30
        except KeyError:
```

Rolling the snowball

```
def snowball round (G. seeds . myspace=False ) :
2
        t0=time()
3
        if myspace:
            seeds=get myspace url(seeds)
5
        sb_data=[]
6
        for s in range(0, len(seeds)):
7
            s sg=get sg(seeds[s])
            new_ego,pen=create_egonet(s_sg)
9
            for p in pen:
10
                    sb_data.append(p)
            if s<1:
11
12
                sb net=nx.compose(G.new ego)
13
                sb_net=nx.compose(new_ego, sb_net)
14
15
            del new ego
            if s==round(len(seeds)*0.2):
16
                sb net.name='20% complete'
17
18
                nx.info(sb net)
19
                print 'AT: '+strftime('%m/%d/%Y, %H:%M:%S', gmtime())
20
                print '
21
22
       # More time keeping, probably a MUCH better way to do this
       sb data=array(sb data)
23
24
        sb data. flatten()
       sb data=unique(sb data)
25
26
       nx. info (sb net)
27
       return sb_net,sb_data
```

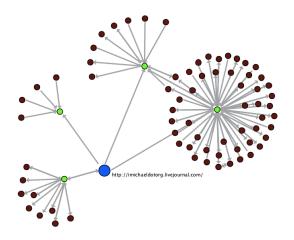
Step	Nodes	Edges	Mean Degree	Density
Seed	5	5	2.0	0.25
k = 2	75	115	3.0	0.02
k = 3	4,938	8,659	3.5	$3.6(10^{-4})$

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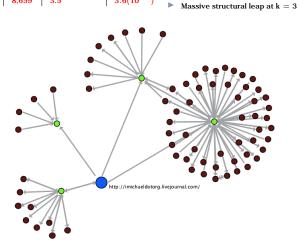
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- Large jump after first snowball

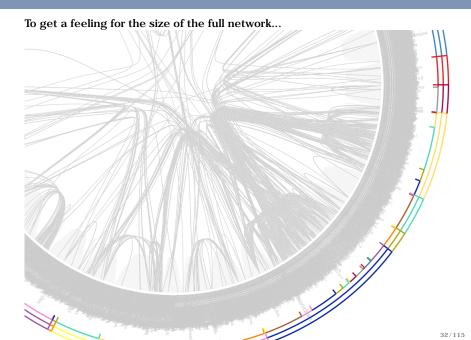


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The full network



Python Dictionaries

The dict type is a data structure that represents a key→value mapping

Working with the dict type

```
# Keys and values can be of any data type

>>> fruit_dict={"apple":1,"orange":[0.23,0.11],"banana":True}

# Can retrieve the keys and values as Python lists (vector)

>>> fruit_dict.keys()

["orange","apple","banana"]

# Or create a (key,value) tuple

>>> fruit_dict.items()

[("orange",[0.23,0.11]),("apple",1),("Banana",True)]

# This becomes especially useful when you master Python ''list comprehension''
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Now, try creating a di ct of your own

Using dictionaries for network analysis

From the documentation...

networkx.closeness centrality Compute closeness centrality for nodes.

closeness centrality(G, v=None, weighted edges=False)

Closeness centrality at a node is 1/average distance to all other nodes

Parameters: G : graph

A networkx graph

v : node, optional

Return only the value for node v.

weighted_edges : bool, optional

Consider the edge weights in determining the shortest paths. If False, all edge weights are considered equal.

Returns: nodes : dictionary

> Dictionary of nodes with closeness centrality as the value.

NetworkX's metric's make extensive use of the dict type

▶ In this case the key→value mapping is of the form: { node-l abel : metric}

Let's look at an example:

```
1 >>> in cen=nx.in degree centrality(hartford)
2 >>> in cen
3 \{1: 0.014218009478672987, 2: 0.018957345971563982....
  90: 0.0047393364928909956, 293: 0.0\}
```

We can see that node #90 has in-degree centrality 0.0047

But we can do so much more!

Running multiple measures

For our first analysis in NetworkX, we will do some basic network manipulation, then run multiple measures to find highest centrality nodes

First, we will need to convert to an undirected network, and extract the main component

```
# We will symmetrize for simplcity

2 >>> hartford_ud=hartford.to_undirected()

3 # The network also has many small components, but for

4 # this analysis we are interested in the largest

5 >>> hartford_mc=hartford_main=nx.connected_component_subgraphs(hartford_ud)[0]
```

Next, we will calculate multiple measures

```
1 # Betweenness centrality
2 >>> bet_cen=nx.betweenness_centrality(hartford_mc)
3 # Closeness centrality
4 >>> clo_cen=nx.closeness_centrality(hartford_mc)
5 # Eigenvector centrality
6 >>> eig_cen=nx.eigenvector_centrality(hartford_mc)
```

Finding most central actors

To find the most central actors we will use Python's list comprehension technique to do basic data manipulation on our centrality dictionaries

```
1 def highest_centrality(cent_dict):
2    """Returns node key with largest value from
3    NX centrality dict""
4    # Create ordered tuple of centrality data
5    cent_items=cent_dict.items()
6    # List comprehension!
7    cent_items=[(b, a) for (a, b) in cent_items]
8    # Sort in descending order
9    cent_items.ort()
10    cent_items.reverse()
11    return cent_items[0][1]
```

Now, just ask for the answer

Finding Most central actors

```
1 >>> print("Actor "+str(highest_centrality(bet_cen))+" has the highest Betweenness centrality")
2 Actor 82 has the highest Betweenness centrality
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List comprehension

```
    Given a dict: d={1: 0.15, 2: 0.67}
    d.items() → [(1, 0.15), (2, 0.67)]
    d=[(b, a) for (a, b in d)] → [(0.15, 1), (0.67, 2)]
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Here, we use list comprehension in order to use Python's built-in sort and reverse list functions

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Calculating degree distribution

One of the most popular network level statistical description of a network is its degree distribution

▶ In NetworkX this is a simply one-line operation

Get list of degree rank frequency 1 # Create a Barabasi-Albert network 2 >>> ba_net=barabasi_albert_graph(1000,2) 3 # Built-in function for degree distribution 4 >>> dh=degree_histogram(ba_net)

Calculating degree distribution

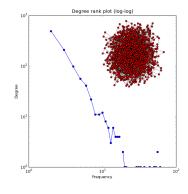
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- ► As we will see next, we can use matplotlib to take this data and create publication ready plots
- ► Ex. from http://networkx.lanl.gov/examples/ drawing/degree_histogram.html



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Cliques

Maximal cliques are the largest complete subgraph containing a given point. There are several algorithms for finding cliques, including Bron Kerbosch (1973), Tomita, Tanaka and Takahashi (2006), Cazals and Karande (2008)

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For each node find the fraction of possible triangles that exist, $c_v = \frac{2T(v)}{\deg(v)(\deg(v)-1)}$, where T(v) is the number of triangles through node v.

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We will use clustering coefficients to identify community structure in the Hartford drug network

- 1 # Calculate clustering coefficients of each node (return as dict)
- 2 clus=clustering(hartford_mc, with_labels=True)
- 3 # Get counts of nodes membership for each clustering coefficient, and clean up
- 4 unique_clus=list(unique(clus.values()))
- 5 clus_counts=zip(map(lambda c: clus.values().count(c),unique_clus),unique_clus)
- 6 clus_counts.sort()
- 7 clus counts.reverse()
- 8 # Create a subgraph from nodes with most frequent clustering coefficient
- 9 mode_clus_sg=subgraph(hartford_mc,[(a) for (a,b) in clus.items() if b==clus_counts[0][1]])

Calculating clustering coefficients

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Use the with-labels to return a dict keyed by node label

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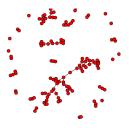
- Use the with-labels to return a dict keyed by node label
- ► The zi p function takes two lists and returns a tuple

```
1 # Calculate clustering coefficients of each node (return as dict)
2 clus=clustering(hartford_mc, with_labels=True)
3 # Get counts of nodes membership for each clustering coefficient, and clean up
4 unique_clus=list (unique(clus.values()))
5 clus_counts=zip(map(lambda c: clus.values().count(c).unique_clus), unique_clus)
6 clus_counts.sort()
7 clus_counts.reverse()
8 # Create a subgraph from nodes with most frequent clustering coefficient
9 mode_clus_sg=subgraph(hartford_mc,[(a) for (a,b) in clus.items() if b==clus_counts[0][1]])
```

- Use the with-labels to return a dict keyed by node label
- ► The zi p function takes two lists and returns a tuple
- More complex list comprehension with logic operator

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# Calculate clustering coefficients of each node (return as dict)
clus=clustering (hartford_mc, with_labels=True)
# Get counts of nodes membership for each clustering coefficient, and clean up
unique_clus=list (unique(clus.values()))
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clus_counts.sort()
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Calculating clustering coefficients

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- ► Use the with—labels to return a dict keyed by node label
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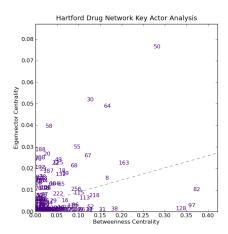
Later, we'll learn how to create a network visualization like the one above

Recall Python's scientific computing trinity: NumPy, Sci Py and mat pl ot lib

While NumPy and Sci Py do most of the behind the scenes work, you will interact with mat pl ot l i b frequently for when doing network analysis

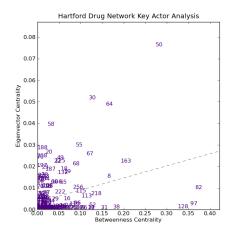
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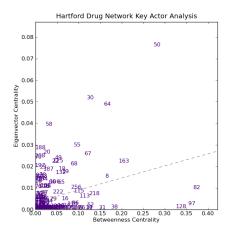
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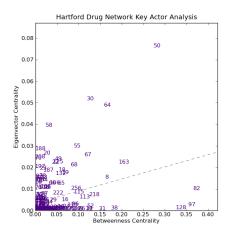


We will need to create a function that takes two centrality di ct and generates this plot

1. Create a matpl otlib figure

Recall Python's scientific computing trinity: NumPy, Sci Py and mat pl ot lib

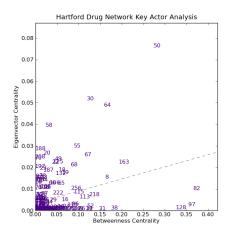
While NumPy and Sci Py do most of the behind the scenes work, you will interact with mat pl ot l i b frequently for when doing network analysis



- 1. Create a matpl otlib figure
- 2. Plot each node label as a point

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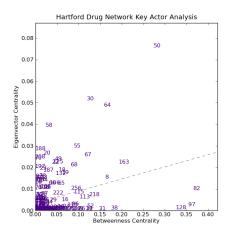
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- 1. Create a matplotlib figure
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- 3. Add a "best fit" line

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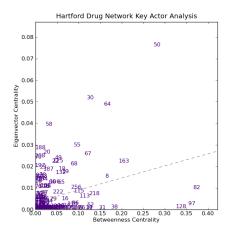
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- 1. Create a matplotlib figure
- 2. Plot each node label as a point
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- 4. Add axis and title labels

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- 1. Create a matplotlib figure
- 2. Plot each node label as a point
- 3. Add a "best fit" line
- 4. Add axis and title labels
- 5. Save figure as a PNG file

The central ity_scatter function, part one

```
def centrality scatter(met dict1, met dict2, path="", vlab="", xlab="", title="", reg=False):
       # Create figure and drawing axis
2
       fig=P. figure (figsize = (7,7))
3
       ax1=fig.add subplot(111)
       # Create items so actors can be sorted properly
       met items1=met dict1.items()
7
       met items2=met dict2.items()
       met items1.sort()
8
       met items2.sort()
9
       # Grab data
10
       xdata=[(b) for (a,b) in met_items1]
11
       ydata=[(b) for (a,b) in met_items2]
12
       # Add each actor to the plot by ID
13
       for p in xrange(len(met items1)):
14
           ax1.text(x=xdata[p],y=ydata[p],s=str(met_items1[p][0]),color="indigo")
15
```

The central ity_scatter function, part one def centrality scatter(met dict1, met dict2, path="", vlab="", xlab="", title="", reg=False): # Create figure and drawing axis 2 fig=P. figure (figsize = (7,7))3 ax1=fig.add subplot(111) # Create items so actors can be sorted properly met items1=met dict1.items() 7 met items2=met dict2.items() met items1.sort() 8 met items2.sort() # Grab data 10 xdata=[(b) for (a,b) in met_items1] 11 ydata=[(b) for (a,b) in met_items2] 12 # Add each actor to the plot by ID 13 for p in xrange(len(met items1)): 14 ax1.text(x=xdata[p],y=ydata[p],s=str(met_items1[p][0]),color="indigo") 15

Create a canvas to draw on

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- Create a canvas to draw on
- manipulate and store centrality data

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```

- Create a canvas to draw on
- manipulate and store centrality data
- ► Add points to plot as node labels

The central ity_scatter function, part two

```
def centrality scatter (met dict1, met dict2, path="", vlab="", vlab="", title="", reg=False):
2
3
       # If adding a best fit line, we will use NumPv to calculate the points.
       if reg:
4
           # Function returns y-intercept and slope. So, we create a function to
5
           # draw LOBF from this data
           slope, vint=polyfit (xdata, ydata, 1)
7
           xline=P. xticks()[0]
8
           vline=map(lambda x: slope*x+vint.xline)
9
           # Add line
10
           ax1. plot(xline, yline, ls='---', color='grey')
11
       # Set new x- and y-axis limits to data
12
       P. xlim((0.0.max(xdata)+(.15*max(xdata))))
                                                     # Give a little buffer
13
14
       P. ylim((0.0, max(ydata) + (.15*max(ydata))))
15
       # Add labels
16
       ax1.set title(title)
17
       ax1.set xlabel(xlab)
       ax1.set ylabel(ylab)
18
       # Save figure
19
       P. savefig (path, dpi=100)
20
```

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Add a best fit line

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- Add a best fit line
- Resize figure to fit data

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18
       # Save figure
19
       P. savefig (path, dpi=100)
20
```

- Add a best fit line
- Resize figure to fit data
- Add labels, and save the figure as a PNG file

Exporting network data and analytics

As powerful as NetworkX and the complementing scientific computing packages in Python are, it may often be useful or necessary to output your data for additional analysis

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In most cases this will entail either exporting the raw network data, or metrics from some network analysis ${\bf r}$

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- 2. When you want to export metrics we can use Python's built-in XML and CSV libraries
- 3. Depending on your needs you may prefer one, the other or both

Next, we will review how to save data in different formats and export metrics to a CSV file using the Hartford drug net data



```
NX syntax for writing a network file

>>> nx.write_format(G, "path/to/file.txt", ...options...)

↑ ↑ ↑

NX function, net variable File to be written Nodes/edge data, etc.
```

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↑ ↑ ↑ ↑

NX function, net variable File to be written Nodes/edge data, etc.
```

The syntax for exporting network data follows exactly the syntax for loading it



Let's try!

- Output the Hartford drug net data as an adjacency list
- Add metric data to each node of the network
- Output new network in Pajek format with node attributes

Saving network data and adding node attributes

As shown, this is a simple one line operation

Output Hartford drug net data as an adjacency list

```
1 nx.write_adjlist(hartford_mc,"../../data/hartford_mc_adj.txt")
```

Next, we will add the Eigenvector centrality of each node to the graph object

Adding node attributes

```
def add_metric(G, met_dict):
    """Adds metric data to G from a dictionary keyed by node labels"""
if(G.nodes().sort()==met_dict.keys().sort()):
    for i in met_dict.keys():
        G.add_node(i, metric=met_dict[i])
    return G
else:
    raise ValueError("Node labels do not match")
```

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As shown, this is a simple one line operation

Output Hartford drug net data as an adjacency list

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Quick error checking

Saving network data and adding node attributes

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```

- Quick error checking
- Add node attribute as "metric"

Using the Python CSV library

Python has powerful built-in tools for reading and writing standard data formats

One of the most useful, and frequently used, is the CSV library and the Di ctWri ter

```
import csv
  def csv exporter(data dict.path):
       """Takes a dict of centralities keyed by column headers and exports
       data as a CSV file """
       # Create column header list
       col headers=["Actor"]
7
       col headers.extend(data dict.kevs())
       # Create CSV writer and write column headers
       writer=csv. DictWriter(open(path, "w"), fieldnames=col headers)
10
       writer.writerow(dict((h,h) for h in col headers))
11
       # Write each row of data
12
       for j in data dict[col headers[1]].keys():
13
           # Create a new dict for each row
14
           row=dict.fromkeys(col_headers)
15
           row["Actor"]= i
16
           for k in data dict.keys():
17
               row[k]=data dict[k][i]
18
           writer . writerow (row)
19
```

The results of CSV export

We can now open the CSV file in our favorite spreadsheet program

- Perform traditional data exploration
- ► Load into other analytics platforms for additional analysis (e.g., R)
- Store for latter use

3 2 0.12475634 0.01767427 0.0002596 4 3 0.12565445 0.05687441 0.0002318 5 4 0.10223642 0.03108639 1.44E-0 6 5 0.1443609 0 0.0031315 7 6 0.09943035 0.01041667 1.49E-0 8 7 0.11340815 0.04362093 6.78E-0 9 8 0.20512821 0.16354003 0.0147188 10 9 0.11267606 0.00741624 0.000111 11 10 0.13983977 0.05258239 0.0009545 12 11 0.1703638 0.01250999 0.0032333 13 13 0.13892909 0 1.79E-0 14 14 0.17219731 0.11848775 0.0002973 15 15 0.13521127 0.00079897 2.11E-0					
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4 3 0.12565445 0.05687441 0.0002318 5 4 0.10223642 0.03108639 1.44E-0 6 5 0.1443609 0 0.0031315 7 6 0.09943035 0.01041667 1.49E-0 8 7 0.11340815 0.04362093 6.78E-0 9 8 0.20512821 0.16354003 0.0147188 10 9 0.11267606 0.00741624 0.000110 11 10 0.13983977 0.05258239 0.00099545 12 11 0.1703638 0.01250999 0.003233 13 13 0.13892909 0 1.79E-0 14 14 0.17219731 0.11848775 0.0002973 15 15 0.13521127 0.00079897 2.11E-0	2	1	0.12467532	0.0072576	0.00025176
5 4 0.10223642 0.03108639 1.44E-0 6 5 0.1443609 0 0.0031315 7 6 0.09943035 0.01041667 1.49E-0 8 7 0.11340815 0.04362093 6.78E-0 9 8 0.20512821 0.16354003 0.0147188 10 9 0.11267606 0.00741624 0.000110 11 10 0.13983977 0.05258239 0.0009545 12 11 0.1703638 0.01250999 0.003233 13 13 0.13892909 0 1.79E-0 14 14 0.17219731 0.11848775 0.0002973 15 15 0.13521127 0.00079897 2.11E-0	3	2	0.12475634	0.01767427	0.00025964
6 5 0.1443609 0 0.0031315 7 6 0.09943035 0.01041667 1.49E-0 8 7 0.11340815 0.04362093 6.78E-0 9 8 0.20512821 0.16354003 0.0147188 10 9 0.11267606 0.00741624 0.000110 11 10 0.3983977 0.0528239 0.0009545 12 11 0.1703638 0.01250999 0.03233 13 13 0.133892909 0 1.79E-0 14 14 0.17219731 0.1184875 0.0002973 15 15 0.13521127 0.00079897 2.11E-0		3	0.12565445	0.05687441	0.00023185
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8 7 0.11340815 0.04362093 6.78E-0 9 8 0.20512821 0.16354003 0.0147188 10 9 0.11267606 0.00741624 0.000110 11 10 0.13983977 0.05258239 0.0009545 12 11 0.1703638 0.01250999 0.003233 13 13 0.13892909 0 1.79E-0 14 14 0.17219731 0.11848775 0.0002973 15 15 0.13521127 0.00079897 2.11E-0		5	0.1443609	0	0.00313152
9 8 0.20512821 0.16354003 0.0147188 10 9 0.11267606 0.00741624 0.000110 11 10 0.13983977 0.0528239 0.0009545 12 11 0.1703638 0.01250999 0.03233 13 13 0.13892909 0 1.79E-0 14 14 0.17219731 0.11848775 0.0002973 15 15 0.13521127 0.00079897 2.11E-0	7	6	0.09943035	0.01041667	1.49E-07
10 9 0.11267606 0.00741624 0.000110 11 10 0.13983977 0.05258239 0.0009545 12 11 0.1703638 0.01250999 0.003233 13 13 0.13892909 0 1.79E-0 14 14 0.17219731 0.11848775 0.0002973 15 15 0.13521127 0.00079897 2.11E-0		7	0.11340815	0.04362093	6.78E-05
11 10 0.13983977 0.05258239 0.0009545 12 11 0.1703638 0.01250999 0.03233 13 13 0.13892909 0 1.79E-0 14 14 0.17219731 0.11848775 0.0002973 15 15 0.13521127 0.00079897 2.11E-0		8	0.20512821	0.16354003	0.01471888
12 11 0.1703638 0.01250999 0.003233 13 13 0.13892909 0 1.79E-0 14 14 0.17219731 0.11848775 0.0002973 15 15 0.13521127 0.00079897 2.11E-0	10	9	0.11267606	0.00741624	0.0001101
13 13 0.13892909 0 1.79E-0 14 14 0.17219731 0.11848775 0.0002973 15 15 0.13521127 0.00079897 2.11E-0	11	10	0.13983977	0.05258239	0.00095456
14 14 0.17219731 0.11848775 0.0002973 15 15 0.13521127 0.00079897 2.11E-0.00079897	12	11	0.1703638	0.01250999	0.0032333
15 15 0.13521127 0.00079897 2.11E-0	13	13	0.13892909	0	1.79E-05
	14	14	0.17219731	0.11848775	0.00029737
16 0 45007200 0 06202647 0 0042202	15	15	0.13521127	0.00079897	2.11E-05
16 0.1590/208 0.0620364/ 0.0043283	16	16	0.15907208	0.06203647	0.00432838

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- Scale up to very large graphs
- Display nodal- (centrality) of network-level (community structure) information

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- ► Maximize "visibility" of network
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NetworkX was designed as a data manipulation and analysis tool, and therefore is not meant as a visualization platform

▶ It is, however, still capable of making very nice visualization



Visualization algorithms in NetworkX - Random & Circle

The most basic visualization techniques are the random and circular layouts

- The random layout places nodes in...random positions
- ► The circular layout places nodes in...a circle





```
# Use subplots to draw random and circular layouts
# of drug net side-by-side
fig1=P. figure (figsize = (9,4))
fig1.add_subplot(121)
s nx. draw_random (hartford_mc, with_labels=False, node_size=60)
fig1.add_subplot(122)
nx. draw_circular (hartford_mc, with_labels=False, node_size=60)
P. savefig("..././images/networks/rand_circ.png")
```

$\label{thm:condition} \textbf{Visualization algorithms in } \texttt{NetworkX-Spring \& Spectral}$

More commonly used visualization techniques include the spring and spectral layouts

Visualization algorithms in NetworkX - Spring & Spectral

More commonly used visualization techniques include the spring and spectral layouts

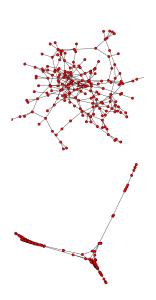
The spring layout is a version of the Fruchterman-Reingold force-directed algorithm, which attempts to minimize overlapping edges



Visualization algorithms in NetworkX - Spring & Spectral

More commonly used visualization techniques include the spring and spectral layouts

- The spring layout is a version of the Fruchterman-Reingold force-directed algorithm, which attempts to minimize overlapping edges
- The spectral layout finds node position using the eigenvectors of the graph Laplacian, which is useful for quickly visualizing structural clustering



The shell layout draws nodes as concentric circles

- Two dimensional extension of the circle layout
- We may have some reason to isolate certain nodes

```
1 P.figure (figsize = (8,8))
2 # Find actors in 25th percentile
3 max_eig=max([(b) for (a,b) in eig_cen.items()])
4 s1=[(a) for (a,b) in eig_cen.items() if b>=.25*max_eig]
5 s2=hartford_mc.nodes()
6 # setdiff1d is a very useful NumPy function!
7 s2=list (setdiff1d (s2,s1))
8 shells=[s1,s2]
9 # Calculate position and draw
10 shell_pos=shell_layout (hartford_mc, shells)
11 draw_networkx(hartford_mc, shell_pos, with_labels=False, node_size=60)
12 P.savefig("../.images/networks/shell.png")
```

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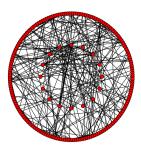
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draw_networkx(hartford_mc, shell_pos, with_labels=False, node_size=60)
P.savefig("../../images/networks/shell.png")
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25th percentile Eigenvector centrality actors

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12 P.savefig(".../,images/networks/shell.png")
```



Beyond layout, we may also want to add analytical data to our visualization

Changing node and edge size and colors

NetworkX allows you to alter the size, color and shape of the nodes and edges in any visualization

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In our final exercise, we will add the following analysis to the Hartford drug network

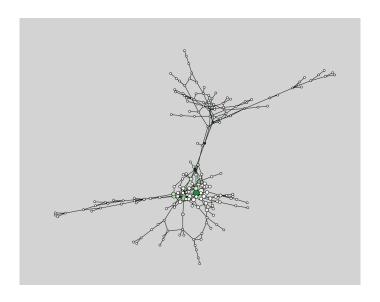
- Node size by Eigenvector centrality
- Intensity of node color by betweenness centrality
- Edge thickness by edge betweenness

The code to add analysis to visualization

More list comprehension and matplotlib colormaps

```
1 # Adding analysis to visualization
2 P. figure (figsize = (15.15))
3 P. subplot(111,axisbg="lightgrey")
  spring pos=nx.spring layout(hartford mc,iterations=1000)
  # Use betweeneess centrality for node color intensity
  bet color=bet cen.items()
7 bet color.sort()
8 bet color=[(b) for (a,b) in bet color]
  # Use Eigenvector centrality to set node size
  eig size=eig cen.items()
  eig size.sort()
12 eig size = [((b)*2000)+20 for (a,b) in eig size]
13 # Use matplotlib's colormap for node intensity
   draw_networkx(hartford_mc, spring_pos, node_color=bet_color, . . .
       ...cmap=P.cm, Greens, node_size=eig_size, with_labels=False)
16 P. savefig("../../images/networks/analysis.png")
```

Final visualization



Basic Analysis

 How to load local data, and an example of building networks from data streamed directly from the Internet

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- ▶ A brief review of the Python di ct data type

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Questions?