

## Module IV - Basic Analysis

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## Agenda for Module IV

### Loading data from multiple sources

- ▶ Local network data files
- ▶ Connecting to a database
- ▶ Building directly from the Internet

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Brief review of Python dictionaries

- ▶ Why is the `dict` so useful?
- ▶ How `NetworkX` utilizes it?

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### Running basic centralities

- ▶ Degree, Closeness, Betweenness Eigenvector
- ▶ Calculating degree distribution
- ▶ Plotting statistics using matplotlib
- ▶ Calculating cliques, clustering and transitivity

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- ▶ Writing network data
- ▶ Saving network analysis statistics

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### Basic visualization

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

As we have seen, one of the main advantages of working with NetworkX is that it can read many different network formats

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

```
>>> G = read_format("path/to/file.txt", ...options...)
      ↑           ↑           ↑
Net variable  NX function, file directory path  Graph type, nodes type, etc.
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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 `info()` to check that data has been loaded correctly

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- ▶ 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 `info()` to check that data has been loaded correctly

It's time to fire up your console and load Python!

# Loading the Hartford drug users network

## Starting NetworkX and loading data

```
>>> from networkx import *
>>> hartford=read_edgelist("../data/hartford_drug.txt",create_using=DiGraph(),nodetype=int)
>>> info(hartford)
Name:
Type:          DiGraph
Number of nodes: 212
Number of edges: 337
Average in degree: 1.5896
Average out degree: 1.5896
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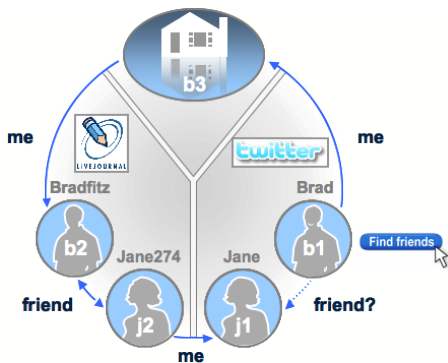
Some formats may have more or less options, **always check the documentations!**

## Building a network from a database

As data sets become larger and persistently changing, it may make more sense to store them in a database rather than a single file

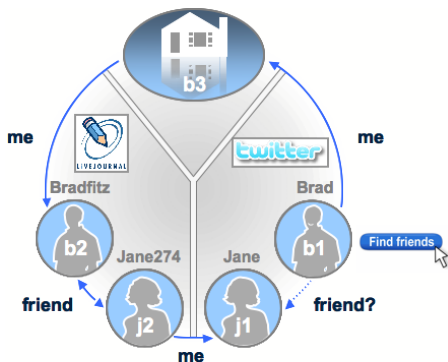
- ▶ As we have seen, Python provides binding to many modern database frameworks

## Building the social network among LiveJournal users



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

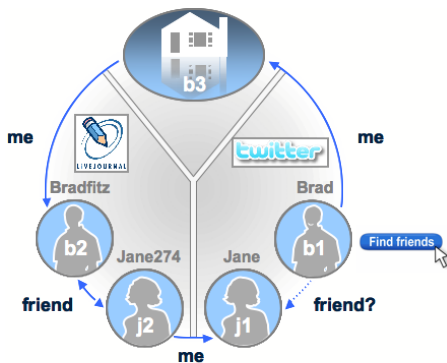
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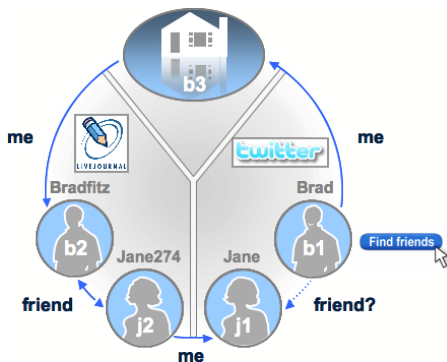


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- ▶ In Python, use NetworkX, cjson and a other standard scientific libraries to parse Google's SocialGraph data
- ▶ Using a "seed" user, we will build out a network



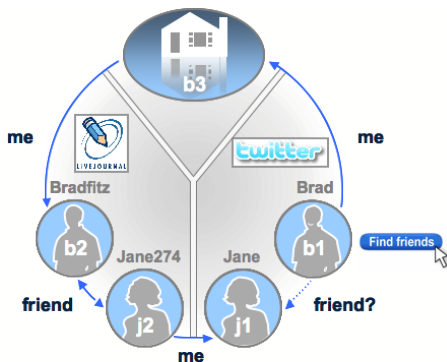
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- ▶ Through a process called "k-snowball searching"  
 $seed \rightarrow friend \rightarrow \dots \rightarrow friend_k$

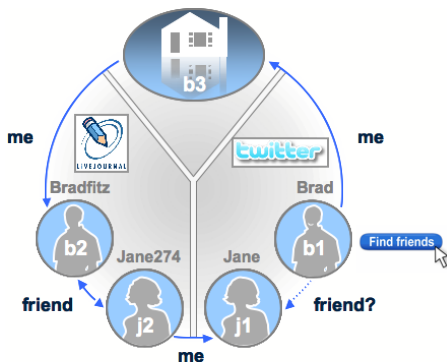
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  - ▶ Seed: imichaeldotorg.livejournal.com
  - ▶  $k = 3$

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  - ▶ Seed: imichaeldotorg.livejournal.com
  - ▶  $k = 3$
- ▶ Note the low value of  $k$

# The code, part 1

## Loading the libraries and setting things up

```
from cjson import *
from urllib import *
from networkx import *
from time import *
from scipy import array,unique
...
if __name__ == "__main__":
    seed_url='http://imichaeldotorg.livejournal.com'
    sg=get_sg(seed_url)
    net,newnodes=create_egonet(sg)
    info(net)
```

## Get the JSON from SocialGraph

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

Name:	['http://imichaeldotorg.livejournal.com/']
Type:	DiGraph
Number of nodes:	5
Number of edges:	5
Average in degree:	1.0
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## Get the JSON from SocialGraph

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def get_sg(seed_url):
    sgapi_url="http://socialgraph.apis.google.com/lookup?q="+seed_url+"&edo=1&edi=1&fme=1&pretty=0"
    try:
        furl=urlopen(sgapi_url)
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```

# Build egonet and snowball

## Creating the egonet

```
def create_egonet(s):
    try:
        raw=decode(s)
        G=DiGraph()
        pendants=[]
        n=raw['nodes']
        nk=n.keys()
        G.name=str(nk)
        pendants=[]
        for a in range(0,len(nk)):
            for b in range(0,len(nk)):
                if a!=b:
                    G.add_edge(nk[a],nk[b])
        for k in nk:
            ego=n[k]
            ego_out=ego['nodes_referenced']
            for o in ego_out:
                G.add_edge(k,o)
                pendants.append(o)
            ego_in=ego['nodes_referenced_by']
            for i in ego_in:
                G.add_edge(i,k)
                pendants.append(i)
        pendants=array(pendants, dtype=str)
        pendants.flatten()
        pendants=unique(pendants)
        return G, pendants
    except DecodeError:
        ...
    except KeyError:
```

## Rolling the snowball

```
def snowball_round(G, seeds, myspace=False):
    t0=time()
    if myspace:
        seeds=get_myspace_url(seeds)
    sb_data=[]
    for s in range(0,len(seeds)):
        s_sg=get_sg(seeds[s])
        new_ego, pen=create_egonet(s_sg)
        for p in pen:
            sb_data.append(p)
        if s<1:
            sb_net=compose(G, new_ego)
        else:
            sb_net=compose(new_ego, sb_net)
    del new_ego
    if s==round(len(seeds)*0.2):
        sb_net.name='20% complete'
        sb_net.info()
        print 'AT: ' +strtime('%m/%d/%Y, %H:%M:%S', gmtime())
        print ''
    ...
    # More time keeping, probably a MUCH better way to do this
    sb_data=array(sb_data)
    sb_data.flatten()
    sb_data=unique(sb_data)
    sb_net.info()
    return sb_net, sb_data
```

# Build the whole network

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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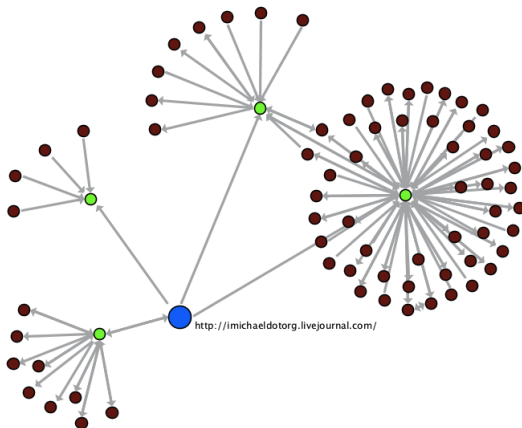
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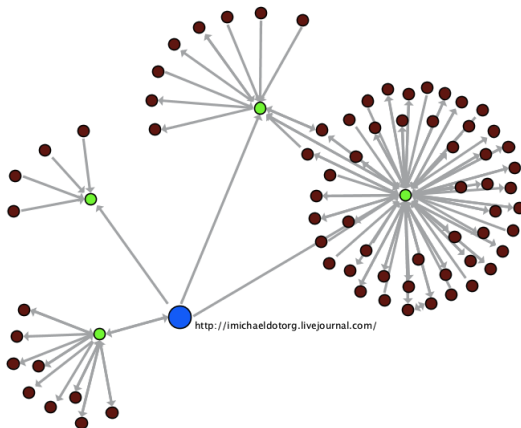
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- ▶ Our seed is abnormally isolated, with only four neighbors
- ▶ Large jump after first snowball
- ▶ Massive structural leap at  $k = 3$

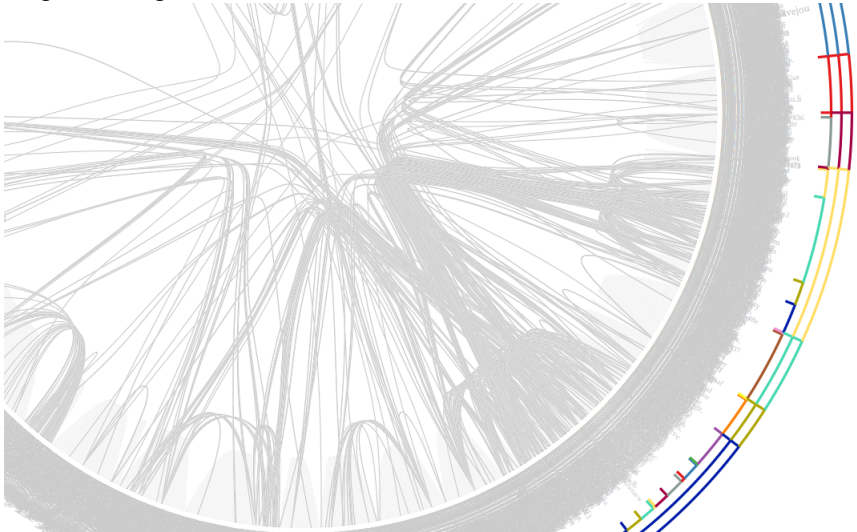


Loading data from multiple sources  
The Python dict  
Running basic centralities  
Getting things out of NetworkX

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## The full network

To get a feeling for the size of the full network...



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["orange","apple","banana"]
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Now, try creating a dict of your own

# Using dictionaries for network analysis

From the documentation...

## networkx.closeness centrality

```
closeness centrality(G, v=None, weighted_edges=False)
```

Compute closeness centrality for nodes.

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\_label: metric}

Let's look at an example:

## In-degree centrality of Hartford data

```
>>> in_cen=in_degree_centrality(hartford)
>>> in_cen
{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!

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## 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

### Find main component & symmetrize

```
# Many of the centrality metrics require undirected graphs, so we will symmetrize
>>> hartford_ud=hartford.to_undirected()
# The network also has many small components, but for
# this analysis we are interested in the largest
>>> hartford_mc=hartford_main=connected_component_subgraphs(hartford_ud)[0]
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Next, we will calculate multiple measures

### Computing multiple centralities

```
# Betweenness centrality
>>> bet_cen=betweenness_centrality(hartford_mc)
# Closeness centrality
>>> clo_cen=closeness_centrality(hartford_mc)
# Eigenvector centrality
>>> eig_cen=eigenvector_centrality(hartford_mc)
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```

Next, we will calculate multiple measures

### Computing multiple centralities

```
# Betweenness centrality
>>> bet_cen=betweenness centrality(hartford_mc)
# Closeness centrality
>>> clo_cen=closeness centrality(hartford_mc)
# Eigenvector centrality
>>> eig_cen=eigenvector centrality(hartford_mc)
```

## 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

### Find main component & symmetrize

```
# Many of the centrality metrics require undirected graphs, so we will symmetrize
>>> hartford_ud=hartford.to_undirected()
# The network also has many small components, but for
# this analysis we are interested in the largest
>>> hartford_mc=hartford_main=connected_component_subgraphs(hartford_ud)[0]
```

Next, we will calculate multiple measures

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```



## 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

### Function for finding most central actor

```
def highest centrality(cent_dict):  
    """Returns node key with largest value from  
    NX centrality dict"""  
    # Create ordered tuple of centrality data  
    cent_items=cent_dict.items()  
    # List comprehension!  
    cent_items=[(b,a) for (a,b) in cent_items]  
    # Sort in descending order  
    cent_items.sort()  
    cent_items.reverse()  
    return cent_items[0][1]
```

Now, just ask for the answer

### Finding Most central actors

```
>>> print("Actor "+str(highest centrality(bet_cen))+" has the highest Betweenness centrality")  
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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## Calculating basic community structure

Often in network analysis we are interested in estimating the cohesiveness of a network, or the communities that exists within the structure

### 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)

### Clustering

- ▶ 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  $v$ .

### Transitivity

- ▶ The fraction of all possible triangles which are in fact triangles. Or,  $Trans = 3 \left( \frac{T}{t} \right)$ , where  $T = \#$  of possible triangles and  $t = \#$  of actual triads

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

## Toy community detection example (not a good one)

### Calculating clustering coefficients

```
# 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()))
clus_counts=zip(map(lambda c: clus.values().count(c),unique_clus),unique_clus)
clus_counts.sort()
clus_counts.reverse()
# Create a subgraph from nodes with most frequent clustering coefficient
mode_clus_sg=subgraph(hartford_mc,[(a,b) for (a,b) in clus.items() if b==clus_counts[0][1]])
```

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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
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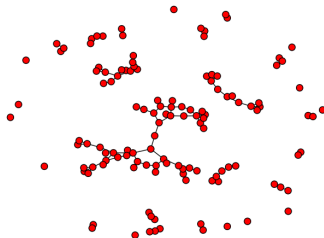
- Use the `with_labels` to return a dict keyed by node label
- The `zip` function takes two lists and returns a tuple
- More complex list comprehension with logic operator

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## Introduction to matplotlib

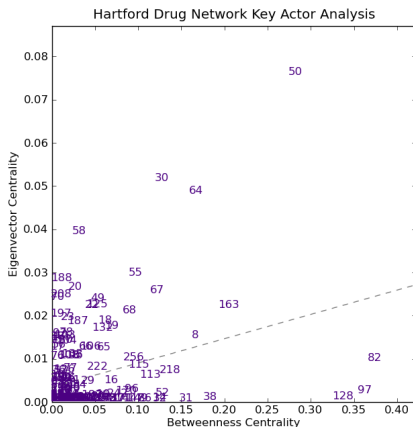
Recall Python's scientific computing trinity: NumPy, SciPy and matplotlib

- ▶ While NumPy and SciPy do most of the behind the scenes work, you will interact with matplotlib frequently for when doing network analysis

## Introduction to matplotlib

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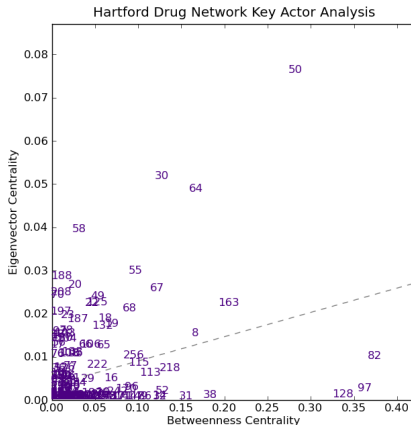
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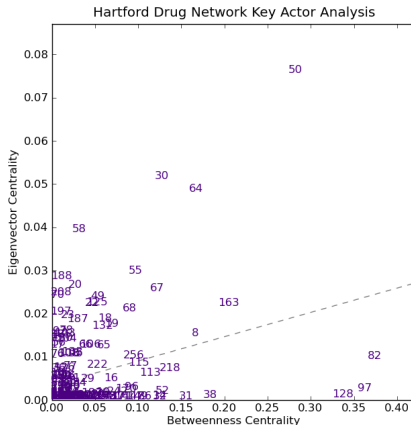


We will need to create a function that takes two centrality dict and generates this plot

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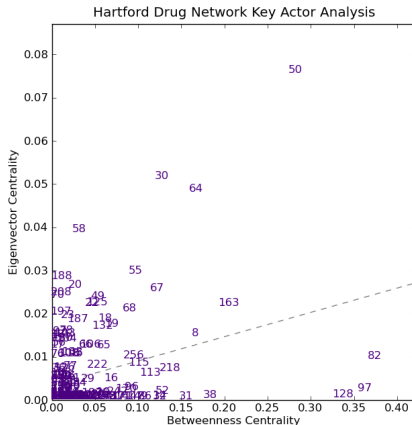
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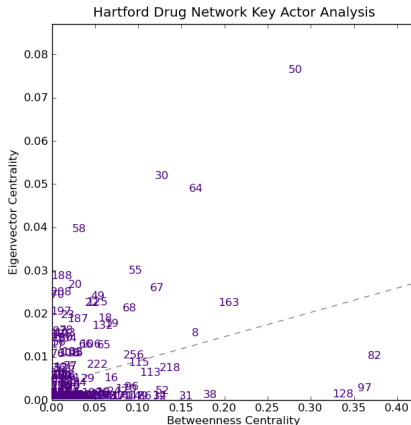
1. Create a matplotlib figure
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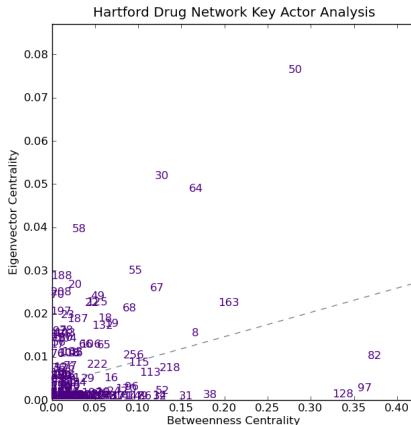
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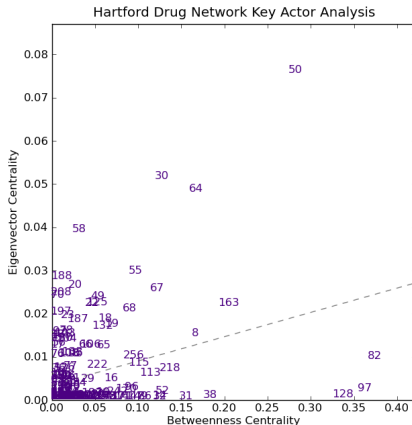
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We will need to create a function that takes two centrality dict and generates this plot

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

## Creating a key actor plot in matplotlib

### The centrality\_scatter function, part one

```
def centrality_scatter(met_dict1,met_dict2,path="",ylab="",xlab="",title="",reg=False):  
    # Create figure and drawing axis  
    fig=P.figure(figsize=(7,7))  
    ax1=fig.add_subplot(111)  
    # Create items so actors can be sorted properly  
    met_items1=met_dict1.items()  
    met_items2=met_dict2.items()  
    met_items1.sort()  
    met_items2.sort()  
    # Grab data  
    xdata=[(b) for (a,b) in met_items1]  
    ydata=[(b) for (a,b) in met_items2]  
    # Add each actor to the plot by ID  
    for p in xrange(len(met_items1)):  
        ax1.text(x=xdata[p],y=ydata[p],s=str(met_items1[p][0]),color="indigo")
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- ▶ Create a canvas to draw on
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- ▶ Create a canvas to draw on
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- ▶ Add points to plot as node labels

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def centrality_scatter(met_dict1,met_dict2,path="",ylab="",xlab="",title="",reg=False):
    ...
    # If adding a best fit line, we will use NumPy to calculate the points.
    if reg:
        # Function returns y-intercept and slope. So, we create a function to
        # draw LOBF from this data
        slope,yint=polyfit(xdata,ydata,1)
        xline=P.xticks()[0]
        yline=map(lambda x: slope*x+yint,xline)
        # Add line
        ax1.plot(xline,yline,ls='--',color='grey')
    # Set new x- and y-axis limits to data
    P.xlim((0.0,max(xdata)+(.15*max(xdata)))) # Give a little buffer
    P.ylim((0.0,max(ydata)+(.15*max(ydata))))
    # Add labels
    ax1.set_title(title)
    ax1.set_xlabel(xlab)
    ax1.set_ylabel(ylab)
    # Save figure
    P.savefig(path,dpi=100)
```



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- ▶ Add a best fit line
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    # Save figure
    P.savefig(path,dpi=100)
```

- ▶ 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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3. Depending on your needs you may prefer one, the other or both

## 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

- ▶ Suite of tools lacks your specific need
- ▶ Require alternate visualization
- ▶ Storage for later analysis

In most cases this will entail either exporting the raw network data, or metrics from some network analysis

1. NetworkX can write out network data in as many formats as it can read them, and the process is equally straightforward
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











## Saving network data in different formats

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

### NX syntax for loading a file

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

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

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

```
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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- ▶ 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 DictWriter

### Exporting centrality data to CSV

```
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"]
    col_headers.extend(data_dict.keys())
    # Create CSV writer and write column headers
    writer=csv.DictWriter(open(path,"w"),fieldnames=col_headers)
    writer.writerow(dict((h,h) for h in col_headers))
    # Write each row of data
    for j in data_dict[col_headers[1]].keys():
        # Create a new dict for each row
        row=dict.fromkeys(col_headers)
        row["Actor"]=j
        for k in data_dict.keys():
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## 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

◇	A	B	C	D
1	Actor	Closeness	Betweenness	Eigenvector
2	1	0.12467532	0.0072576	0.00025176
3	2	0.12475634	0.01767427	0.00025964
4	3	0.12565445	0.05687441	0.00023185
5	4	0.10223642	0.03108639	1.44E-05
6	5	0.1443609	0	0.00313152
7	6	0.09943035	0.01041667	1.49E-07
8	7	0.11340815	0.04362093	6.78E-05
9	8	0.20512821	0.16354003	0.01471888
10	9	0.11267606	0.00741624	0.0001101
11	10	0.13983977	0.05258239	0.00095456
12	11	0.1703638	0.01250999	0.0032333
13	13	0.13892909	0	1.79E-05
14	14	0.17219731	0.11848775	0.00029737
15	15	0.13521127	0.00079897	2.11E-05
16	16	0.15907208	0.06203647	0.00432838