

# Networks Structure and Dynamics

## 2. Graph algorithms

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## Outline

- 1 Definitions and metrics
  - Reminder
  - Distributions
    - Benefit
    - Cumulative distributions
- 2 Algorithms
  - Connected components
  - Local density
  - Distances and diameter
  - Centralities

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## Definitions and notations

A graph  $G = (V, E)$  is a couple of sets.

- $V$  is the set of *nodes*
- $E \subseteq (V \times V)$  is the set of *edges*

We denote :

- $n = |V|$  the number of nodes
- $m = |E|$  the number of edges

$u$  and  $v$  are **neighbors** if there is an edge between them.

**Degree:** ( $k_i$  or  $d^\circ(i)$ ) number of neighbors of  $i$

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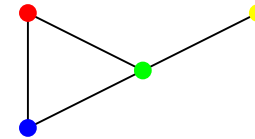
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## Average degree, density

- Average degree of a graph,  $d^\circ(G) = \frac{\sum_v d^\circ(v)}{n}$
- density of a graph,  $\delta = \frac{2m}{n(n-1)}$

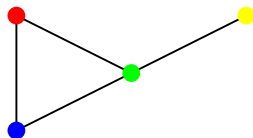


degrees : 2, 2, 3, 1 ; average degree 2

$$n = 4, m = 4, \delta = \frac{8}{12} = 0.66..$$

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## Connectedness

**Path** from  $u$  to  $v$ : sequence of edges

$(u, v_1), (v_1, v_2), \dots, (v_{\alpha-1}, v)$

**Length**: number of edges in the path (here  $\alpha$ )

**Connected component**: maximal set of nodes such that  $\exists$  a path between any pair of nodes

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## Distributions

Distribution : synthetic way to represent a **sequence of values**.

→ how many times a value occurs in the sequence?

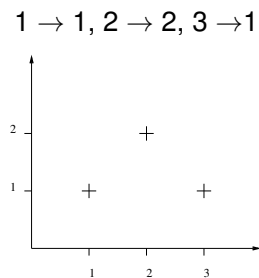
## Distributions

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→ how many times a value occurs in the sequence?

Example/reminder of the degree distribution:

4 nodes, degrees : **2 2 3 1**



## Distribution characteristics

One benefit: characterize **qualitatively** a sequence of values.

### Power law

- degree distribution follows  $N_k \sim k^{-\alpha}$
- line in échelle log-log scale

**heterogeneous** distribution: non-homogeneous (various types of behaviors), in practice often **close** to a power law

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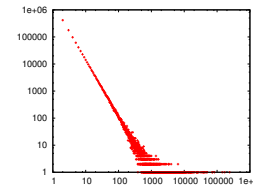
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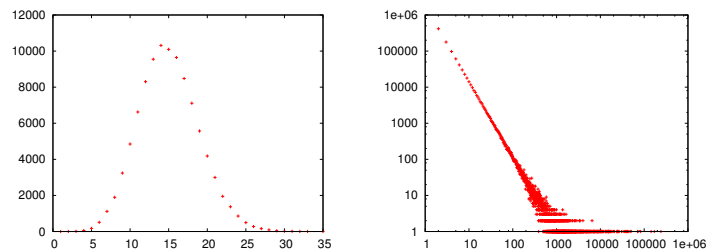
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## Heterogeneous vs homogeneous distributions



### Homogeneous

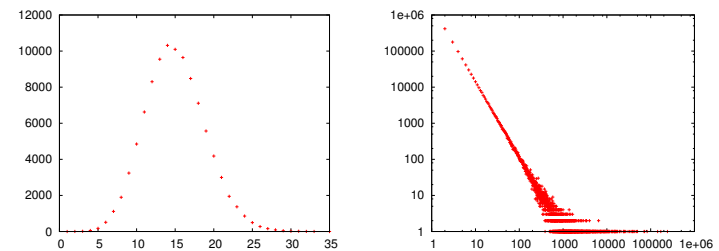
Idea of normality (and of **exceptions**)

### Heterogeneous

Any kind of behaviours → no notion of normality

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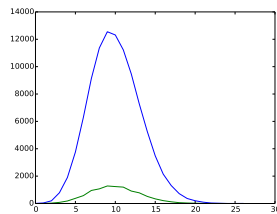
**Two choices:**

- $N_k$ : number of occurrences of value  $k$  in the sequence
- $p_k$ : **proportion** of the value  $k$  in the sequence  
→ **Normalized** distribution

$$p_k = \frac{N_k}{n}$$

Just a change of the value on the Y-axis.

Allow to compare graphs with different sizes:



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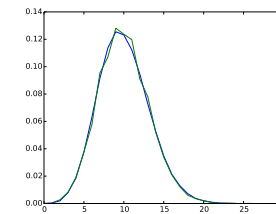
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**Example of homogeneous distribution: the normal law****Normal law (Gaussian law)**

$$P(x) \sim \exp\left(-\frac{(x - \langle x \rangle)^2}{2\sigma^2}\right)$$

Normal law often found in “nature”, good model for:

- asymptotic limit of binomial law: repeated coin flipping experiments (heads:+1, tails:-1), Galton board, ...



- complex phenomena: human height distribution etc.

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**Characterizing a distribution****First and second moments**

- $\langle k \rangle = \sum k p_k$ : first moment (**mean**)
- $\langle k^2 \rangle = \sum k^2 p_k$ : second raw moment  
we rather use the second central moment or **variance** ( $\sigma^2$ ):

$$\sum (k - \langle k \rangle)^2 p_k = \langle k^2 \rangle - \langle k \rangle^2$$

and higher order moments  
(skewness - *dissymétrie*, kurtosis - *aplatissement*, ...)

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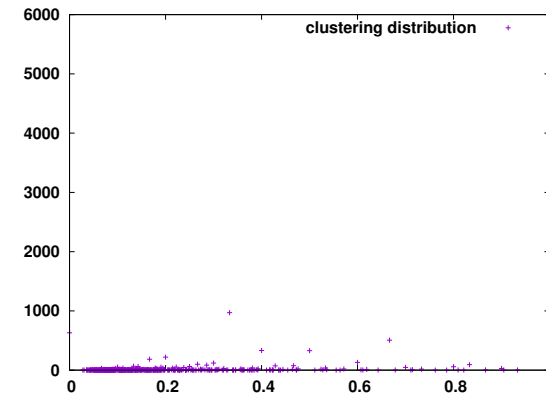
## Notion of cumulative distributions

- **Distribution of  $k$ :**  
 $N_k$ : number of occurrences **equal to  $k$**
- **Cumulative distribution of  $k$ :**  
 $C_k$ : number of occurrences **lower or equal to  $k$**
- **Inverse cumulative distribution of  $k$ :**  
 $IC_k$ : number of occurrences **greater or equal to  $k$**

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## Relevance of cumulative distributions

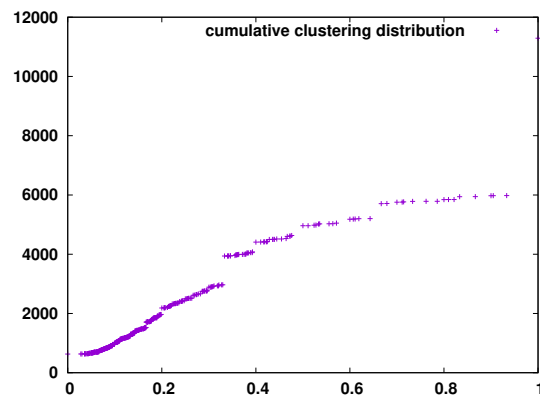
Illustration: clustering distribution of a coauthoring network



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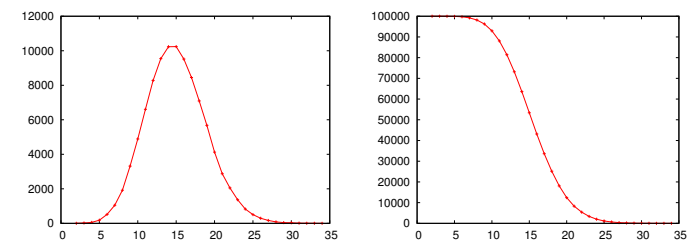


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$N_k$  and  $IC_k$  for a **homogeneous** distribution:



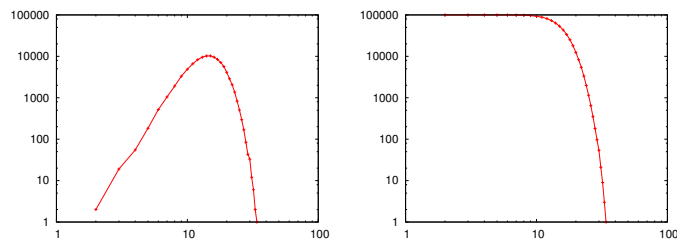
linear scale

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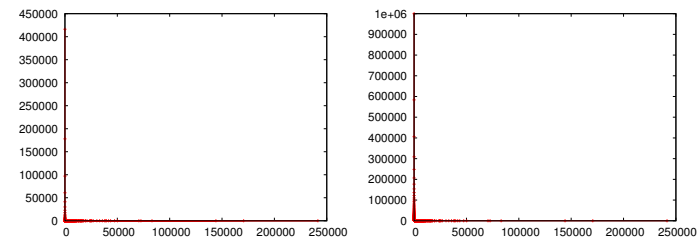
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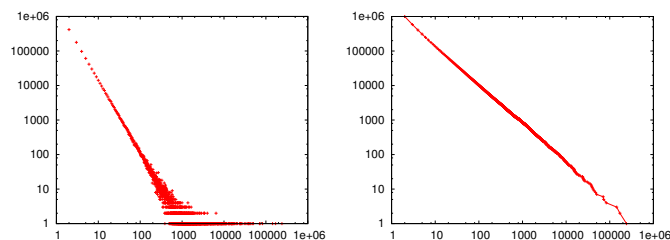
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## Homogeneous and heterogeneous

- can be distinguished on both normal and cumulative distributions

## Ex: power-law

- $N_k \sim k^{-\alpha} \Rightarrow C_k \sim k^{-\alpha+1}$   
Remark: same idea as  $\int x^{-\alpha} dx \sim x^{-\alpha+1}$

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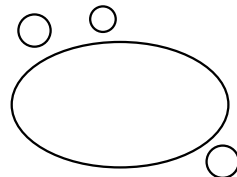
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## Connectedness (reminder)

### For complex networks

In general, **giant** component  
→ contains most nodes



**Q** : How to identify the giant component? How to count the connected components?

## Breadth First Search algorithm (BFS)

*Parcours en largeur*

**Algorithm 1:** Breadth First Search of a graph  $G$  from node  $s$ .

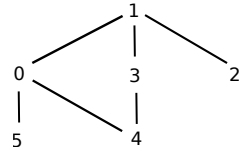
```

begin
  F ← CreateEmptyQueue()
  Enqueue(F,s)
  Mark(s)
  while F not empty do
    u ← DequeueFirstElement(F)
    Display u
    for v neighbor of u in G do
      if Unmarked(v) then
        Enqueue(F,v)
        Mark(v)
      end
    end
  end
end
end

```



## Example



- 1 Apply this algorithm to the previous example from node 3.
- 2 How to modify it so that it returns the tree of shortest paths from a node? Draw the corresponding tree.

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## Breadth First Search algorithm (BFS)

Properties of a BFS:

- From a node : we detect **its connected component**  
→ 1 BFS per component
- Complexity:  $\mathcal{O}(m)$
- With the alternative with parentage memorization:  
**spanning tree** (fr: *arbre couvrant*) of shortest paths

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## Going back to local density

Several means to capture this idea, for example:

- **clustering coefficient**:  $cc(G) = \frac{\sum_v \frac{\Delta(v)}{\Lambda(v)}}{n}$
- **transitive ratio**:  $tr(G) = \frac{3\Delta(G)}{\Lambda(G)}$

In other words:

- clustering coefficient: compute a value for each node (with degree  $\geq 2$ ) then average
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## Transitive ratio vs Clustering coefficient

We have  $n \in \mathbb{N}$ , let's  $G_n$  be the graphs of  $2n + 1$  nodes and  $3n$  edges such that:

- a unique node  $n_0$  is connected to all other nodes in the network
- all other nodes have degree 2

Exercise :

- 1 Draw the cases of  $G_3, G_4$ .
- 2 Compute the local density coefficients for  $G_4$ .
- 3 How do these coefficients evolve when  $n$  goes to  $\infty$ ?
- 4 Deduce how to interpret these coefficients in terms of probability.

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## Computing the number of triangles

Both coefficients rely on the number of triangles. How to enumerate the number of triangles that node  $n$  belongs to?

**Naive answer:** For any pair of neighbors  $(u_1, u_2)$  of  $v$ , test if  $(u_1, u_2)$  exists.

Questions :

- 1 Complexity of the algorithm for node  $v$ ?
- 2 In which case is it expensive?
- 3 Is it a problem for the networks we are working on?

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- 3 Is it a problem for the networks we are working on?  
yes, because heterogeneous degree distribution

## (Faster) computation of the number of triangles

**Other point view:** Considering edge  $(u, v)$ , how many triangles contain it?

**Idea:** There is a triangle if there exists a node  $w$  neighbor of  $u$  and  $v$ .

**Solution:** We need to compute the size of the intersection of the neighborhood of  $u$  and  $v$ . **Efficient if lists of neighbors are ordered.**

**Algorithm 2:** Size of the intersection of two ordered lists  $U$  and  $V$ , of size  $d^o(u)$  and  $d^o(v)$

```

i_u = 0
i_v = 0
nb = 0
while (i_u < d^o(u)) and (i_v < d^o(v)) do
  if U[i_u] < V[i_v] then
    i_u++
  else
    if U[i_u] > V[i_v] then
      i_v++
    else
      nb++ // triangle found
      i_u++
      i_v++
    end
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    end
  end
end
end

```

## (Even faster) computation of the number of triangles

### Remarks:

- Algorithm slow if  $u$  or  $v$  have high degree
- Triangle  $(u, v, w)$  may be detected from the 3 list intersections:  
 $U$  with  $V$ ,  $U$  with  $W$ ,  $V$  with  $W$

**Idea:** Reduce as much as possible the size of the lists involved in the intersection computation:

- sort nodes per decreasing degree and re-index the graph
- consider edges  $(u, v)$  such that  $u < v$
- we look for nodes  $w$  neighbor of  $u$  and  $v$  only if  $w < u$  (and so  $w < v$ ).

## (Fast) computation of the number of triangles

**Algorithm 4:** Compute the number of triangles

**Result:** Table  $tr$  contains the # of triangles of each node

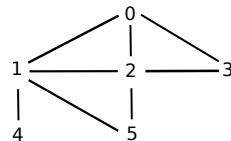
$tr$ : table of size  $n$  initialized at 0;

```

for  $u = 0; u < n; u++$  do
  for  $i = 0; i < d^{\circ}(u); i++$  do
     $v = i$ -th neighbor of  $u$ 
    if  $u < v$  then
       $U =$  list of neighbors of  $u$ 
       $V =$  list of neighbors of  $v$ 
       $i_u = 0, i_v = 0$ 
      while  $(i_u < d^{\circ}(u))$  and  $(i_v < d^{\circ}(v))$ 
        and  $(U[i_u] < u)$  and  $(V[i_v] < u)$  do
          if  $U[i_u] < V[i_v]$  then
             $i_u++$ ;
          else
            if  $U[i_u] > V[i_v]$  then
               $i_v++$ ;
            else
               $tr[u]++$ ;
               $tr[v]++$ ;
               $tr[U[i_u]]++$ ;
               $i_u++$ ;
               $i_v++$ ;
            end
          end
        end
      end
    end
  end
end
end
end
    
```

## Exercise

Apply the previous algorithm to the following graph:



## Higher order cliques

### Clique

Triangles are 3-nodes cliques. A clique is a complete subgraph.

- subgraph:** graph obtained considering a subset of nodes and the edges between these nodes (fr: *sous-graphe*)
- complete:** any node is connected to all others (fr: *complet*)

### Maximal cliques

Decomposition of a graph into its maximal cliques

Obtaining the list of all maximal cliques is known to be computationally hard  
→ we favor search with fixed size

## Higher order cliques

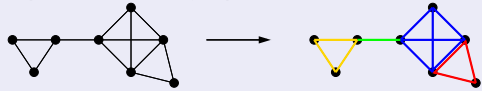
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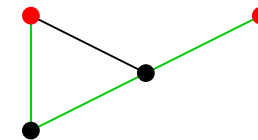


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## Distances and diameter (reminder)

**path** from  $u$  to  $v$  = sequence of edges  $u\dots v$



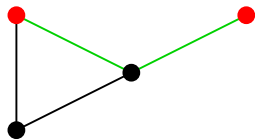
one path of length 3

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## Distances and diameter (reminder)

**path** from  $u$  to  $v$  = sequence of edges  $u\dots v$

**distance**  $d(u, v)$  = length of *one* shortest path



a shortest path; length 2  $\Rightarrow$  distance = 2

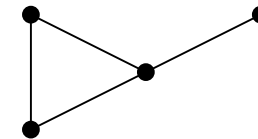
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**average distance** = average of the distances between all pairs of nodes



Average distance =  $8/6$

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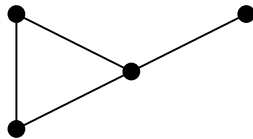
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**average distance** = average of the distances between all pairs of nodes

**diameter**  $\Delta$  = highest distance (on any set of pairs)



diameter = 2

## Distances computation

Distance from a node to all others:  
modified **breadth first search** → Complexity:  $\mathcal{O}(m)$

### Average distance, diameter

Need all distances →  $\mathcal{O}(nm)$   
possible to **approximate** or to **give bounds** on the diameter

### Example: approximation of average distance

Distance is a homogeneous property  
Possible to use a **random sampling**  
!: not if **heterogeneous property**

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!: not if **heterogeneous property**

## Distances computation

Distance from a node to all others:  
modified **breadth first search** → Complexity:  $\mathcal{O}(m)$

### Average distance, diameter

Need all distances →  $\mathcal{O}(nm)$   
possible to **approximate** or to **give bounds** on the diameter

### Example: approximation of average distance

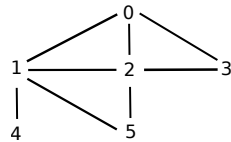
Distance is a homogeneous property  
Possible to use a **random sampling**  
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## Exercise

Whether it be for the average distance or the diameter, one needs to be able to compute the distance between two nodes.

Exercise :

- Propose an algorithm, based on the BFS principle, enabling to compute the distance between one given node towards **all** nodes of the graph.
- Apply the algorithm on the following graph (taking 5 as the initial node):



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## Modified BFS

**Algorithm 5:** Distance from node  $s$  in graph  $G$ .

```

begin
  F ← CreateEmptyQueue()
  Enqueue(F,s)
  ∀v Dist(v) initialized at -1
  Dist(s) ← 0
  while F not empty do
    u ← DequeueFirstElement(F)
    Display u
    for v neighbor of u in G do
      if Dist(v) = -1 then
        Enqueue(F,v)
        Dist(v) ← Dist(u) + 1
      end
    end
  end
end
end
  
```

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## Approximations

Approximation: given a property  $P$ , how to estimate this property on a given graph.

Examples: average degree, average distance, diameter, ...

### One possible approach (sampling)

- Pick a node  $v$  of  $G$  at random
- Estimate the property for  $v$
- Go back to step 1 while the estimation is not good enough

Questions:

- How to express the notion of "good enough"?
- How to know if this approach provides a good approximation or not?

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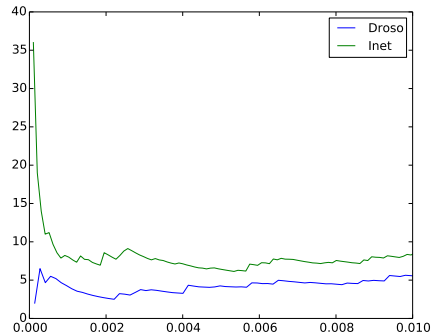
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## Average degree

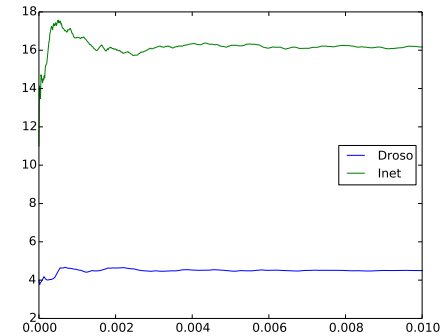
Application of the former method to the average degree:  
X-axis : fraction of measured values



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## Average distance

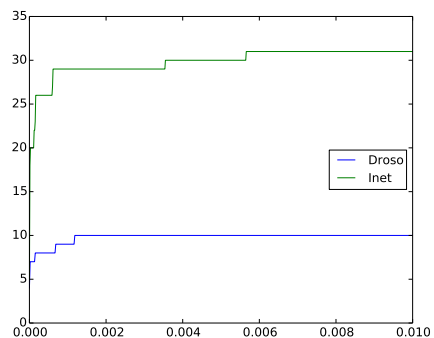
Application of the former method to the average distance:  
X-axis : fraction of measured values



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## Diameter

Application of the former method to the average degree:  
X-axis : fraction of measured values



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## Approximations

Quality of the approximation: depends on the nature of the property.

### Other possible approach:

- Compute (lower and upper) bounds of the property
- Rely on the property to drive the computations

Example: For every node  $v$ , let  $max_v$  be the greatest distance from  $v$  to a node of  $G$ . Then the diameter  $D$  of  $G$  is such that:

$$max_v \leq D \leq 2max_v$$

Exercise : explain why.

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## Centrality measures

**Question: how to quantify the importance of a node in a network?**

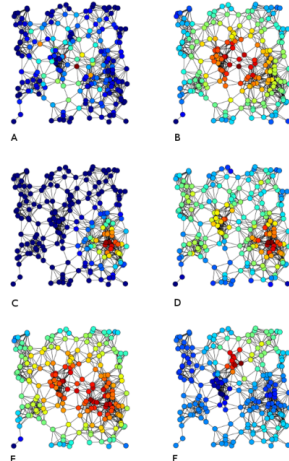
**Centrality measures:** “key” nodes of the network.

- Degree centrality (degree)
- Katz centrality
- Closeness (and harmonic) centralities
- Eigenvector and Pagerank centralities
- **Betweenness centrality**
- etc.

## Centrality measures

**Different metrics highlight different properties**

- A. betweenness
- B. closeness
- C. eigenvector
- D. degree
- E. harmonic
- F. Katz



## About betweenness centrality

Let  $G$  be a graph and  $v$ ,  $s$  and  $t$  be nodes of  $G$ .  
We call *betweenness centrality* the value:

$$BC(v) = \sum_{s \neq t \neq v} \frac{\sigma_{st}(v)}{\sigma_{st}}$$

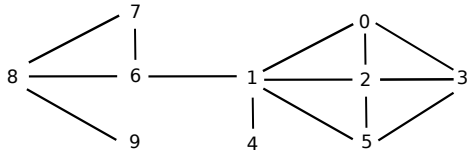
with:

- $\sigma_{st}$  : number of shortest paths between  $s$  and  $t$
- $\sigma_{st}(v)$  : number of shortest paths between  $s$  and  $t$  going through  $v$

## Meaning

Capture the importance of a node  $v$  in a graph as a **relay** for diffusing information:

- propagation of signals/messages/virus
- connectivity of the network



Exercise : compute  $BC(0)$ ,  $BC(1)$  and  $BC(4)$ .

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## Link betweenness centrality

In the former example:

- 1 has a high centrality
- but all links starting from 1 don't have the same importance

There exists a similar definition to compute the importance of the links:

$$BC(u, v) = \sum_{s \neq t \neq u \neq v} \frac{\sigma_{st}(u, v)}{\sigma_{st}}$$

with:

- $\sigma_{st}$  : nb of shortest paths between  $s$  and  $t$
- $\sigma_{st}(u, v)$  : nb of shortest paths between  $s$  and  $t$  using  $(u, v)$

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## Relation to communities

What do you think about the relation between link betweenness and the communities of a network?

Can you think of an algorithm that would use betweenness to infer communities?

- 1 (Re)compute the centrality of all links
- 2 Delete the link with highest centrality
- 3 Test if an isolated connected component appears
- 4 Repeat steps 1 – 3 until there are no more links.

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## About sparse matrix product

Many measurement on graphs demand matrix product  
ex: *Katz, PageRank, eigenvector centralities*

### Matrix product

Given a matrix  $M (n \times n)$  and a vector  $V$   
we compute the vector  $W (n \times 1)$  such that:

$$W = M \times V$$

where  $\forall i \in [1, n], W_i = \sum_{j=1}^n M_{ij} \times V_j$

Note that for a  $n \times n$  sparse matrix  $M$  with  $m$  nonzero values,  
we take into account only the nonzero values  
 $\Rightarrow$  compute  $W$  is in  $\mathcal{O}(m + n)$  and not in  $\mathcal{O}(n^2)$ .

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## Classic Version

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**Algorithm 6:** Classic sparse Matrix/Vector multiplication

---

**input:**  $M, V$ ;

**for**  $i$  from 1 to  $n$  **do**

$W[i] \leftarrow 0$

**for each**  $j$  such that  $M[i][j] \neq 0$  **do**

$W[i] = W[i] + M[i][j] \times V[j]$

**end**

**end**

**return**  $W$ ;

---

Should be implemented with “the adjacency list” graph data structure.

**to be continued in the course on PageRank**