

# **Graphs**Part: Structure



Algorithms and Data Structures 2, 340300 Lecture – 2023W Univ.-Prof. Dr. Alois Ferscha, teaching@pervasive.jku.at

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#### **Definition**

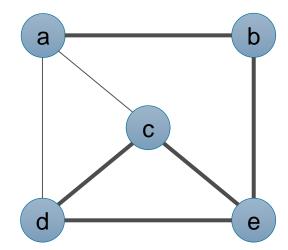
A graph G = (V, E) consists of

*V* ... a quantity of vertices/nodes and*E* ... a quantity of edges/arcs.

An edge e = (u, v) is a pair of vertices.

#### **Example:**

$$V = \{ a, b, c, d, e \}$$
  
 $E = \{ (a,b), (a,c), (a,d), (b,e), (c,d), (c,e), (d,e) \}$ 





#### **Definitions of Terms**

Two vertices are adjacent to each other, if they are connected by an edge.

■ Example: a and c are adjacent to each other

An edge connecting two adjacent vertices is called incident to these vertices.

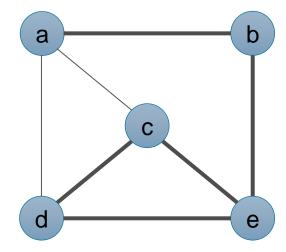
The **degree** of a vertex **deg(v)** is defined as the number of vertices adjacent to it.

• We have: 
$$\sum_{v \in V} \deg(v) = 2 * (\#edges)$$

• Example: deg(c) = 3

A **path** is a **sequence** of **vertices**  $(v_1, v_2, ..., v_k)$  in which successive vertices  $v_i$  and  $v_{i+1}$  are adjacent to each other.

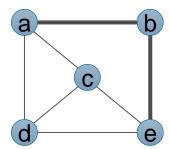
Example: path (a, b, e, d, c, e)



#### **Definitions of Terms**

In a **simple path** no vertex occurs more than once.

Example: a, b, e



A cyclic path (short: cycle) is a simple path with the exception, that the **first vertex** in the path is **identical** to the **last vertex** in the path.

Example: a, b, e, c, a



A **subgraph** is a subset of vertices and edges of a graph, which in turn form a graph.



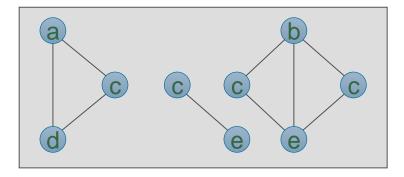
#### **Definitions of Terms**

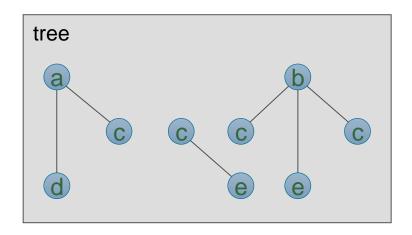
A connected component in a graph is a maximum connected subgraph.

Example:Graph with 3 connected components

A tree is a connected graph without cycles.

A forest is a set of trees.







### Connectivity

In a **complete graph each pair** of vertices is **adjacent** to each other.

Let

**n** ... number of vertices and

*m* ... number of **edges** 

• in a complete graph we have:  $m = \frac{1}{2} \sum_{v \in V} \deg(v) = \frac{1}{2} \sum_{v \in V} (n-1) = \frac{n(n-1)}{2}$ 

• in a **non-complete** graph we have:  $m < \frac{n(n-1)}{2}$ 

• in a tree we have: m = n - 1

If m < n-1, then the graph is **not connected** (consists of more than one connected component).

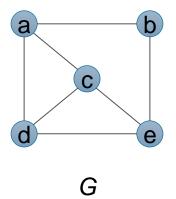
### **Spanning Tree**

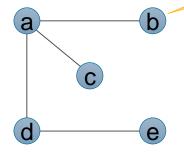
A **spanning tree ST** in a graph *G* is a subgraph of *G* for which we have:

• ST is a tree

• ST contains all vertices of G

not unambiguous!





Spanning tree of G

If a **single edge** is **removed**, the graph is **no longer connected**.



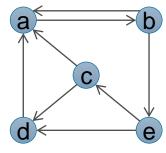
### **Directed Graph**

An edge (v, w) is defined as **directed**, if it leads v to w, but not vice versa. Then (v, w) is an ordered pair.

Illustration: as arrow



A graph is directed (digraph), if it contains directed edges.



A directed graph without cycles is called a **directed acyclic graph** (**DAG**).

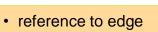
A digraph is strongly connected, if each vertex can be reached from any other vertex.

A **strongly connected component** in a digraph is a subgraph, in which each vertex can be reached from any other vertex.

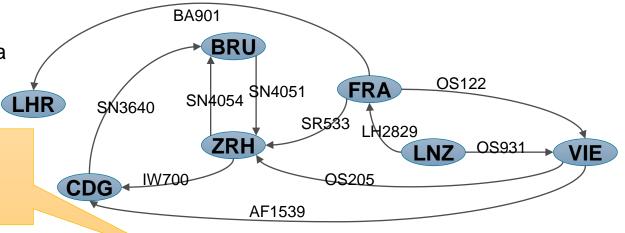


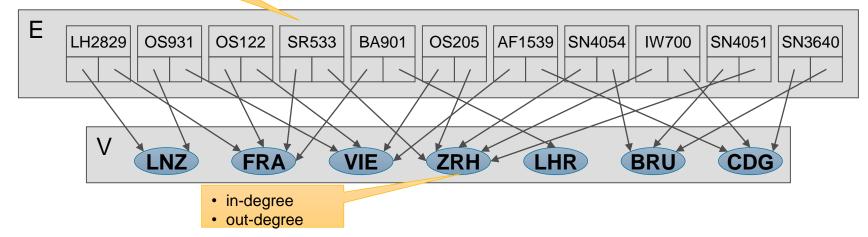
#### **Edge list**

**Edges** are stored **in a list**, from each edge a **reference leads** to the end vertices stored in a separate data structure.



- (un)directedreference to vertex
- position in container

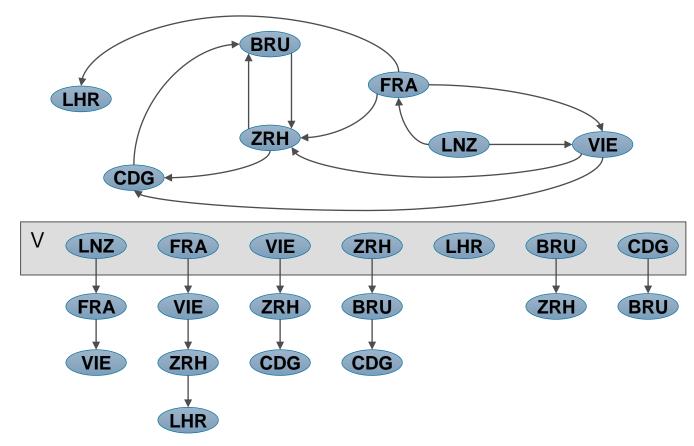






#### **Adjacency list (traditional)**

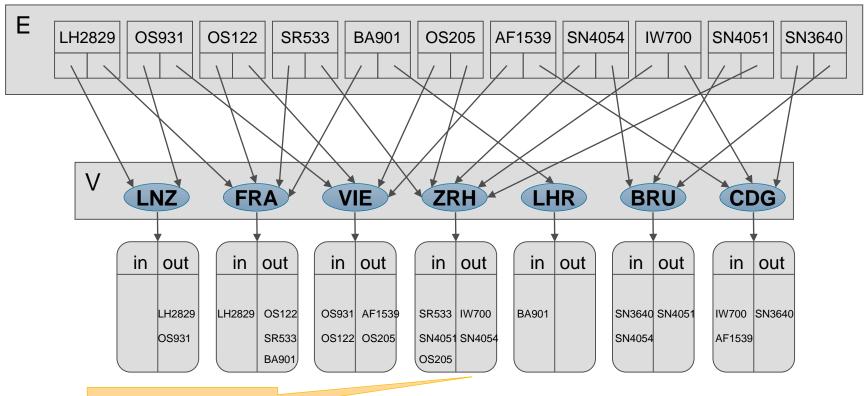
- For each vertex its **adjacent vertices** are stored in a **list**.
- All vertices with adjacency lists are again stored in a list.





#### Adjacency list (modern)

■ Extension of the edge list by a list of the incident edges for each vertex



- Incidency container
- (in/out count)



#### **Adjacency matrix (traditional)**

■ Matrix *M* where *M[i,j]* is **true** if an **edge** from **vertex** *i* to **vertex** *j* exists, or **false** otherwise.

from	LNZ	FRA	VIE	ZRH	LHR	BRU	CDG
LNZ	F	Т	Т	F	F	F	F
FRA	F	F	Т	Т	Т	F	F
VIE	F	F	F	Т	F	F	Т
ZRH	F	F	F	F	F	Т	Т
LHR	F	F	F	F	F	F	F
BRU	F	F	F	Т	F	F	F
CDG	F	F	F	F	F	Т	F



#### **Adjacency matrix (modern)**

Matrix elements store a reference to an edge object

from	LNZ	FRA	VIE	ZRH	LHR	BRU	CDG
LNZ	F	LH2829	OS931	F	F	F	F
FRA	F	F	OS122	SR533	BA901	F	F
VIE	F	F	F	OS205	F	F	AF1539
ZRH	F	F	F	F	F	SN4054	IW700
LHR	F	F	F	F	F	F	F
BRU	F	F	F	SN4051	F	F	F
CDG	F	F	F	F	F	SN3640	F



# **Complexity Comparison**

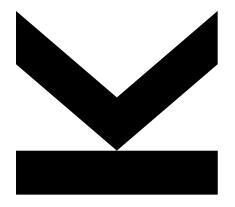
#### Assumption:

- Edges, vertices as doubly linked list
- Container with position information

	Complexity				
	Edge list	Adjacency list	Adjacency matrix		
Space	O(m+n)	O(m+n)	O(n²)		
Operations:					
size, isEmpty	O(1)	O(1)	O(1)		
vertices	O(n)	O(n)	O(n)		
edges	O(m)	O(m)	O(m)		
endVertices, opposite, isDirected	O(1)	O(1)	O(1)		
incidentEdges	O(m)	O(deg(v))	O(n)		
areAdjacent	O(m)	O(min(deg(u),deg(v)))	O(1)		
insertVertex	O(1)	O(1)	O(n²)		
removeVertex	O(m)	O(deg(v))	O(n²)		
<pre>insertEdge, insertDirectedEdge, removeEdge</pre>	O(1)	O(1)	O(1)		



# **Graph Structure Analysis**



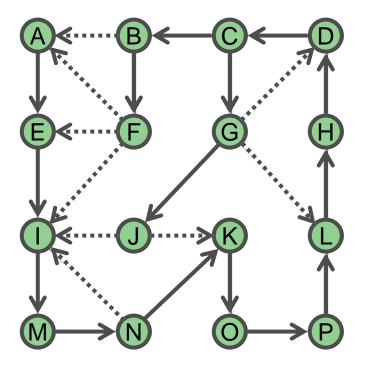


### **Graph Traversal – Depth-First Search (DFS)**

```
Algorithm DFS(v);
    Input: A vertex v in a graph
    Output: A labeling of edges as "discovery" and "back" edges
    for each edge e incident to v do
        if edge e is unexplored then
             let w be the other endpoint of e
            if vertex w is unexplored then
                 label e as a discovery edge
                 recursively call DFS(w)
            else
                 label e as a back edge
```



# **DFS :: Example**



discovery edge

back edge



### **Properties of DFS**

For an **undirected graph G** in which a **DFS starting** with **vertex s** is executed, we have:

- The traversal **visits all vertices** in the connected component (=maximum connected subgraph), which contains s.
- The set of discovery edges form a spanning tree for this connected component.

#### **Complexity of DFS**

If  $n_s$  is the number of **vertices** in a connected component within s and if  $m_s$  is the number of **edges** in the connected component of s,

then the complexity of DFS is  $O(n_s + m_s)$  on the assumptions:

- Graph is stored so that access to the vertices and edges is O(1)
- Marking and testing of the edges is O(1)
- There is a mechanism that systematically searches the edges of a node without looking at an edge more than once.



#### **DFS :: Finding a Path**

Each **vertex** is labeled initially as **UNEXPLORED** and can be relabeled as **VISITED**.

Each **edge** is labeled initially as **UNEXPLORED** and can be relabeled as **DISCOVERY** or **BACK**.

A **stack S** is used to keep track of the **path between** the **start vertex** and the **current vertex**.

When the destination vertex z is reached, the stack content is returned as the path between v and z.

Based on Goodrich et al., Data Structures & Algorithms in Python

```
Input: A graph G and the vertices v (start) and z
(destination) in graph G
Output: Path as a sequence of vertices
setLabel(v, VISITED)
S.push(v)
if v = z
     return S. elements() and terminate algorithm
for all e ∈ G.incidentEdges(v)
     if getLabel(e) = UNEXPLORED
           w \leftarrow opposite(v,e)
           if getLabel(w) = UNEXPLORED
                setLabel(e, DISCOVERY)
                S.push(e)
                pathDFS(G, w, z)
                S.pop(e)
           else
                setLabel(e, BACK)
S.pop(v)
```

**Algorithm** pathDFS(G,v,z);

### **DFS :: Finding a Cycle**

Each **vertex** is labeled initially as **UNEXPLORED** and can be relabeled as **VISITED**.

Each **edge** is labeled initially as **UNEXPLORED** and can be relabeled as **DISCOVERY** or **BACK**.

A **stack S** is used to keep track of the **path between** the **start vertex** and the **current vertex**.

A **cycle** is detected when a back edge to an already visited vertex is recognized, which is not a parent.

Based on Goodrich et al., Data Structures & Algorithms in Python

```
Algorithm cycleDFS(G,v);
     Input: A vertex v (start) in a graph G
     Output: Path of in the graph representing the cycle
     setLabel(v, VISITED)
     S.push(v)
     for all e \in G.incidentEdges(v)
           if getLabel(e) = UNEXPLORED
                 \mathbf{w} \leftarrow \text{opposite}(v,e)
                 S.push(e)
                 if getLabel(w) = UNEXPLORED
                       setLabel(e, DISCOVERY)
                       cycleDFS(G, w)
                       S.pop(e)
                 else
                       T ← new empty stack
                       repeat
                             o \leftarrow S.pop()
                             T.push(o)
                       until o = w
                       return T.elements() and terminate
     S.pop(v)
```

### **Graph Traversal – Breadth-First Search (BFS)**

#### **Principle:**

- Let s be the start node of BFS, set s to level 0
- In the first step, visit all (not yet visited) vertices that are adjacent to s and set them to level 1.
- In the next step visit for all vertices on level 1 all not yet visited adjacent vertices and set them to level 2.
- Repeat this step as long as all vertices have been reached.

#### **Result:**

- Traversal of the graph
- Level of a vertex *v* shows the length of the shortest path from *v* to *s*.



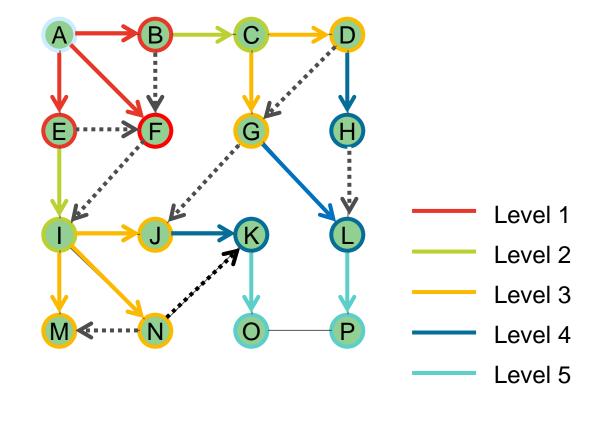
# **Graph Traversal – Breadth-First Search (BFS)**

Pseudo code example:

```
Algorithm BFS(v);
      Input: A vertex s in a graph
      Output: A labeling of edges as "discovery" and "cross" edges
      initialize container L<sub>0</sub> to contain vertex s
      i \leftarrow 0
      while L<sub>i</sub> is not empty do
             create container L<sub>i+1</sub> to initially be empty
             for each vertex v in L<sub>i</sub> do
                   for each edge e incident on v do
                          if edge e is unexplored then
                                 let w be the other endpoint of e
                                 if vertex w is unexplored then
                                       label e as discovery edge
                                        insert w into L<sub>i+1</sub>
                                 else
                                       label e as cross edge
             i ← i+1
```



### **BFS** :: Example



····· cross edge



### **Properties of BFS**

Let *G* be an **undirected graph** in which BFS is executed starting with vertex *s*, then we have:

- The traversal **visits all vertices** in the connected component, which contains s.
- The set of **discovery edges** form a **spanning tree** (BFS tree) for this connected component.
- For each vertex v on **level** i, the **path** to s **along** the **BFS tree** has **length** i and **any other** path from v to s has **at least length** i.
- If (u,v) is an edge that is not in the BFS tree, then the **levels** of u and v differ by 1 at maximum.

#### **Complexity of BFS (analogue DFS)**

If n is the number of vertices of G and m the number of edges, then the complexity of BFS in G is O(n + m)

With the **same complexity** the following tasks can be solved:

- Test, if G is connected
- Calculation of a spanning tree in G
- Calculation of the connected components in G
- Calculation of a cycle in G (or that G has no cycles)



### **Graph Traversal – DFS for Directed Graphs**

#### Same algorithms as for undirected graphs

The result can also be a set of unconnected DFS trees (forest)

#### Accessibility in directed graphs

A DFS tree with root s contains all vertices that can be reached from s via directed edges.

#### **Transitive closure**

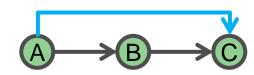
■ The **transitive closure** *G*\* to a graph *G* is obtained by **inserting** a **directed edge** (*v,w*) **if** *v* is **reachable** from *w* (there is a directed path from v to w)

#### Calculation of the **transitive closure** for graph G:

Apply DFS on each vertex of G

■ Complexity: O( n (n+m) )

• Alternative: Floyd-Warshall Algorithm





### Floyd-Warshall Algorithm

Assumption: Operations areAdjacent and insertDirectedEdge have complexity O(1) (Graph is e.g. as adjacency matrix stored)

```
Algorithm FloydWarshall(G); let v_1 \dots v_n be an arbitrary ordering of the vertices of G_0 = G for k = 1 to n do  
// consider all possible routing vertices vk 
G_k = G_{k-1} for each ( i, j = 1,...,n ) ( i != j ) ( i,j != k ) do  
// for each pair of vertices v_i and v_j if G_{k-1}.areAdjacent(v_i, v_k) and G_{k-1}.areAdjacent(v_k, v_j) then G_k. insertDirectedEdge(v_i, v_j, v_k) return G_n
```

**Digraph**  $G_k$  is the **subdigraph** of the **transitive closure** of G, which results from the paths to the **intermediate vertices** of set  $\{v_1, ..., v_k\}$ 

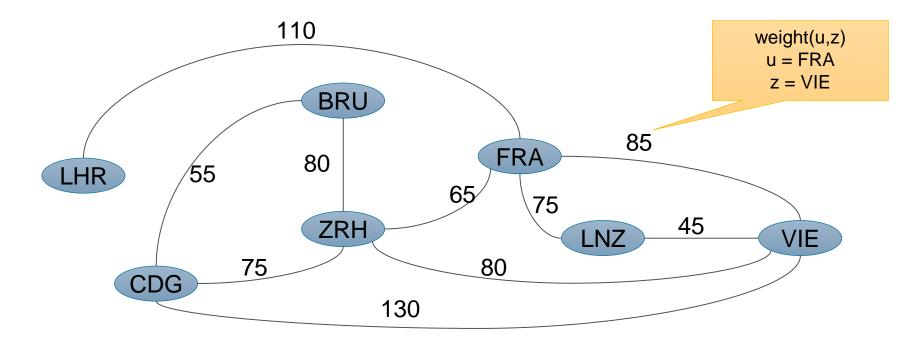
Runtime: O(n³)



#### **Weighted Graph**

An edge is weighted, if a numerical value is assigned to it.

Value can represent for example a distance, travel time or cost.





### **Shortest Path – Dijkstra's Algorithm**

#### BFS finds path with minimum number of edges

Corresponds to the shortest path if all edges have the same weight

#### Dijkstra's algorithm

- Finds shortest path for all vertices z to start vertex s in a graph
  - with undirected edges and
  - with non-negative edge weights
- based on *greedy method*

#### Algorithmic idea

- Set of nodes for which a path has already been found is stored in set C.
- D(z) denotes the **shortest path** from v to s found **so far**.
- When a new node u is visited, the system checks whether a route via this node to an already visited node z has a shorter distance than the shortest route found so far, i.e. whether D[u] + weight(u,z) < D[z]</p>
- If yes, then this path is saved via b as the new shortest path to z and D[z] is updated (relaxation)



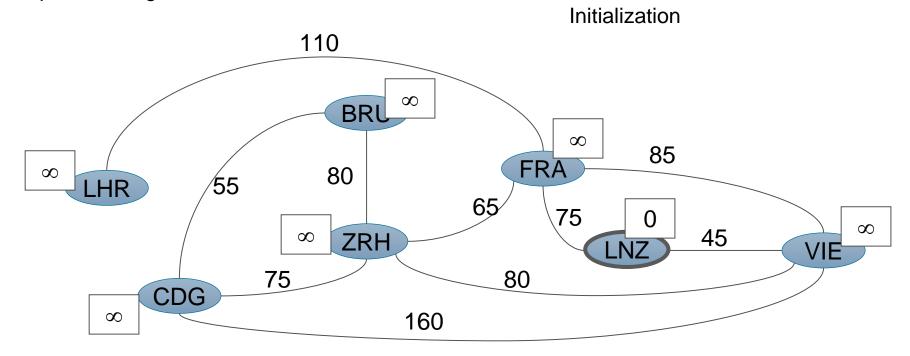
### **Dijkstra's Algorithm**

```
Algorithm ShortestPath(G,s);
           A weighted Graph G and distinguished vertex s of G
    Input:
    Output: labels D[u] for each vertex u of G giving the length of the shortest
             path from u to s in G
    initialize D[s] \leftarrow 0 and D[u] \leftarrow + \infty for each vertex u \neq s
    let Q be a priority queue that contains all of the vertices of G
         using the D labels as keys
    while Q \neq \emptyset do
        // pull u into cloud C
         u \leftarrow Q.removeMinElement()
         for each vertex z adjazent to u such that z is in Q do
             // perform relaxation operation on edge (u,z)
             if D[u] + w((u,z)) < D[z] then
                  D[z] = D[u] + w((u,z))
                  change the key value of z in Q to D[z]
    return label D[u] of each vertex u
```



### **Dijkstra's Algorithm :: Example**

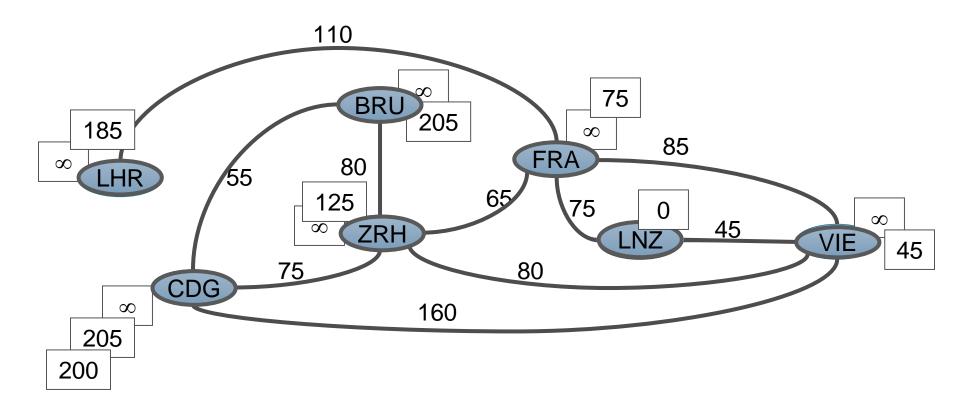
Search shortest path starting from LNZ





# **Dijkstra's Algorithm :: Example**

Search shortest path starting from LNZ





### **Runtime of Dijkstra's Algorithm**

Assumption: **G** is stored as **adjacency list** 

Vertices that are adjacent to u can be visited in O(j), where j is the number of vertices that are adjacent to u

#### Variation 1: Priority Queue Q implemented as heap

Initialization of all vertices in Q O(log n)

Within the loop:

■ removeMin O(log n)

■ **key updates** O(log n) (if locators are used in the heap)

■ Relaxation O(degree(u) log n )

Runtime of the loop

$$\sum_{u \in G} (1 + degree(u)) \log(n)$$

Therefore total runtime of Dijkstra:  $O((n+m) \log n)$ 

To be preferred when number of edges is small:

 $m < n^2 / log n$ 

### **Runtime of Dijkstra's Algorithm**

Assumption: **G** is stored as **adjacency list** 

Vertices that are adjacent to u can be visited in O(j), where j is the number of vertices that are adjacent to u

#### **Variation 2: Priority Queue as unsorted list**

Within the loop

■ removeMin O(n)

key updatesO(1)

Therefore total runtime of Dijkstra:  $O(n^2 + m)$ 

To be preferred if the number of edges is large:

 $m > n^2 / log n$ 



### **Minimum Spanning Tree (MST)**

#### MST is a spanning tree with minimum total edge weight

- Prim-Jarnik algorithm (similar to Dijkstra)
  - Start with (any) start vertex v
  - Tree is constructed vertex by vertex
  - **Insert** edge (*v*,*u*) and vertex *u*, which are:
    - Vertex v is already in the tree
    - Vertex u is not yet in the tree
    - The weight of (v,u) is the **minimum** of the weights **of all possible candidates** u.
  - For each vertex *D[u]* is stored
    - Indicates for a vertex that has not yet been visited, the minimum of the weights of all edges with which u could be connected to the tree.
    - Advantage: shortened search for minimum

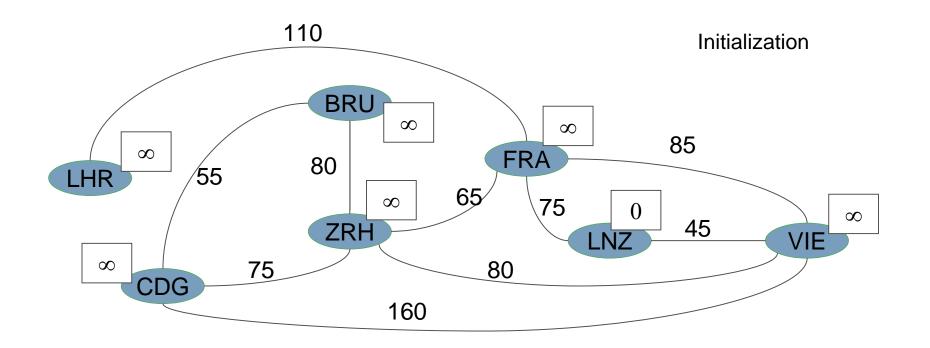


### **Prim-Jarnik Algorithm**

```
Algorithm PrimJarnik(G):
     Input:
                A weighted graph G.
     Output: A minimum spanning tree T for G.
pick any vertex v of G
                                       // grow the tree starting with vertex v
T \leftarrow \{v\}
D[v] \leftarrow 0
E[v] \leftarrow \emptyset
for each vertex u \neq v do
                                        D[u] \leftarrow +\infty
let Q be a priority queue that contains vertices, using the D labels as keys
while Q \neq \emptyset do
                                       // pull u into the cloud C
     u← Q.removeMinElement()
     add vertex u and edge E[u] to T
     for each vertex z adjacent to u do
     if z is in Q
                                        // perform the relaxation operation on edge (u, z)
          if weight(u, z) < D[z] then
                D[z] \leftarrow weight(u, z)
                E[z] \leftarrow (u, z)
                change the key of z in Q to D[z]
                                                                                           runtime
return tree T
                                                                                        O((n+m) \log n)
```

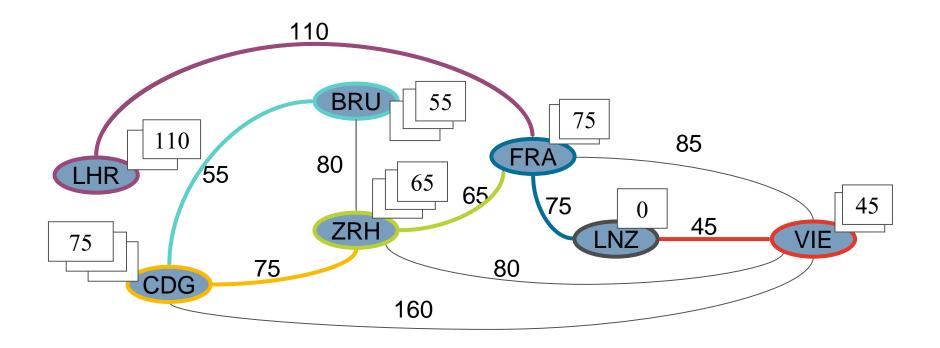


### **Prim-Jarnik Algorithm :: Example**





# **Prim-Jarnik Algorithm :: Example**







# Kruskal's Algorithm

Put one edge after the other into the **MST** under the following conditions:

- Select the edge with the lowest weight
- An edge is only inserted, if no cycle will result from the insertion

#### Data structure:

- Algorithm manages a set of trees (forest)
- An edge is accepted if it connects vertices from different trees

Therefore the data structure must manage disjunctive subsets and support the following operations:

- find(u) returns the set, which *u* contains
- union(u,v) merges the sets, which u and v contain

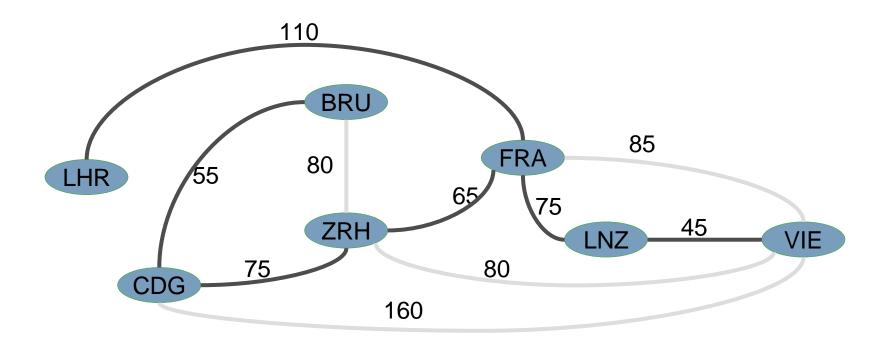


# Kruskal's Algorithm

```
Algorithm Kruskal(G):
    Input:
             A weighted graph G.
    Output: A minimum spanning tree T for G.
let P be a partition of the vertices of G, where each
vertex forms a separate set
let Q be a priority queue storing the edges of G, sorted
by their weights
T \leftarrow \emptyset
while Q \neq \emptyset do
    (u,v) \leftarrow Q.removeMinElement()
    if P.find(u) \neq P.find(v) then
         add edge (u,v) to T
         P.union(u,v)
return T
```



# **Kruskal's Algorithm :: Example**





Edge not in MST



**Distance** between two vertices *u,v* in a Graph *G* (V,E)

- shortest path between them
- G unweighted: d(u,v) number of hops of the shortest path
- G weighted: d(u,v) sum of the weights of the shortest path

### **Eccentricity**

Maximum (shortest) distance from the vertex to any other vertex in a connected graph

$$\sigma(v) = \max_{u \in V} \delta(u, v)$$

#### **Diameter**

is the maximum (shortest) distance between two vertices of a connected graph

$$\Delta = \max_{v \in V} \sigma(v)$$



#### **Radius**

Minimum eccentricity among all vertices

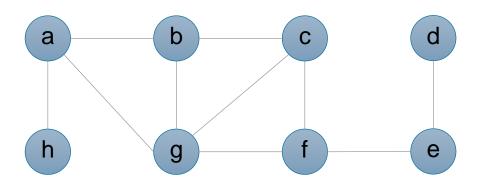
$$rad(G) = \min_{v \in V} \sigma(G)$$

Relation: diameter-radius

$$rad(G) \le diam(G) \le 2rad(G)$$



# **Distance and Centrality :: Example**



### $\Delta(G)$

• h-a-g-f-e-d=5

eccentricities a, b, c, d, e, f, g, h

• 4,4,3,5,4,3,3,5

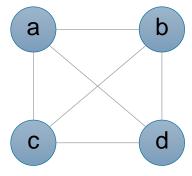
#### radius:

• 3 (min eccentricities)



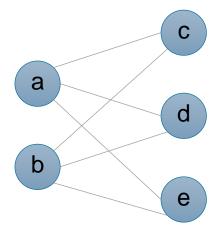
Given a **complete** graph  $K_n$  where  $n \ge 2$ :

• the diameter and radius is 1



Given a **complete bipartite** graph  $K_{m,n}$  where either m or  $n \ge 2$ 

• the diameter and radius is 2





#### **Average distance of a vertex**

The average distance  $d_v(av)$  of a vertex v is defined as the arithmetic mean of the distance of v to all other vertices:

$$d_v(av) = \frac{1}{n-1} \sum_{u \in V} \delta(u, v)$$

### Average distance of a graph

■ The average distance  $d_v(av)$  of a graph G is defined as the arithmetic mean of the distance among all vertices:

$$d_G(av) = \frac{1}{n} \sum_{v \in V} d_v(av)$$



#### Average distance of a graph cont'd.

substituting in equations above:

$$d_G(av) = \frac{1}{n} \sum_{v \in V} d_v(av) = \frac{1}{n} \sum_{v \in V} (\frac{1}{n-1} \sum_{u \in V} \delta(u, v))$$

#### Interpretation

- low average distance => short paths between most of the vertices
- high average distance => general difficult to reach one vertex from another



... vertex degree

$$\delta_{i,j}$$

... (minimum) distance (between vertex *i* and vertex *j*)

$$\Delta = \max(\delta_{i,j} \ 0 \le i, j \le (M-1))$$

... diameter

$$\frac{\sum_{\delta=1}^{\Delta} \delta N_{\delta}}{M-1}$$

... average distance

 $N_{\delta}$ 

... number of distance  $\delta$  vertices

### **Moore Bound**

$$\bar{M} = \begin{cases} 2\Delta + 1 & \text{if } d = 2\\ \frac{d(d-1)^{\Delta} - 2}{d-2} & \text{if } d > 2 \end{cases}$$

d=4

$$\bar{\Delta} = \frac{\log(\frac{\bar{M}+1}{2})}{\log 3}$$



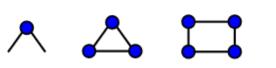
**Moore Graph**: a regular graph of **degree** d and **diameter**  $\Delta$  (also k) whose number of vertices

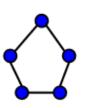
equals the **upper bound** = 
$$\begin{cases} 2\Delta + 1 & if \ d = 2 \\ \frac{d(d-1)^{\Delta} - 2}{d-2} & if \ d > 2 \end{cases}$$

(construction: given d and k, what is the graph with maximum N?)



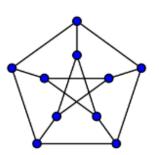




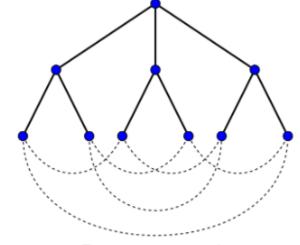


$$d=2$$
  $d=2$   $d=2$   $d=2$   $k=1$   $k=2$   $k=2$ 

N = 3 N = 4 N = 5Moore!



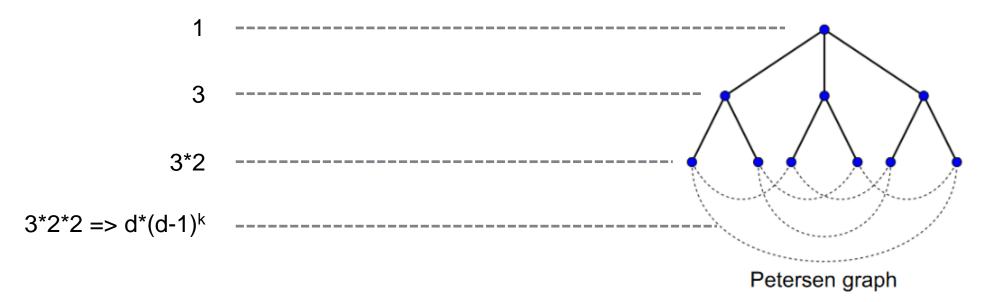
d = 3k = 2N = 10Moore!



Petersen graph

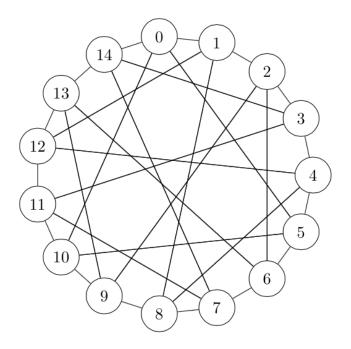
**Moore Graph**: a regular graph of **degree** d and **diameter**  $\Delta$  (also k) whose number of vertices

equals the **upper bound** = 
$$\begin{cases} 2\Delta + 1 & if \ d = 2 \\ \frac{d(d-1)^{\Delta} - 2}{d-2} & if \ d > 2 \end{cases}$$





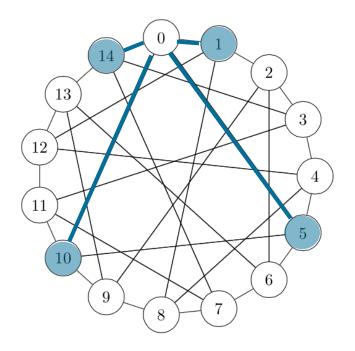
**Moore Bound** 



$$N = 15$$
  $d = 4$   $\Delta = 2$   $Aver.Dist. = 1.71$ 



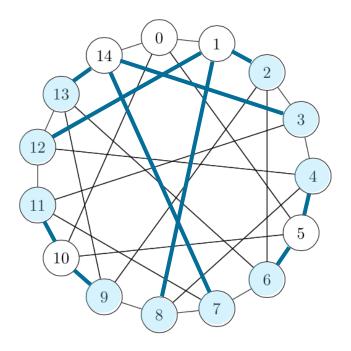
**Moore Bound** 



$$N = 15$$
  $d = 4$   $\Delta = 2$   $Aver.Dist. = 1.71$ 



**Moore Bound** 



$$N = 15$$
  $d = 4$   $\Delta = 2$   $Aver.Dist. = 1.71$ 



#### **Generalized Chordal Rings**

- A graph G is a generalized chordal ring if nodes can be labelled with integers modulo M and there exists a divisor p of M such that node i is joined to node j if node i + p is joined to node j + p.
- G is a chordal ring if all nodes (i,i+1) appear

#### **GCR Construction**

Obtained from a ring by adding chords (= additional link between nodes) by variations over:

- Number c of chords
- Chord length w = number of edges between two nodes (on the ring)
- Period p = number of nodes having the same chord length (same connection pattern)

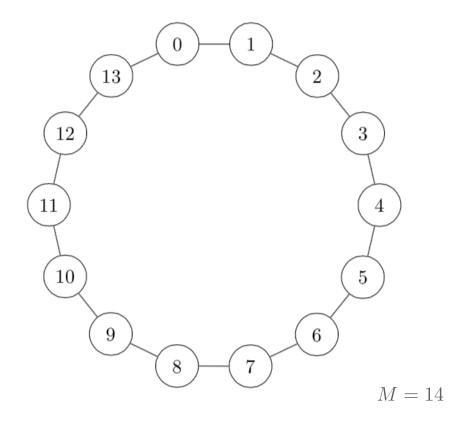
```
0 \le \text{number of chords} \le \text{vertex degree -2}

2 \le \text{chord length} \le \text{M-2}

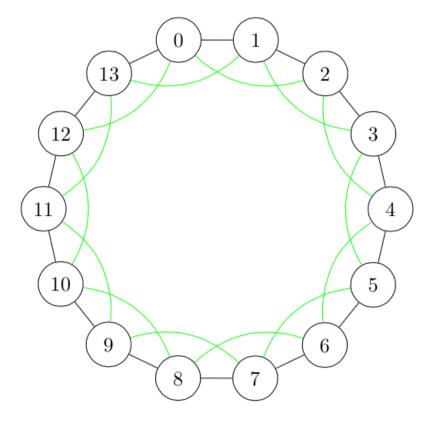
1 \le \text{period} \le \text{M } \text{div } 2

1 \le \text{M } \text{mod period} = 0
```

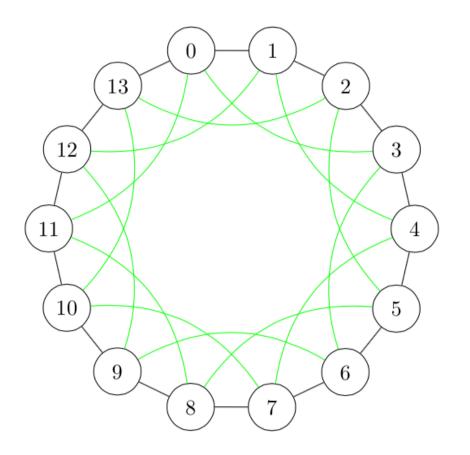




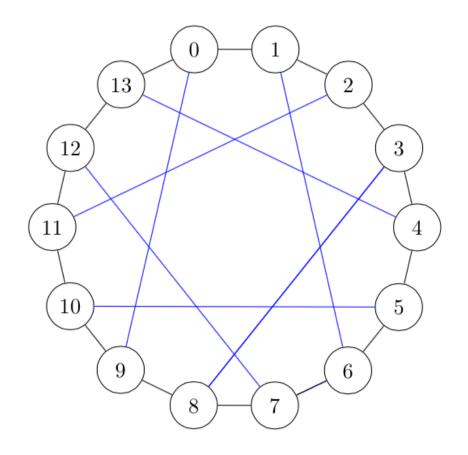




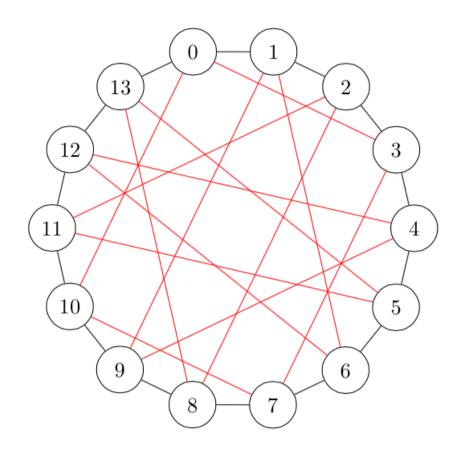
$$P = 1$$
 
$$W = 2$$
 
$$Diameter$$
 
$$\Delta = 4$$

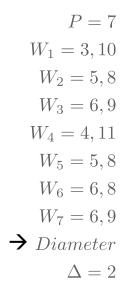








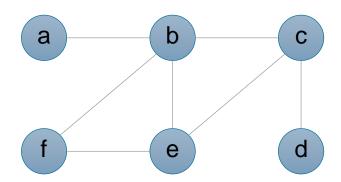




### (Degree) Centrality

Identify the **importance** of specific nodes or edges in the network as a significantly more important node or an edge may be **joining** two **distinctive** parts of the network.

$$C_D(i) = \sum_{j \in V} a_{ij}$$



The equation  $C_D = A \times [1]$  where rows and columns are ordered a,...,f:

$$\begin{bmatrix} 1 \\ 4 \\ 3 \\ 1 \\ 2 \end{bmatrix} = \begin{bmatrix} 0 & 1 & 0 & 0 & 0 & 0 \\ 1 & 0 & 1 & 0 & 1 & 1 \\ 0 & 1 & 0 & 1 & 1 & 0 \\ 0 & 0 & 1 & 0 & 0 & 0 \\ 0 & 1 & 1 & 0 & 0 & 1 \\ 0 & 1 & 0 & 0 & 1 & 0 \end{bmatrix} \times \begin{bmatrix} 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \end{bmatrix}$$

$$\rightarrow \Theta(n^2)$$

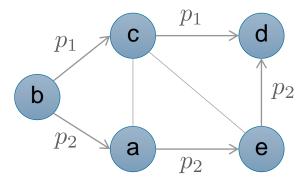


### k-path centrality

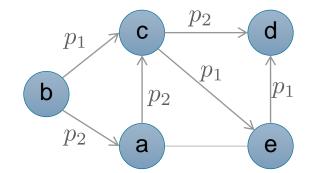
of a vertex v, k- $C_D(v)$  is the **number** of **paths** of **length** k or less emanating from vertex v. When k=1, k-path centrality is equal to degree centrality.

**Edge** disjoint paths **Vertex** disjoint paths

(with common source/sink) do **not** have any **common edges** do **not** have any **common vertices** 



Vertex disjoint paths



Edge disjoint paths (messages from source to sink never have to use the same link)



### Edge disjoint k-path centrality

• of a vertex v is the **number** of **edge disjoint paths** of **length** k or **less** that **start** or **end** at vertex v. The **Ford- Fulkerson Theorem** states that the number of edge disjoint paths between two nodes u and v of a graph is equal to the **minimum number of vertices** that must be **removed to disconnect** u and v.

#### **Vertex disjoint k-path centrality**

- of a node v is the **number of vertex disjoint paths** of **length** k or **less** that start or end at vertex v. **Menger** showed that the number of vertex disjoint paths between two nodes u and v equals the number of **nodes** that must be **removed** to **disconnect** u and v.
- → Major drawback with degree centrality: only **considers local information**.

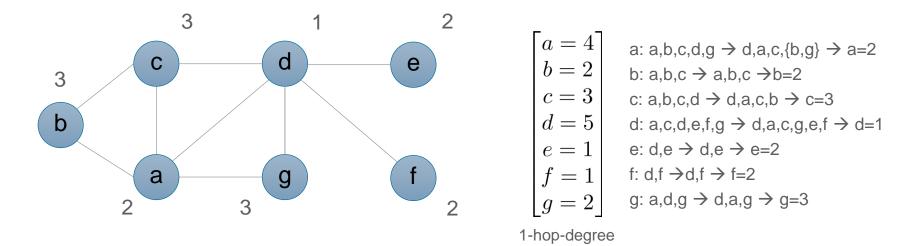
  (a node having few important neighbors may be more influential globally than a node which has many less important neighbors.)



#### k-hop degree (k-hop-C<sub>D</sub>)

• number of neighbors a node has in its k-hop neighborhood. For k=1, k-hop- $C_D$  equals  $C_D$ .

**k-rank** of a node as its position in the descended sorted degree list of neighbors in its k-hop neighborhood.



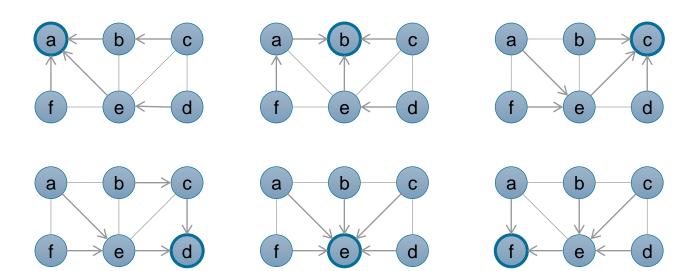
One-rank centrality example



#### **Closeness Centrality**

- based on global rather than local knowledge
- defined as the reciprocal of the total distance from this vertex to all other vertices in a graph G(V,E)

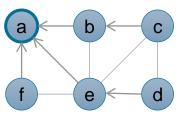
$$C_c(i) = \frac{1}{\sum_{j \in V} \delta(i, j)}$$

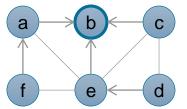


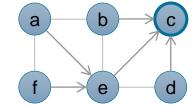
BFS trees for vertex distance calculation

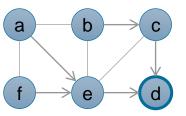


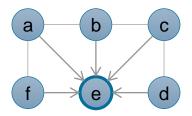
### **Closeness Centrality**

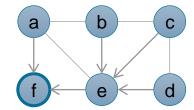












$$\begin{bmatrix} 7 \\ 7 \\ 7 \\ 8 \\ 5 \\ 8 \end{bmatrix} = \begin{bmatrix} 0 & 1 & 2 & 2 & 1 & 1 \\ 1 & 0 & 1 & 2 & 1 & 2 \\ 2 & 1 & 0 & 1 & 1 & 2 \\ 2 & 2 & 1 & 0 & 1 & 2 \\ 1 & 1 & 1 & 1 & 0 & 1 \\ 1 & 2 & 2 & 2 & 1 & 0 \end{bmatrix} \times \begin{bmatrix} 1 \\ 1 \\ 1 \\ 1 \\ 1 \end{bmatrix}$$

$$C_c(a) = \frac{1}{7}$$

$$C_c(b) = \frac{1}{7}$$

$$C_c(c) = \frac{1}{7}$$

$$C_c(d) = \frac{1}{8}$$

$$C_c(e) = \frac{1}{5}$$

$$C_c(f) = \frac{1}{5}$$

Highest closeness centrality

### **Stress Centrality**

total number of all pairs shortest paths that pass through a vertex v

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

where  $\sigma_{st(v)}$  is the number of shortest paths between vertices s and t.

It is an estimation of the **stress** that a vertex in a network bears, assuming all communication will be carried along the shortest path.

### **Betweenness Centrality**

If the paths for **shortest paths between** nodes of a network pass through some vertices **more often** than others, then these vertices are significantly **more important** than others for communication purposes.

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

Simple procedure to find  $C_B$  is to **calculate all shortest paths** in the graph G and **count** the number of paths that pass through v by excluding the paths that start or end at v.

where  $\sigma_{st(v)}$  is the total number of shortest paths between vertices s and t that pass through vertex v

Perform modified BFS on each node  $\rightarrow$   $\Theta(n(n+m))$ 



#### **Betweenness Centrality :: Newman's Algorithm**

```
forall v in V
    b_v \leftarrow 1
     find BFS paths from v to all other nodes
end
forall s in V
     starting from the farthest nodes, move from u towards s along paths
     using vertex v.
     b_v \leftarrow b_v + b_{ii}
     If v has more than one predecessor, then b<sub>v</sub> is devided equally
     between them.
end
```



### **Eigenvalue Centrality**

- a vertex in a network may be considered important if it has important neighbors.
- Given a graph G(V,E) with an adjacency matrix A, the score of a vertex i can be defined as proportional to the sum of all its neighbors' scores:
  1

$$x_i = \frac{1}{\lambda} \sum_{j \in N(i)} x_j = \frac{1}{\lambda} \sum_{j \in V} a_{ij} x_j$$

where N(i) is the set of neighbors of i and  $\lambda$  is a constant

■ The **score** of a node can simply be its **degree**. We are **adding** the **degrees of** the **neighbors** of a vertex *i* to find its **new score**.

$$\overrightarrow{x} = \frac{1}{\lambda} A \overrightarrow{x}$$
 which is the eigenvalue equation  $A \overrightarrow{x} = \lambda \overrightarrow{x}$ 

■ To find the **eigenvalue centrality** of a graph *G*, we need to find the **eigenvalues** of the **adjacency matrix** *A*. We select the largest eigenvalue and its associated **eigenvector**. This eigenvector contains the **eigenvalue centralities** of all vertices.



# **Graphs**Part: Structure



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