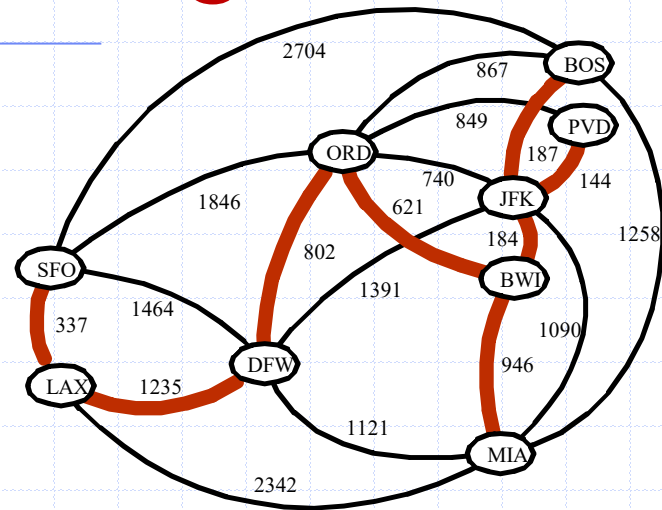


# Lecture 14: Minimum Spanning Trees

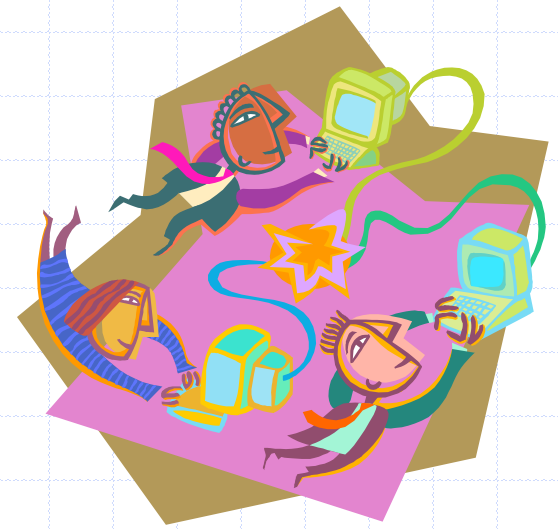
Infinite Correlation



# Wholeness Statement

A minimum spanning tree is a spanning tree subgraph with minimum total edge weight. Efficient greedy algorithms have been developed to compute MST both with and without special data structures. Pure creative intelligence is the source of all creative algorithms. Regular practice of TM improves our ability to make use of our own innate creative potential.

# Outline and Reading



## ◆ Minimum Spanning Trees (§7.3)

- Definitions
- Cycle Property
- Partition Property

## ◆ The Prim-Jarnik Algorithm (1957, 1930) (§7.3.2)

## ◆ Kruskal's Algorithm (1956) (§7.3.1)

## ◆ Baruvka's Algorithm (1926) (§7.3.3)

# Minimum Spanning Tree

## Spanning subgraph

- Subgraph of a graph  $G$  containing all the vertices of  $G$

## Spanning tree

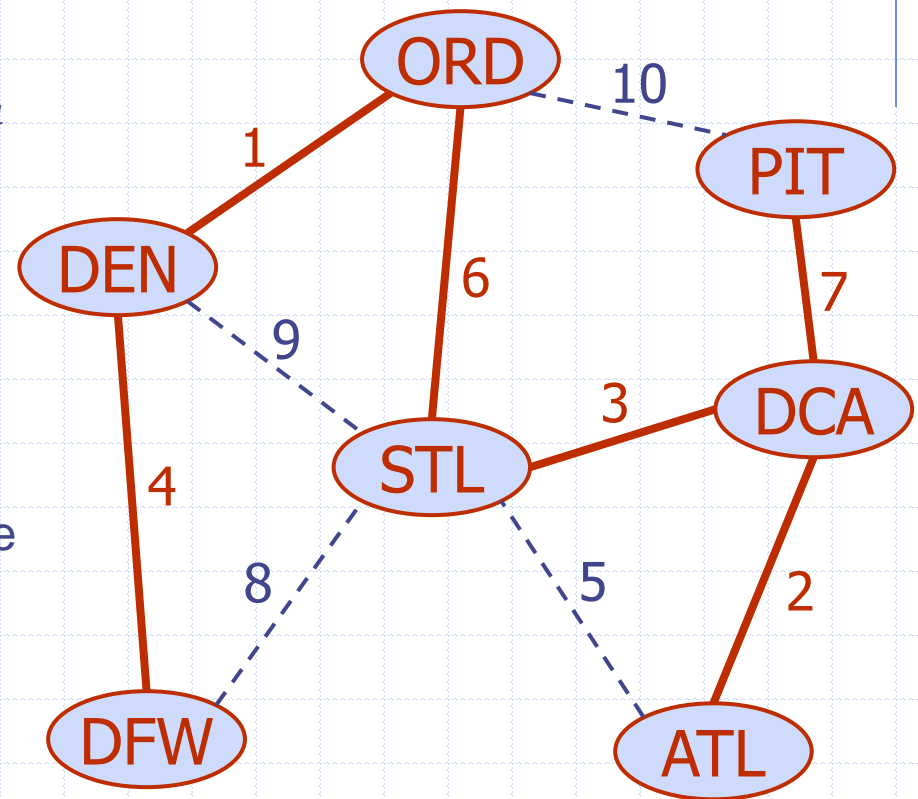
- Spanning subgraph that is itself a (free) tree

## Minimum spanning tree (MST)

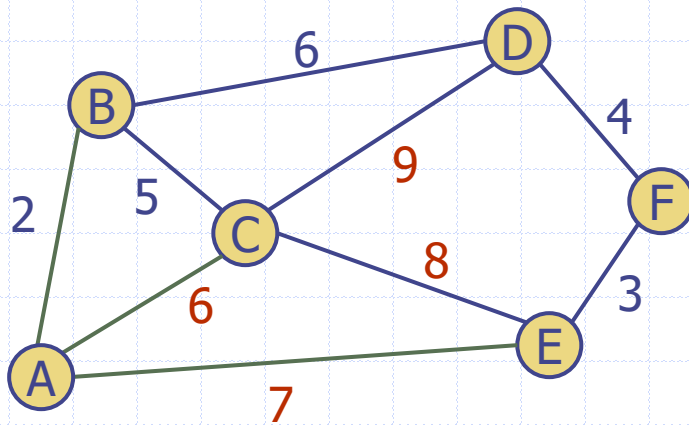
- Spanning tree of a weighted graph with minimum total edge weight

## ◆ Applications

- Communications networks
- Transportation networks

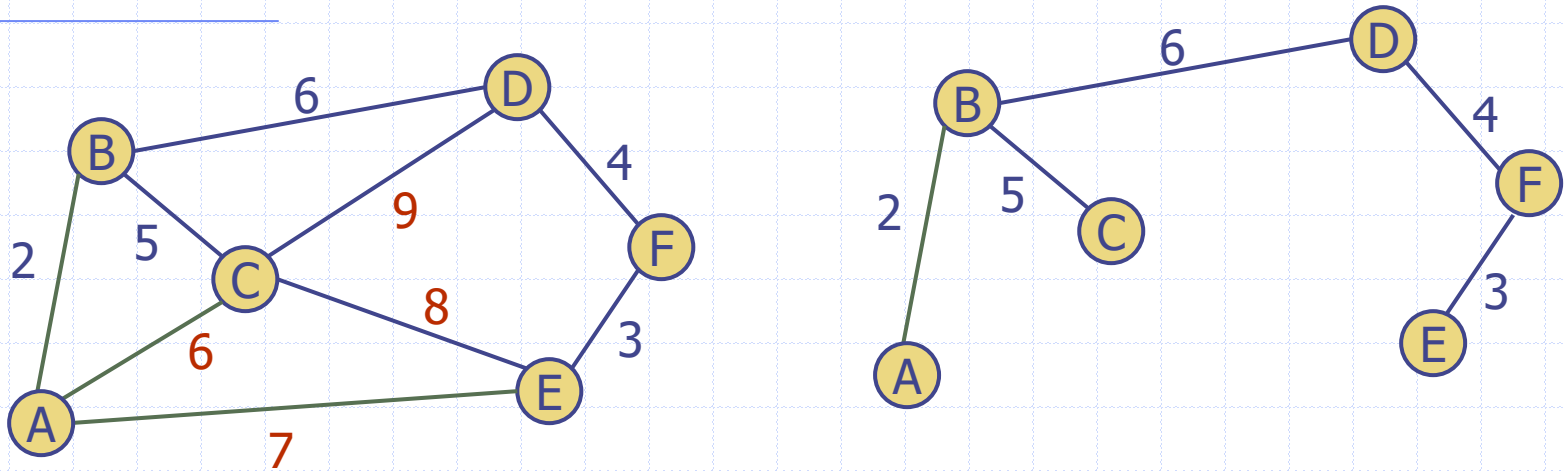


# Cycle Property



If the weight of an edge  $e$  of a cycle  $C$  is larger than the weights of other edges of  $C$ , then this edge cannot belong to a MST.

# Cycle Property



If the weight of an edge  $e$  of a cycle  $C$  is larger than the weights of other edges of  $C$ , then this edge cannot belong to a MST.

# Cycle Property

- ◆ **Cycle Property:** For any cycle  $C$  in a graph, if the weight of an edge  $e$  of  $C$  is larger than the weights of other edges of  $C$ , then this edge cannot belong to a MST.
- ◆ **Proof:** Assume the contrary, i.e. that  $e$  belongs to an MST  $T_1$ ; then deleting  $e$  will break  $T_1$  into two subtrees with the two ends of  $e$  in different subtrees. The remainder of  $C$  reconnects the subtrees, hence there is an edge  $f$  of  $C$  with ends in different subtrees, i.e., it reconnects the subtrees into a tree  $T_2$  with weight less than that of  $T_1$ , because the weight of  $f$  is less than the weight of  $e$ ; thus  $T_1$  cannot be a MST.

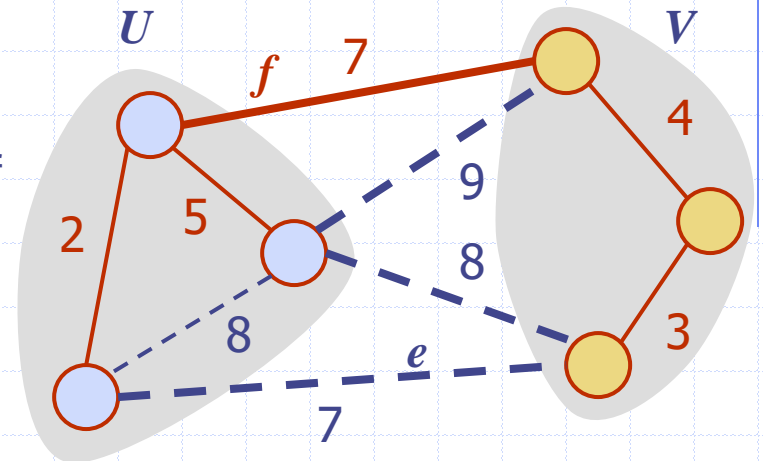
# Partition Property- A Crucial Fact

## Partition Property:

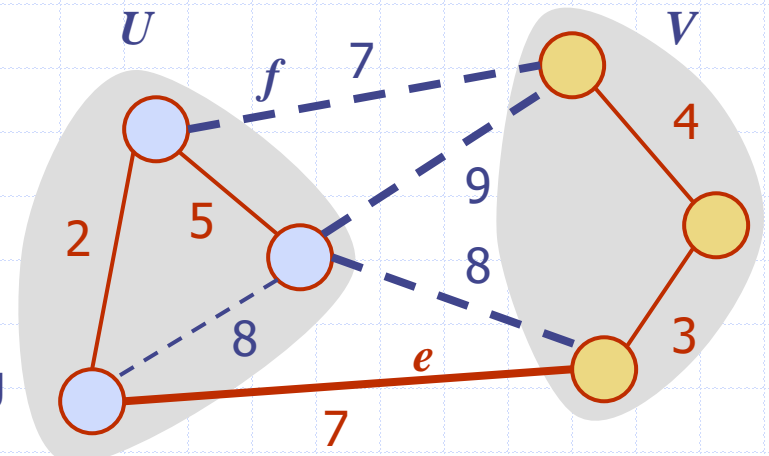
- Consider a partition of the vertices of  $G$  into subsets  $U$  and  $V$
- Let  $e$  be an edge of minimum weight across the partition
- There is a minimum spanning tree of  $G$  containing edge  $e$

## Proof:

- Let  $T$  be an MST of  $G$
- If  $T$  does not contain  $e$ , consider the cycle  $C$  formed by  $e$  with  $T$  and let  $f$  be an edge of  $C$  across the partition
- By the cycle property,  
$$\text{weight}(f) \leq \text{weight}(e)$$
- Thus,  $\text{weight}(f) = \text{weight}(e)$
- We obtain another MST by replacing  $f$  with  $e$



Replacing  $f$  with  $e$  yields another MST





# Generic MST Algorithm

**Algorithm** *GenericMST(G)*

$T \leftarrow$  a tree with all vertices in  $G$ , but no edges

**while**  $T$  *does not form a spanning tree* **do**

$(u, v) \leftarrow$  a safe edge of  $G$

$T \leftarrow (u, v) \cup T$

**return**  $T$

A *safe edge* is one that when added to  $T$  forms a subgraph of a MST

# Main Point

1. A minimum spanning tree algorithm gradually grows a (sub-solution) tree by adding a “safe edge” that connects a vertex in the tree to a vertex not yet in the tree.

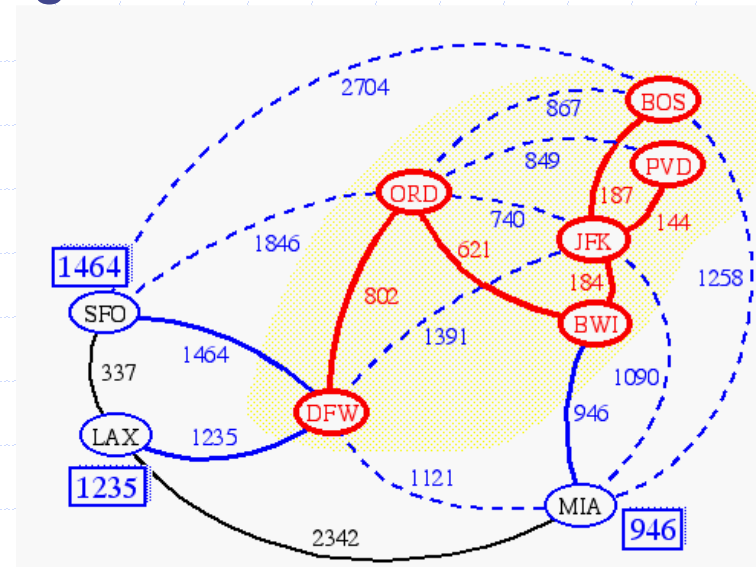
*Science of Consciousness:* The nature of life is to grow and progress to the state of enlightenment, fulfillment.

# Prim(1957)-Jarnik(1930) Algorithm

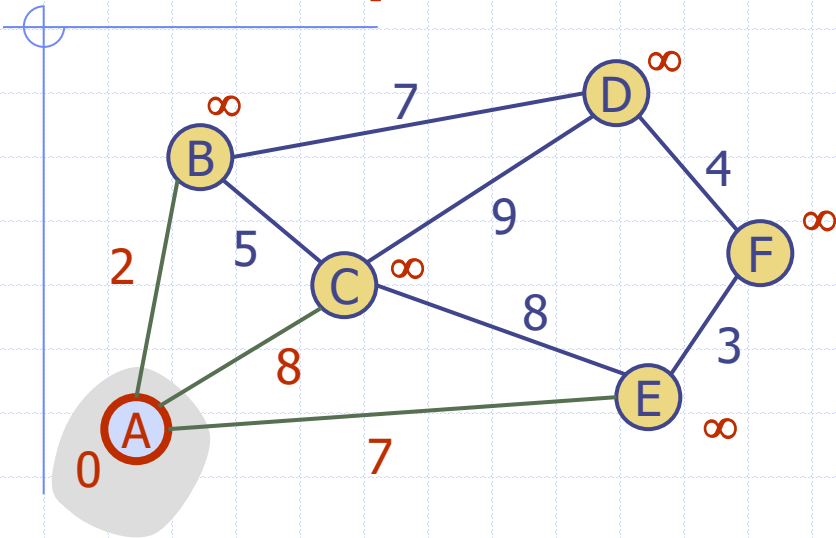
AKA Dijkstra-Prim Algorithm

# Prim-Jarnik's Algorithm

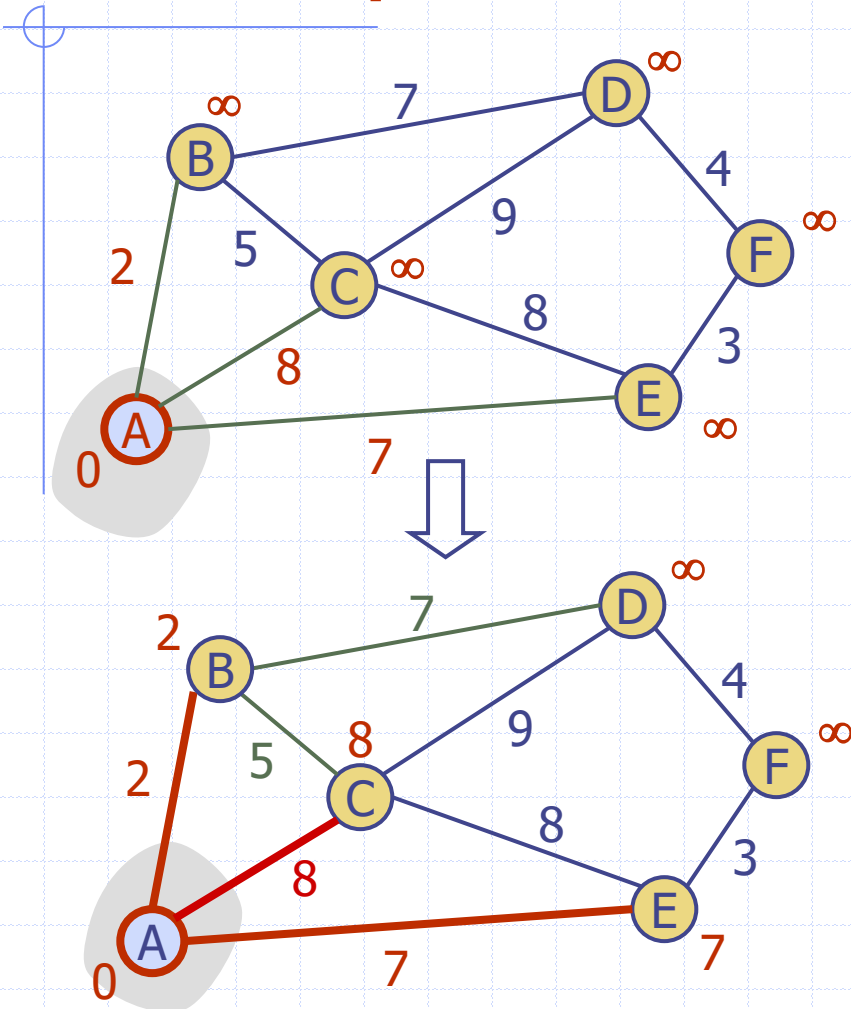
- ◆ Similar to Dijkstra's shortest path algorithm (for a connected graph)
- ◆ We pick an arbitrary vertex  $s$  and we grow the MST as a cloud of vertices, starting from  $s$
- ◆ We store with each vertex  $v$  a label  $d(v)$  = the smallest weight of an edge connecting  $v$  to a vertex in the cloud
- ◆ At each step:
  - We add to the cloud the vertex  $u$  outside the cloud with the smallest distance label
  - We update the labels of the vertices adjacent to  $u$



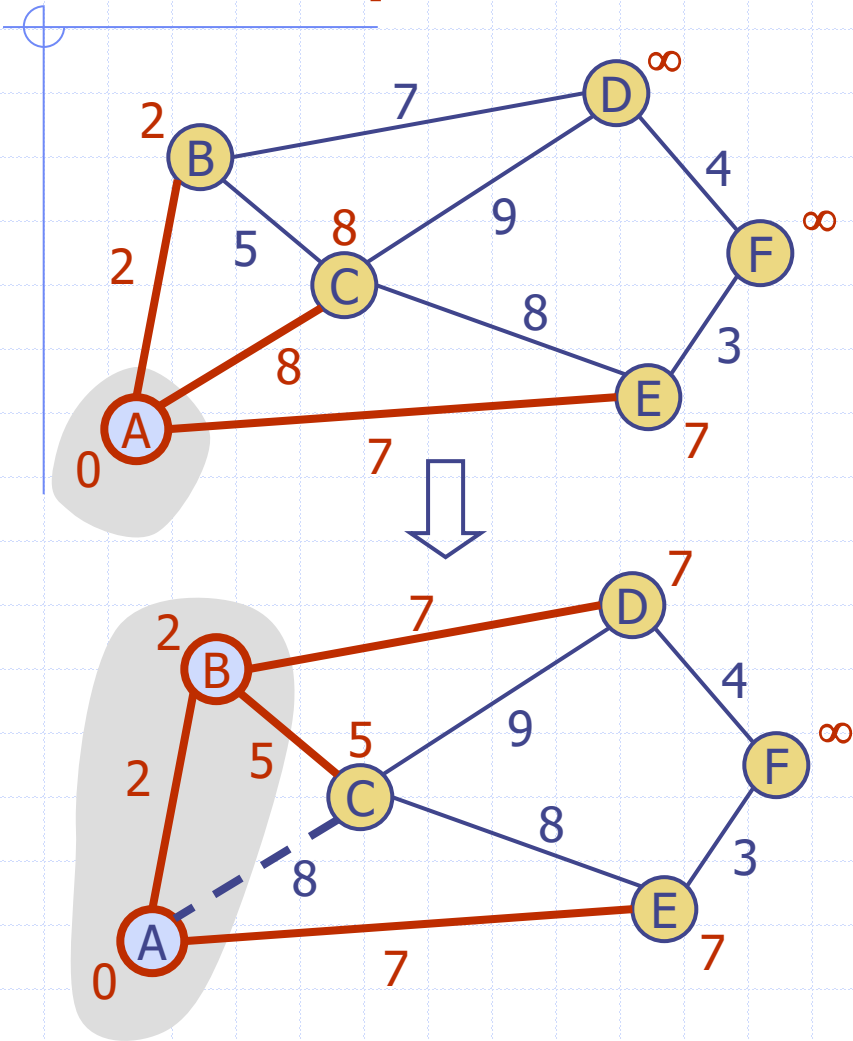
# Example



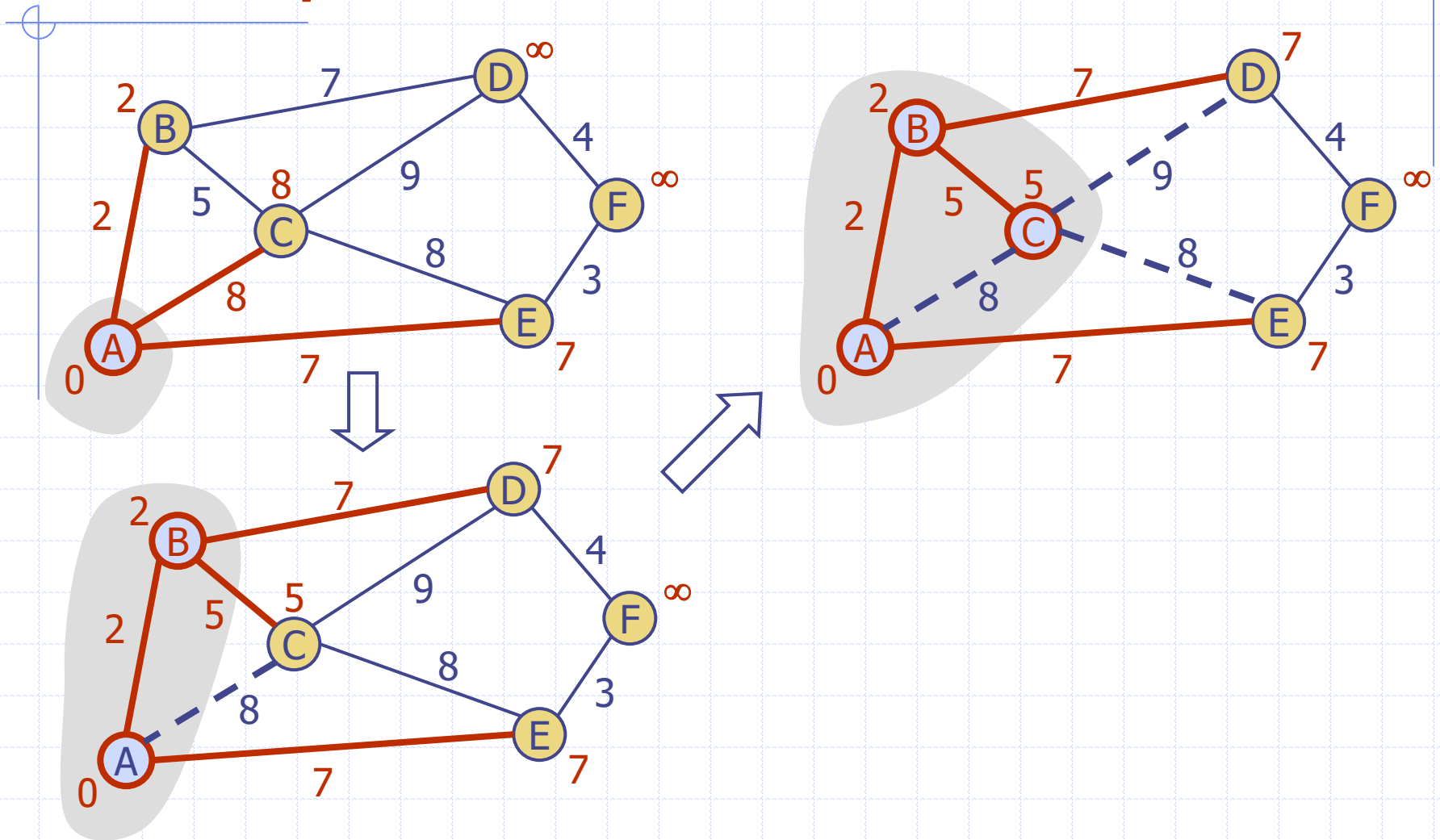
# Example



# Example

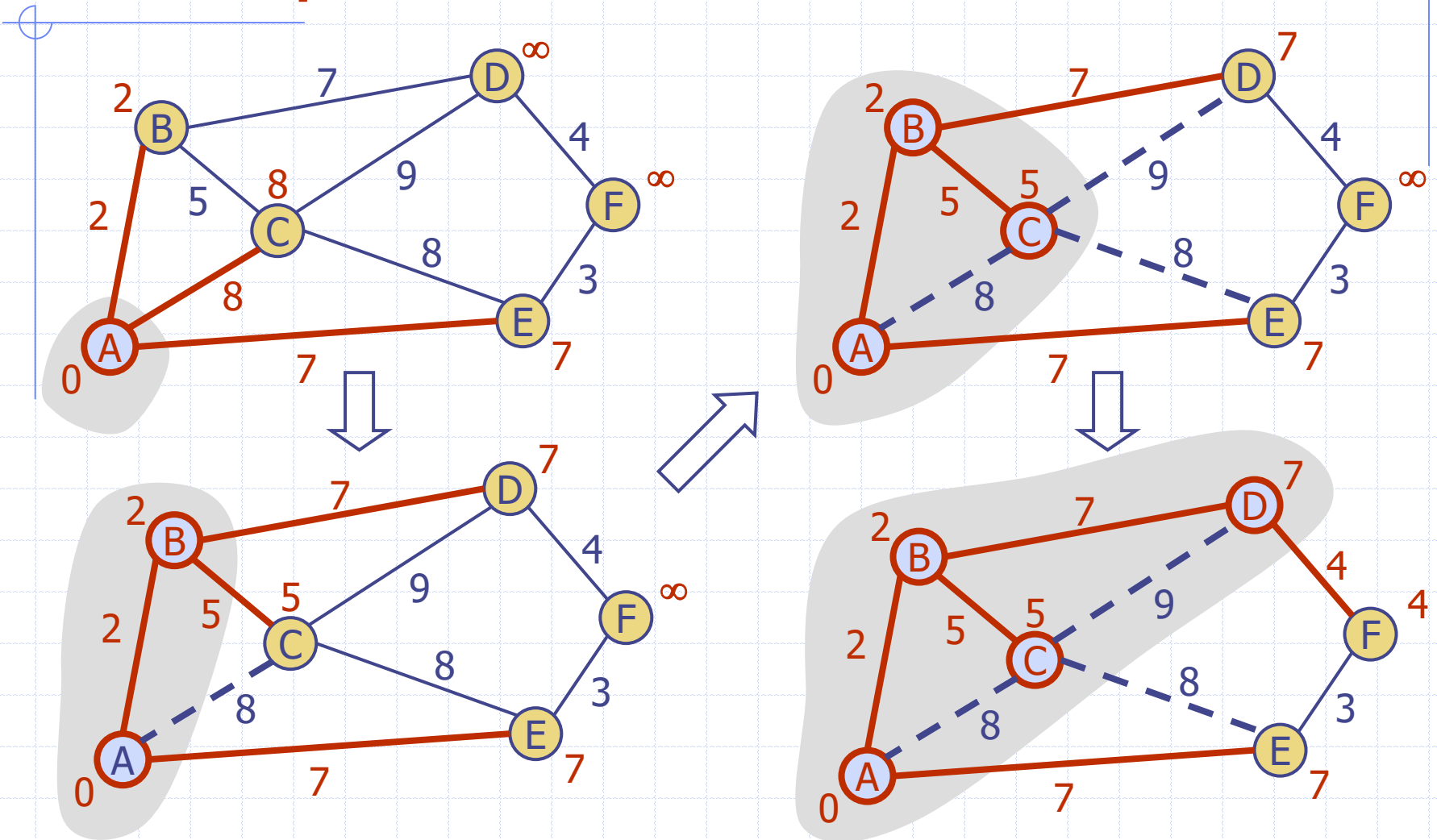


# Example

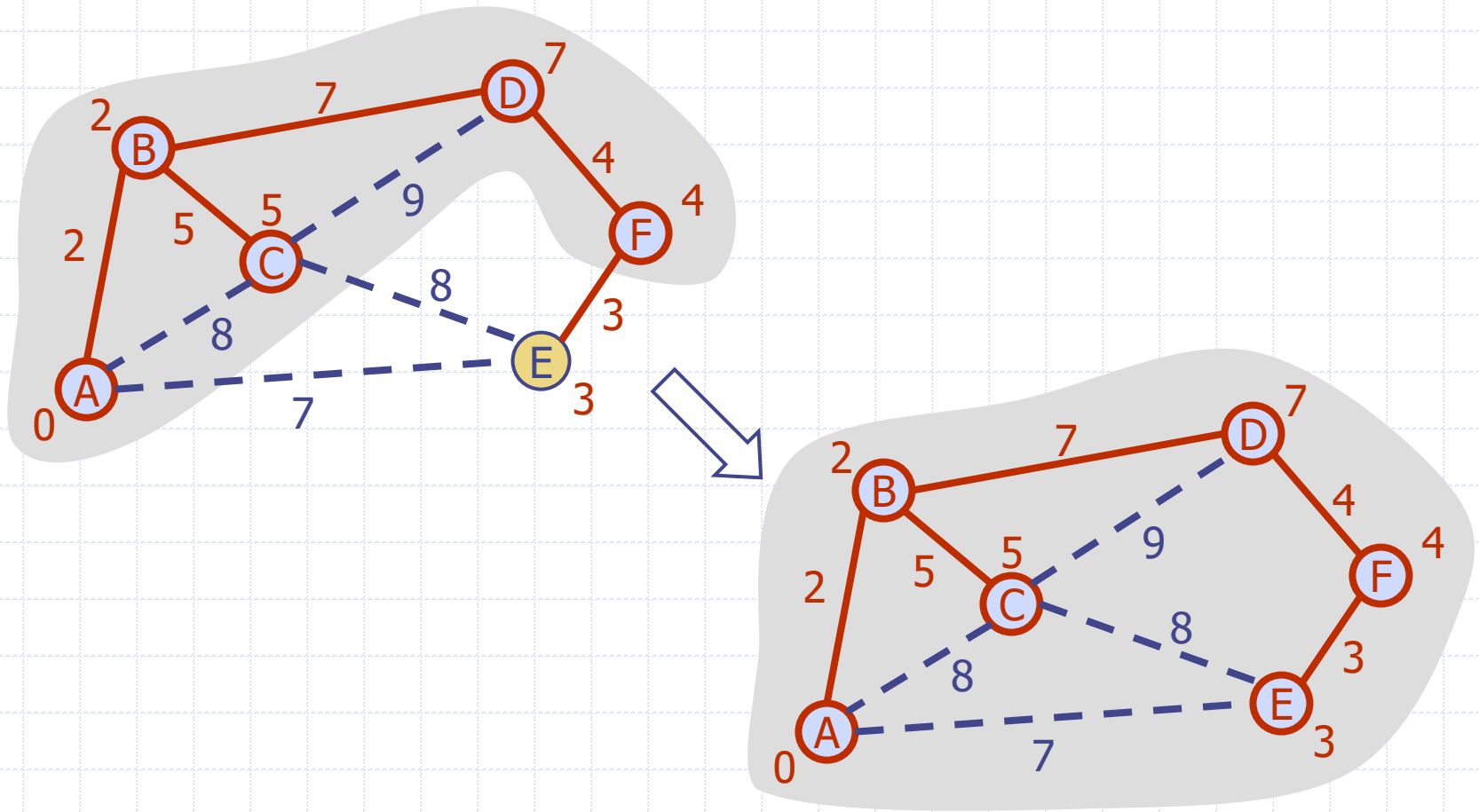




# Example



# Example (contd.)



# Prim-Jarnik's Algorithm (cont.)

- ◆ A priority queue stores the vertices outside the cloud
  - Key: distance
  - Element: vertex
- ◆ Locator-based methods
  - *insert(k, e)* returns a locator position
  - *replaceKey(l, k)* changes the key of an item
- ◆ We store three labels with each vertex:
  - Distance
  - Parent edge in MST
  - Locator in priority queue
- ◆ Correction in red inspired by Bereket Chalew (May 2014)

## Algorithm *PrimJarnikMST(G)*

```
s ← G.aVertex ()
Q ← new heap-based priority queue
for all v ∈ G.vertices() do
  if v = s then
    setDistance(v, 0)
  else
    setDistance(v, ∞)
    setParent(v, ∅)
    l ← Q.insertItem(getDistance(v), v)
    setLocator(v, l)
while ! Q.isEmpty() do
  u ← Q.removeMin()
  setLocator(u, ∅) {u is now in MST}
  for all e ∈ G.incidentEdges(u) do
    z ← G.opposite(u, e)
    r ← weight(e) {diff. from ShortestPath}
    if getLocator(z) ≠ ∅ {z not yet in MST}
      ∧ r < getDistance(z) then
        setDistance(z, r)
        setParent(z, e)
        Q.replaceKey(getLocator(z), r)
```

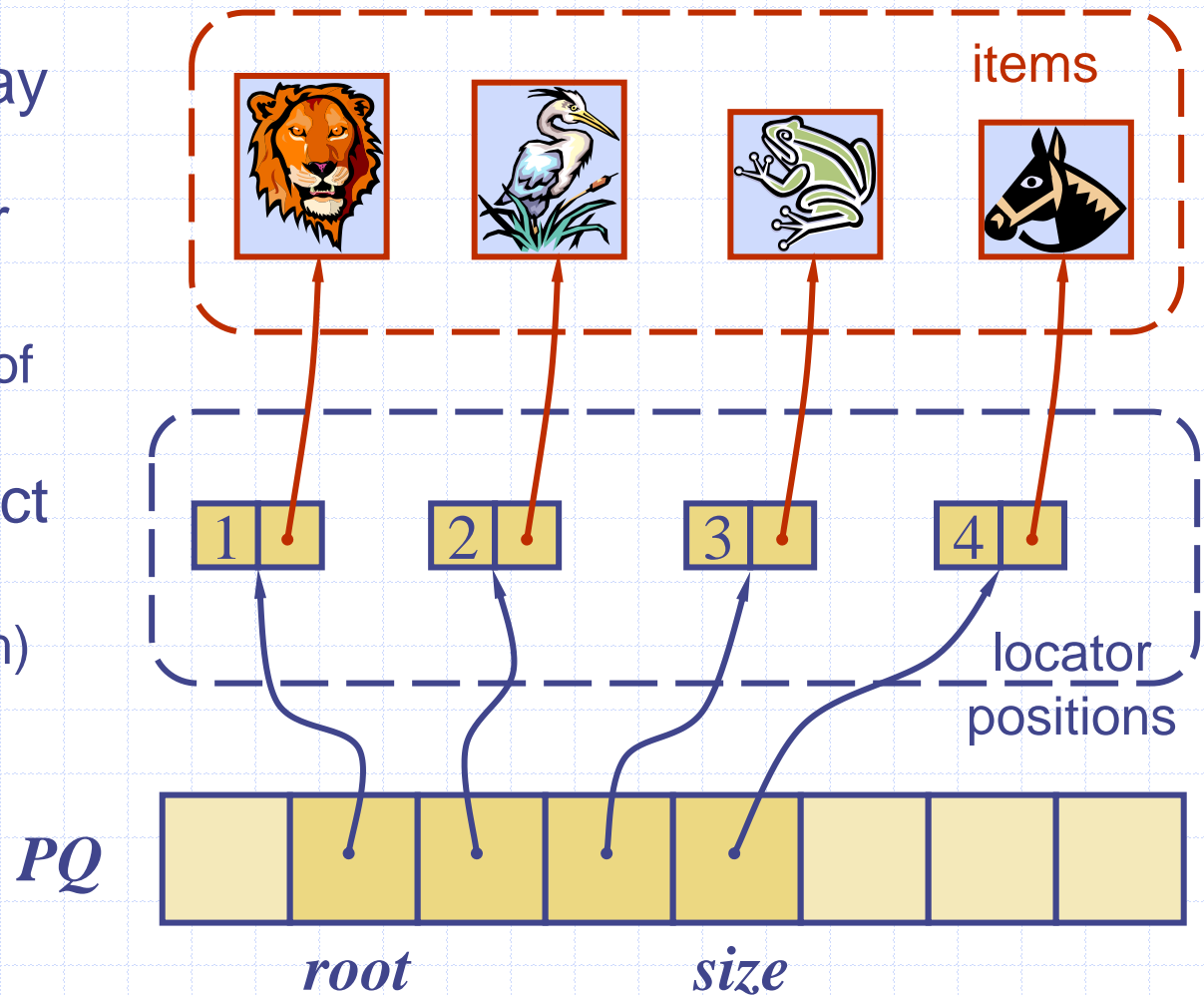
# Array-based Implementation

◆ We use an array storing locator-positions in our Priority Queue

- Another level of indirection

◆ A position object stores:

- Item (key,elem)
- Index



# Analysis

## ◆ Graph operations

- Method `incidentEdges` is called once for each vertex
- Recall that  $\sum_v \deg(v) = 2m$

## ◆ Label operations

- We set/get the distance, parent and locator labels of vertex  $z$   $O(\deg(z))$  times
- Setting/getting a label takes  $O(1)$  time

## ◆ Priority queue operations

- Each vertex is inserted once into and removed once from the priority queue, where each insertion or removal takes  $O(\log n)$  time
- The key of a vertex  $w$  in the priority queue is modified at most  $\deg(w)$  times, where each key change takes  $O(\log n)$  time

## ◆ Prim-Jarnik's algorithm runs in $O((n + m) \log n)$ time provided the graph is represented by the adjacency list structure

## ◆ The running time is $O(m \log n)$ since the graph is connected

# Main Point

2. A defining feature of the Minimum Spanning Tree (and shortest path) greedy algorithms is that once a vertex becomes in-tree (or “inside the cloud”), the resulting subtree is optimal and nothing can change this state.

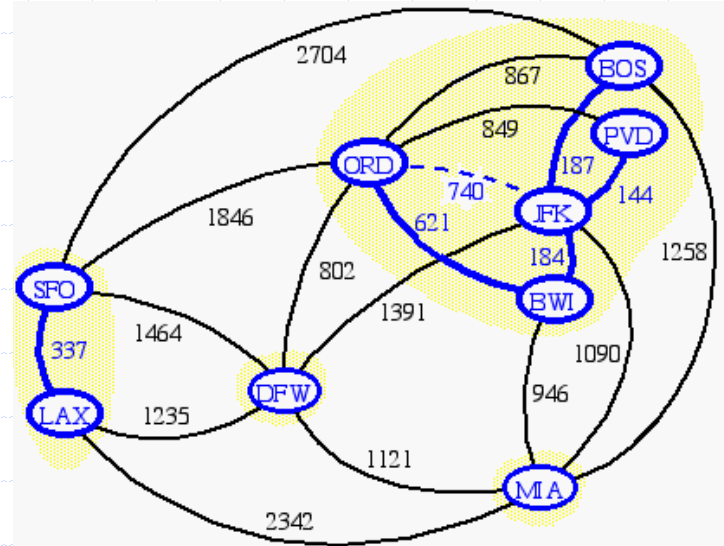
*Science of Consciousness:* A defining feature of enlightenment is that once this state is reached, one’s consciousness is optimal and nothing can change this state.

# Kruskal's Algorithm

Based on the Partition Property

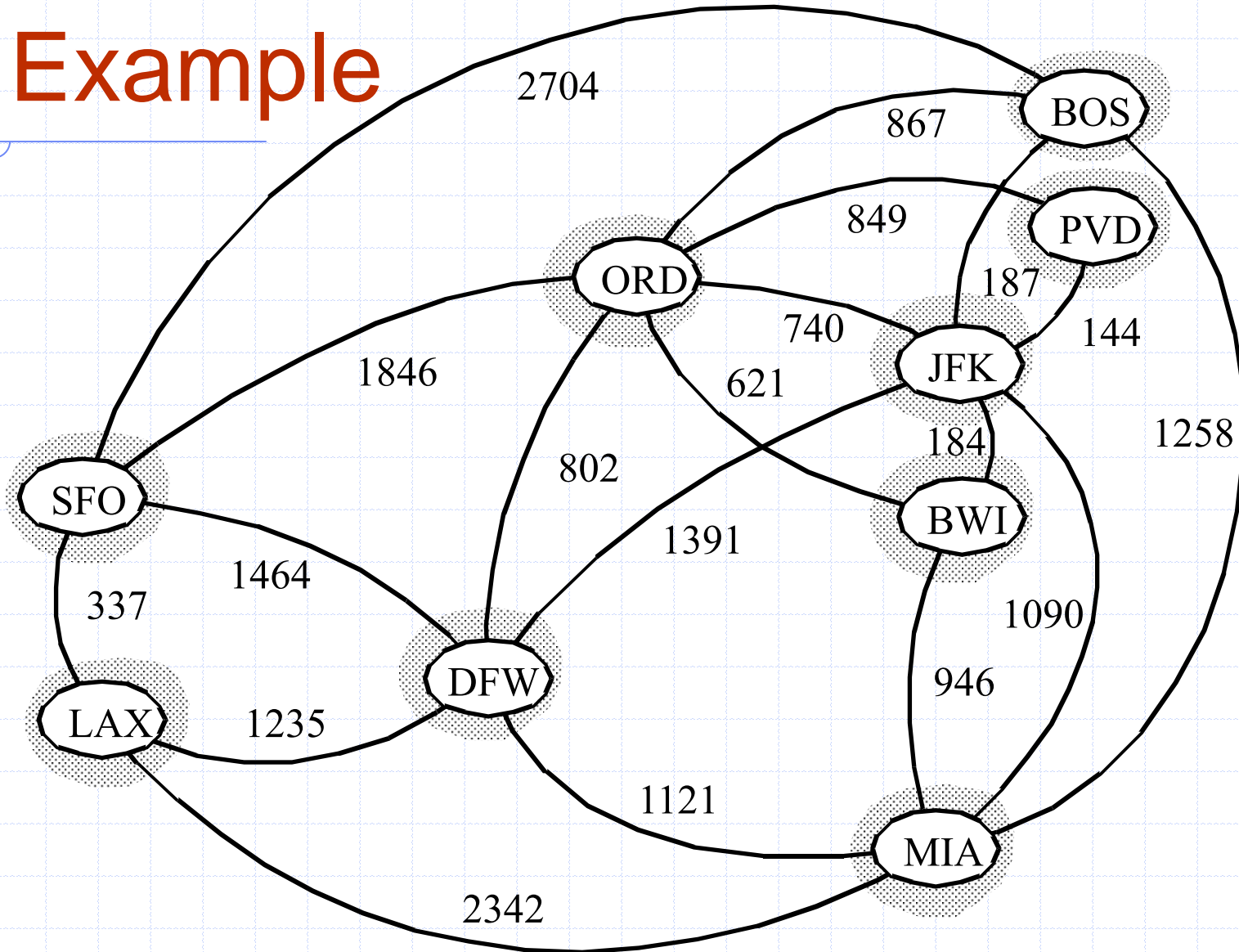
# Basic Idea of the Kruskal Algorithm

- ◆ The algorithm maintains a forest of trees
- ◆ Select the edge with the smallest weight
  - The edge is accepted if it connects distinct trees
- ◆ Keep adding edges until the tree has  $n-1$  edges

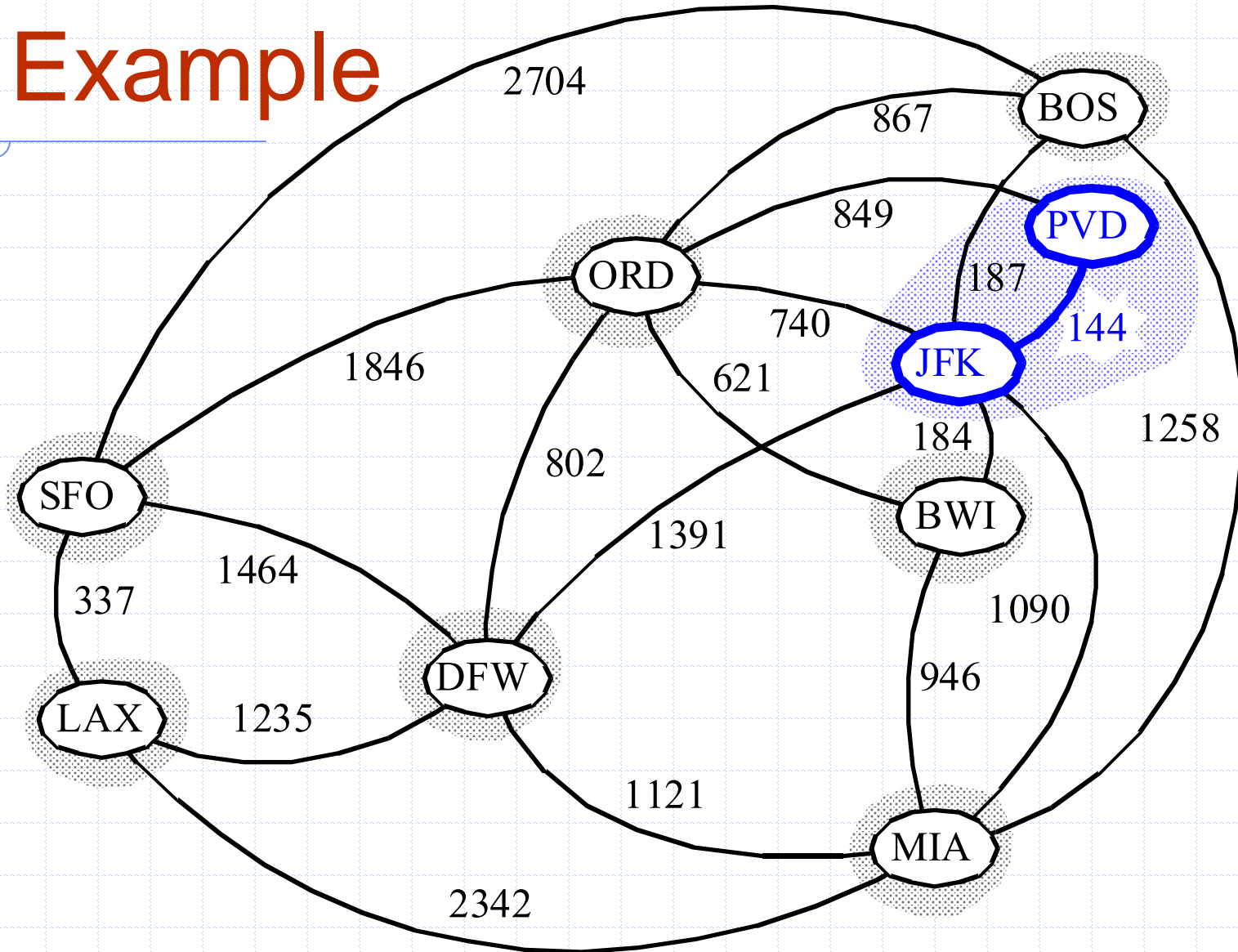




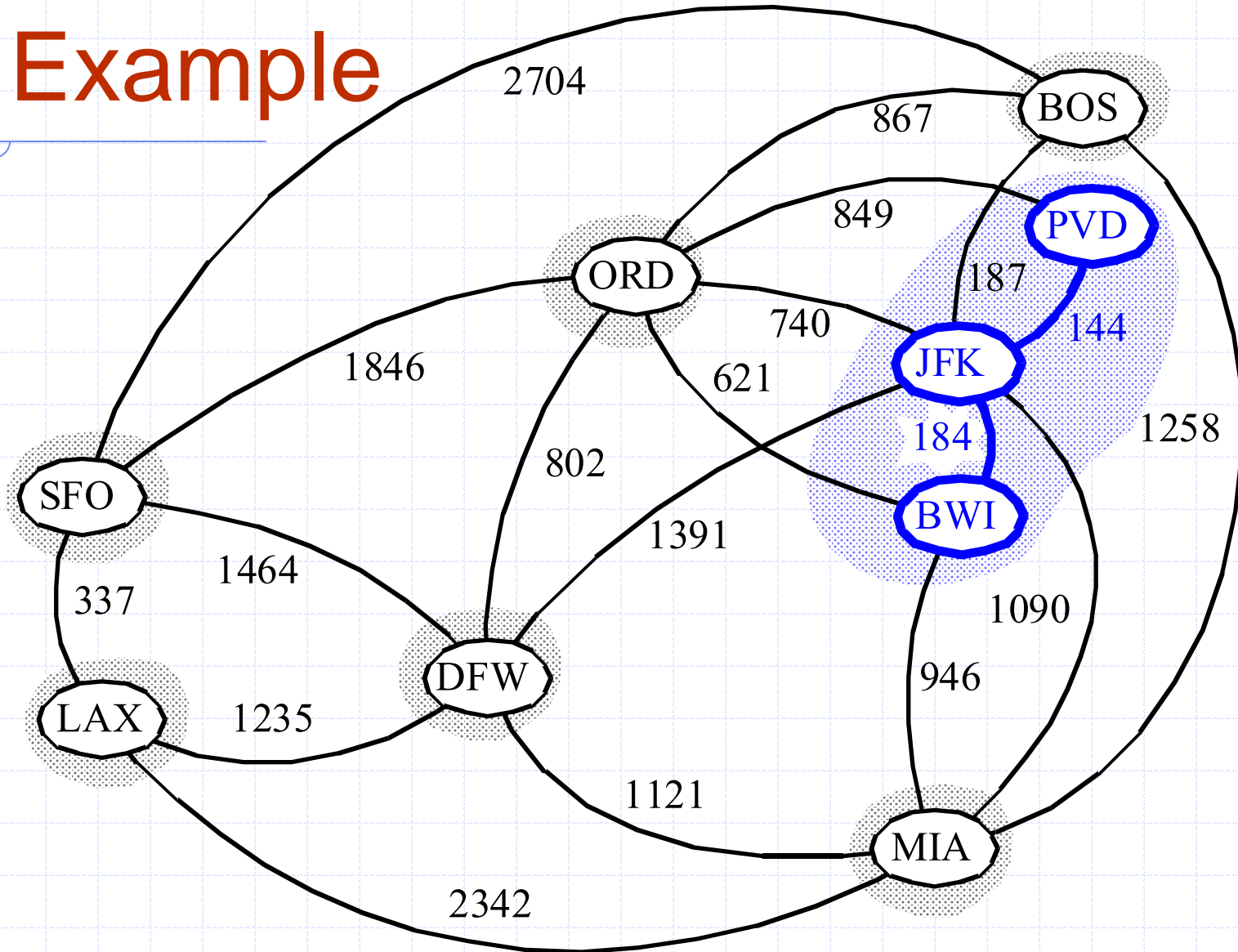
# Kruskal Example



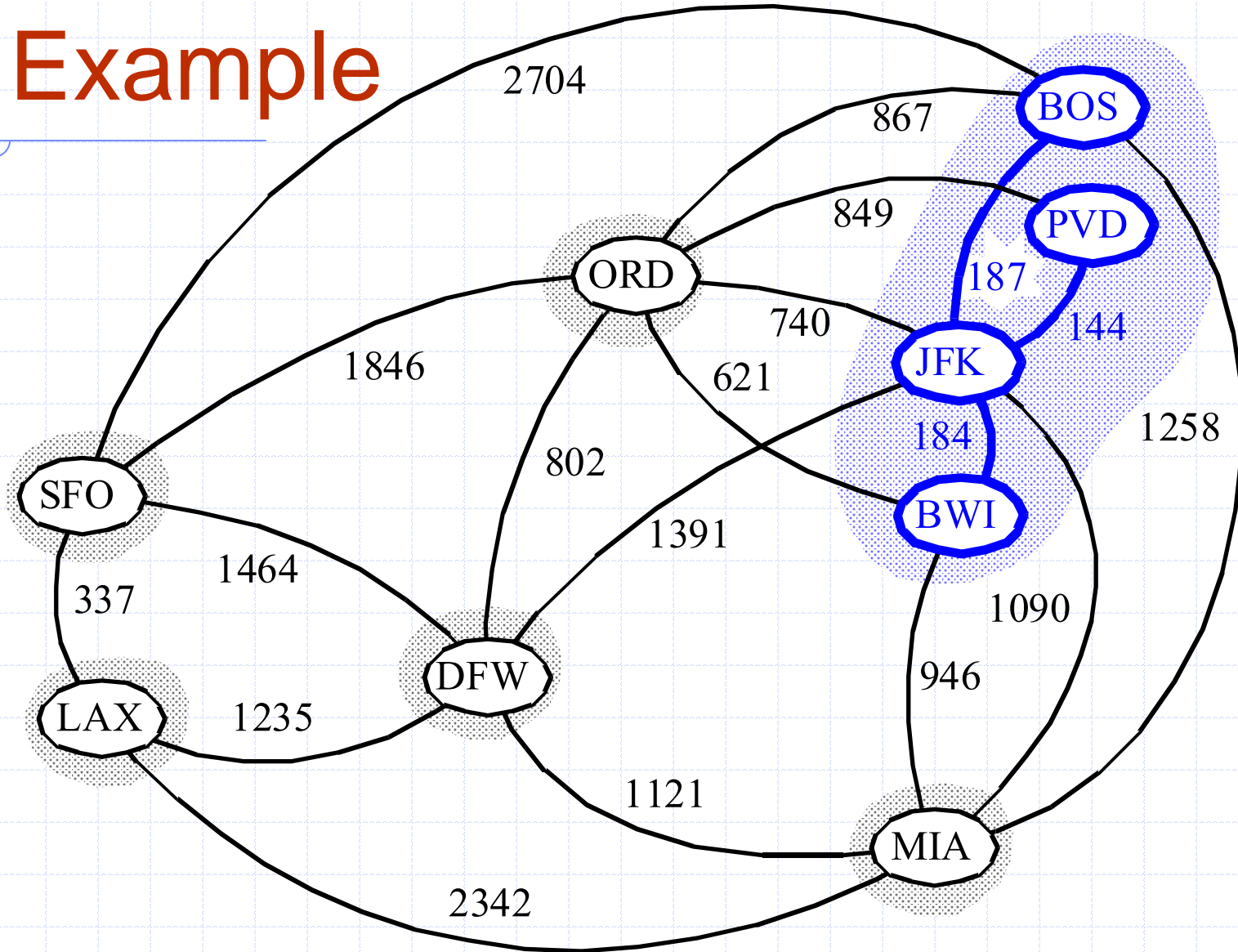
# Example



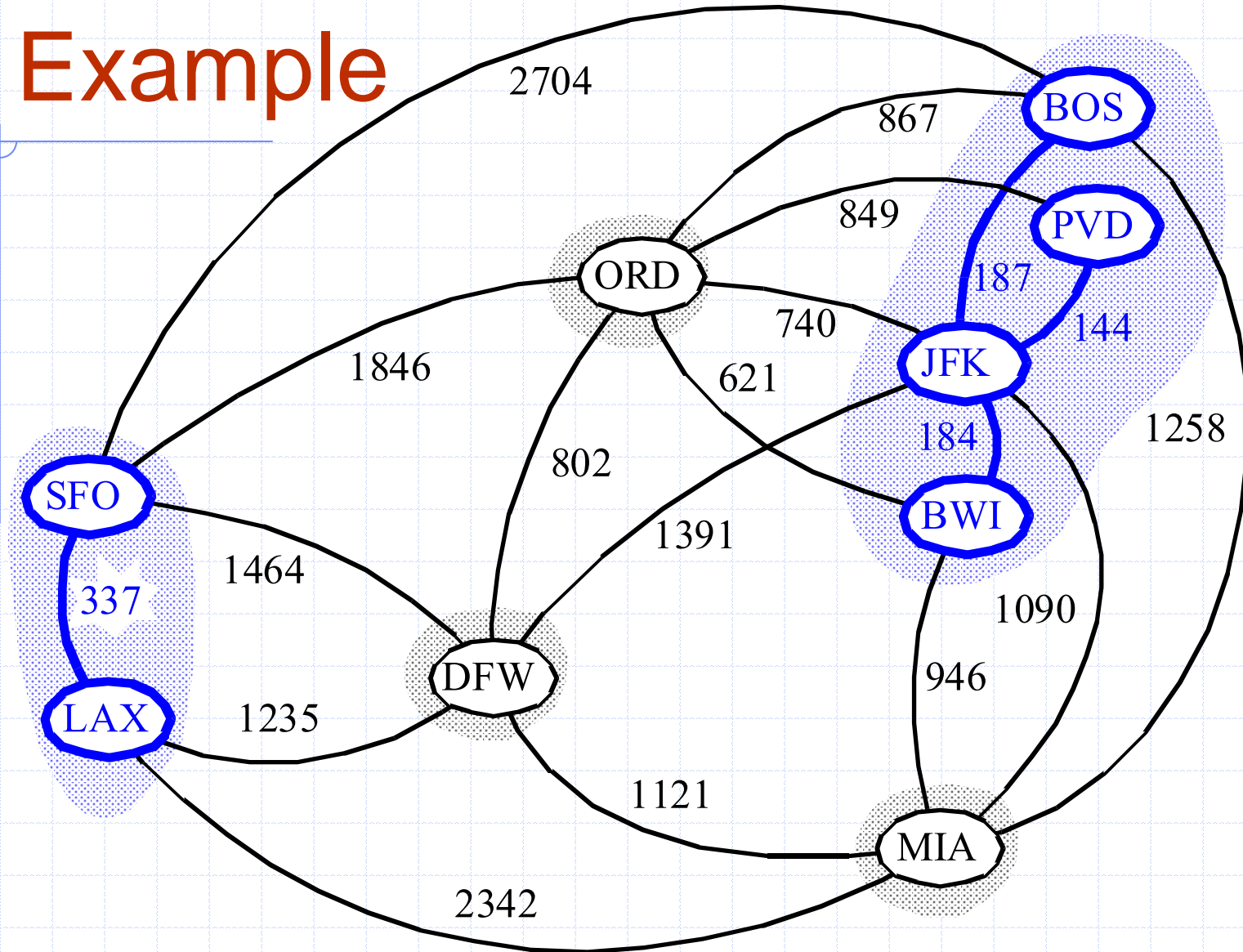
# Example



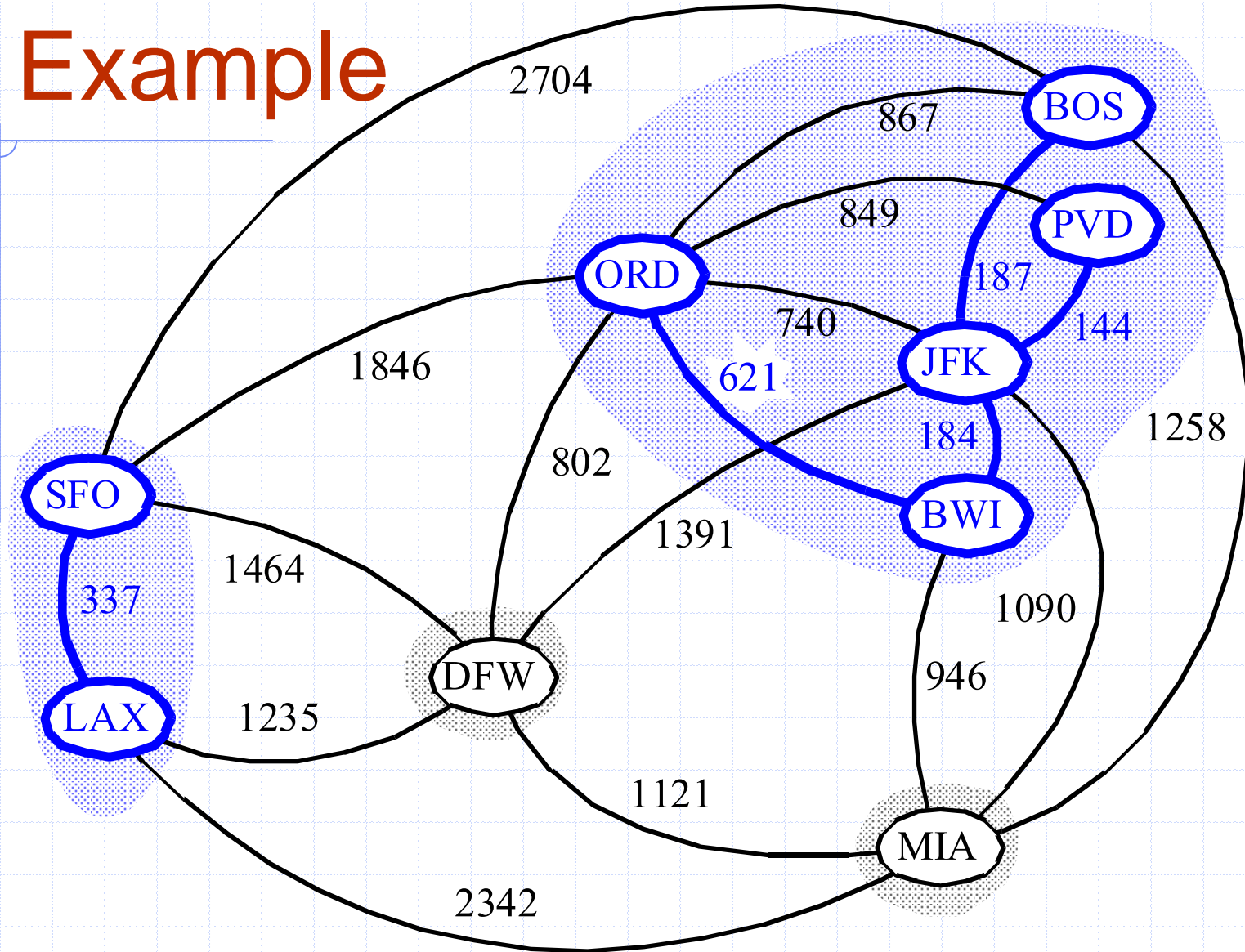
# Example



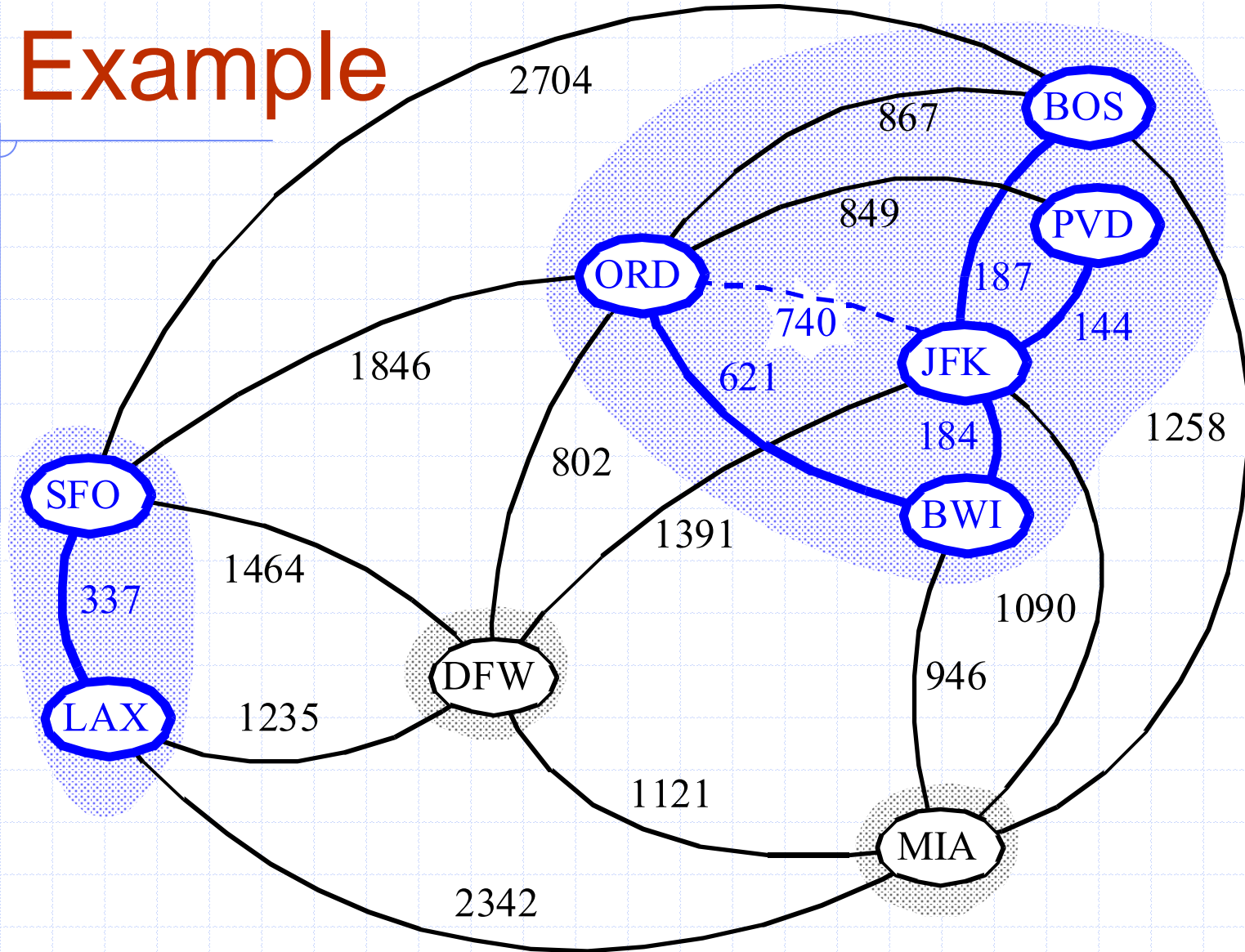
# Example



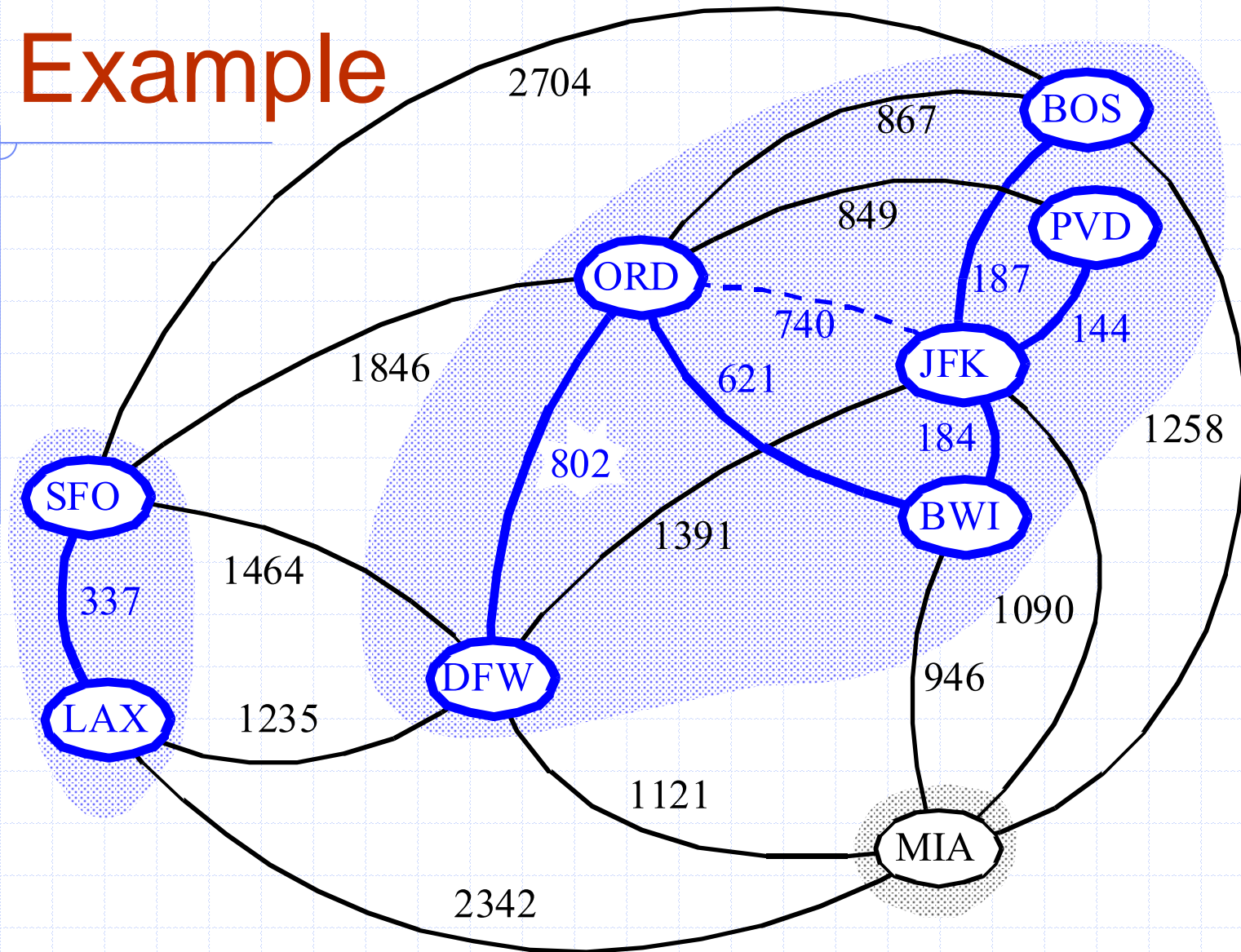
# Example



# Example

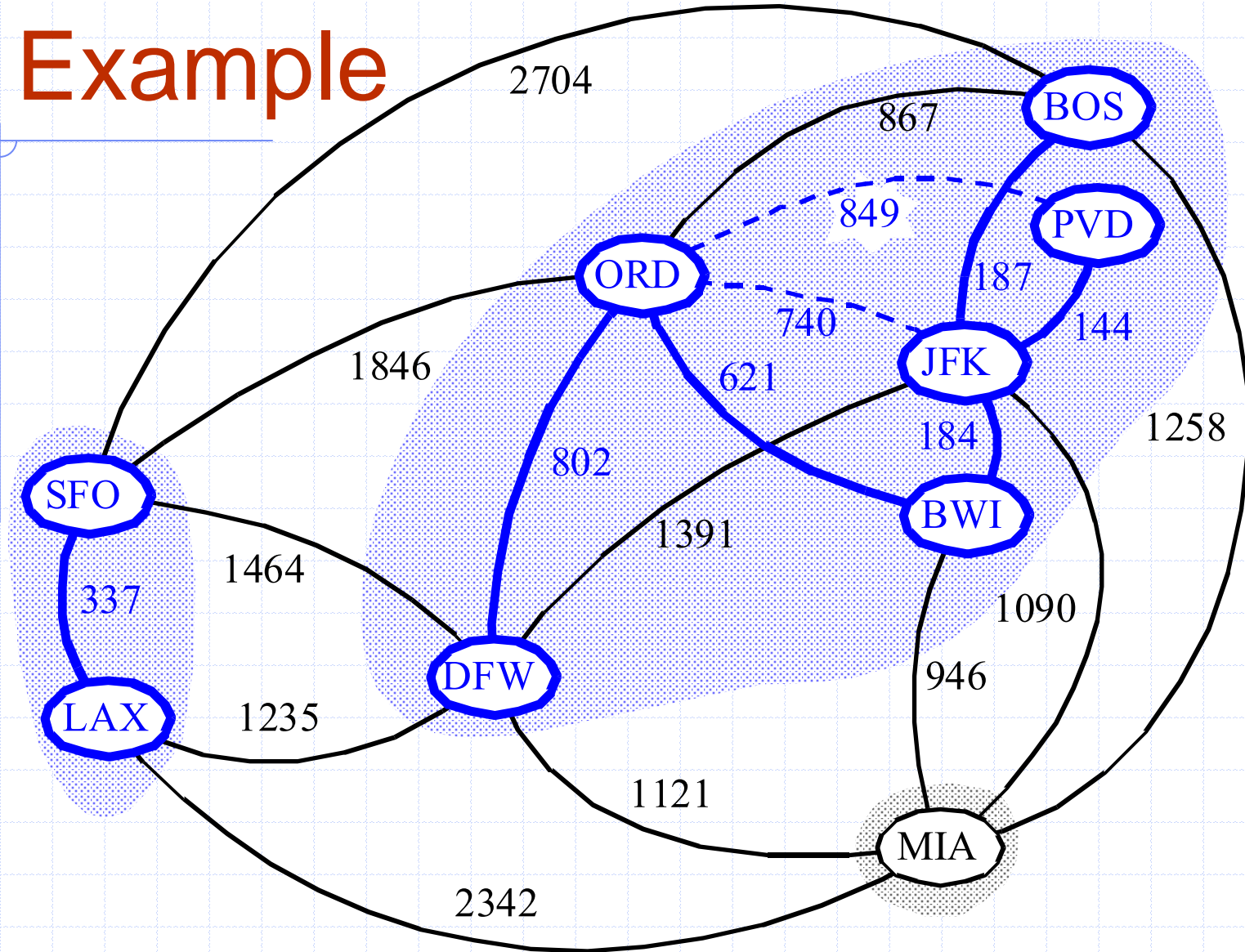


# Example

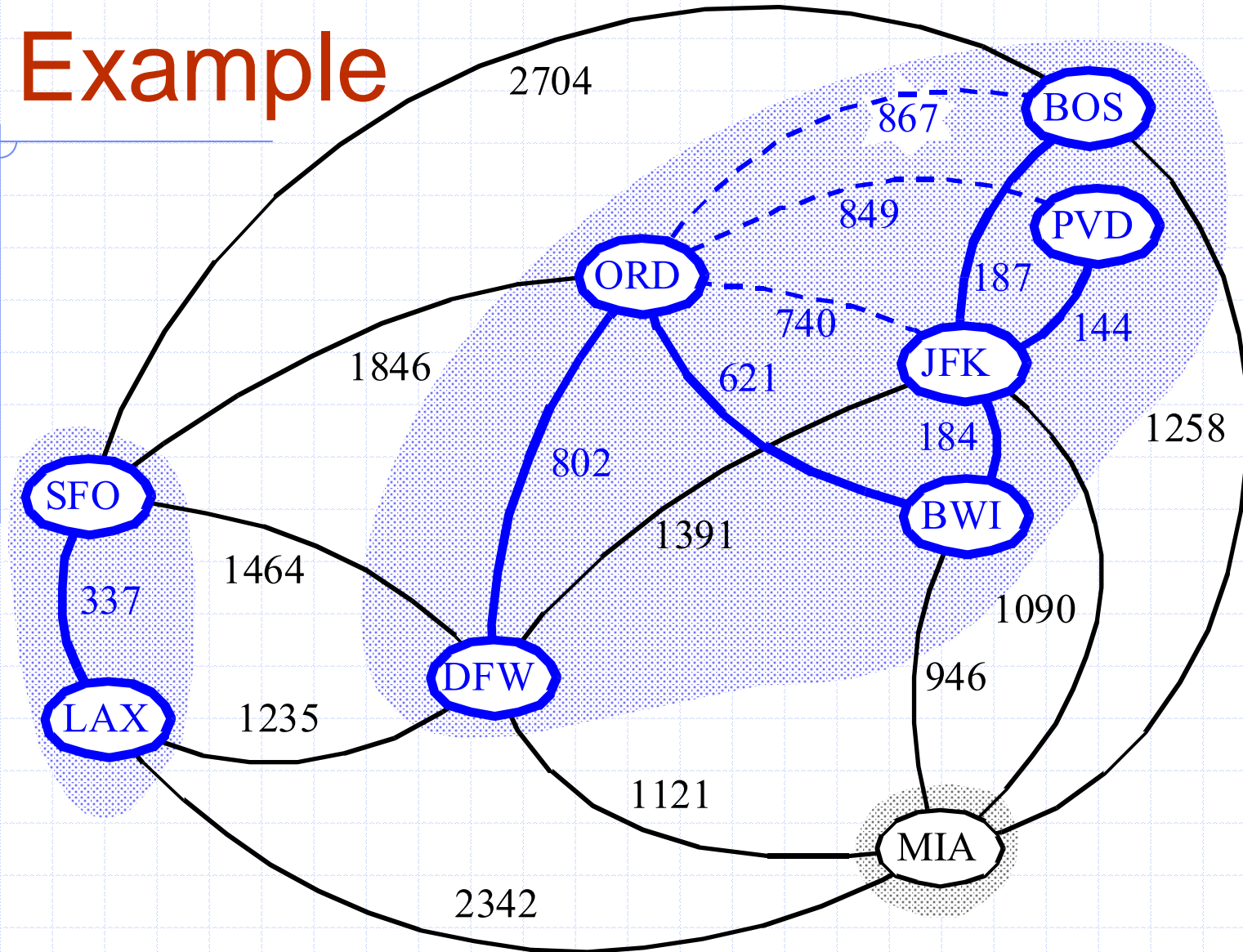




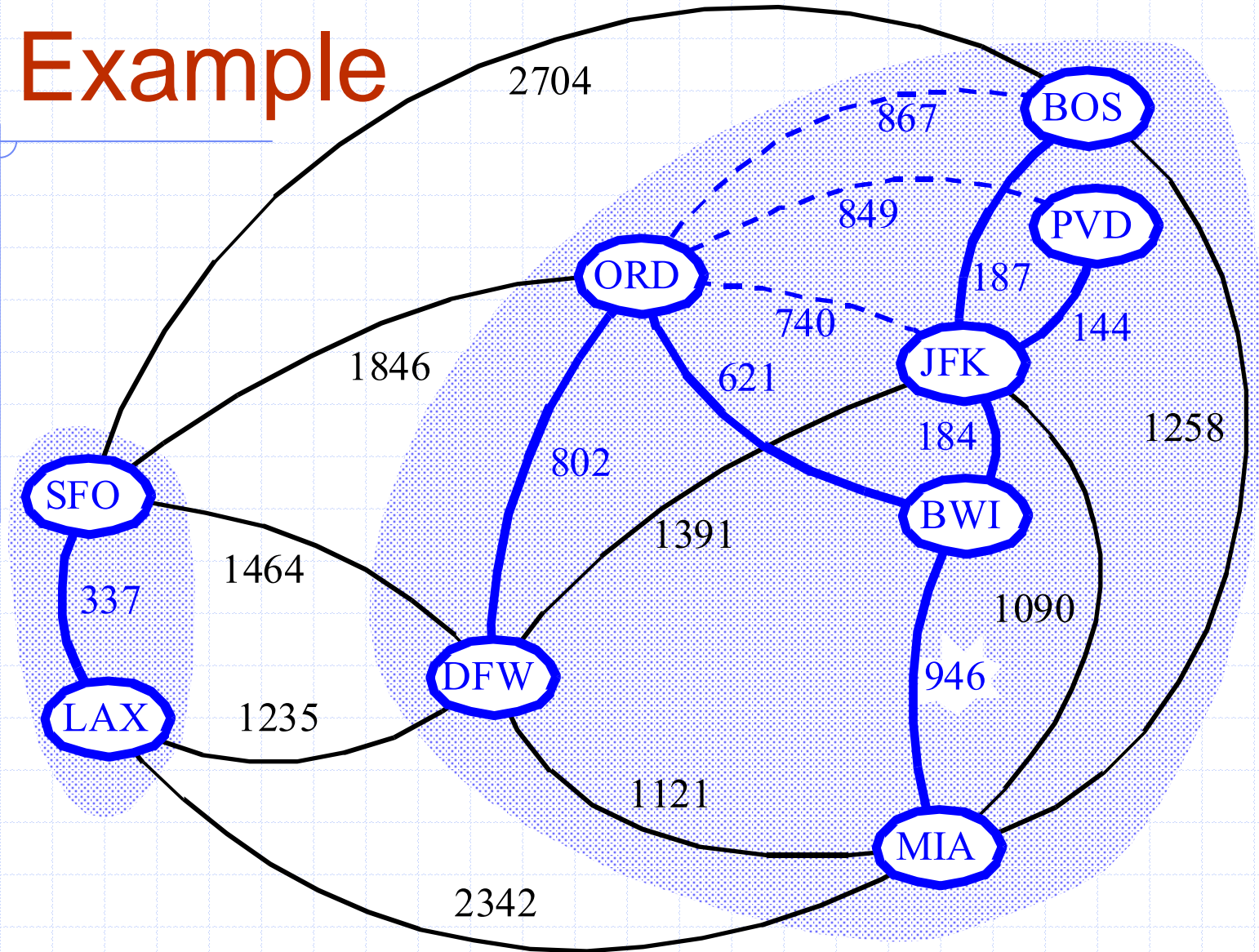
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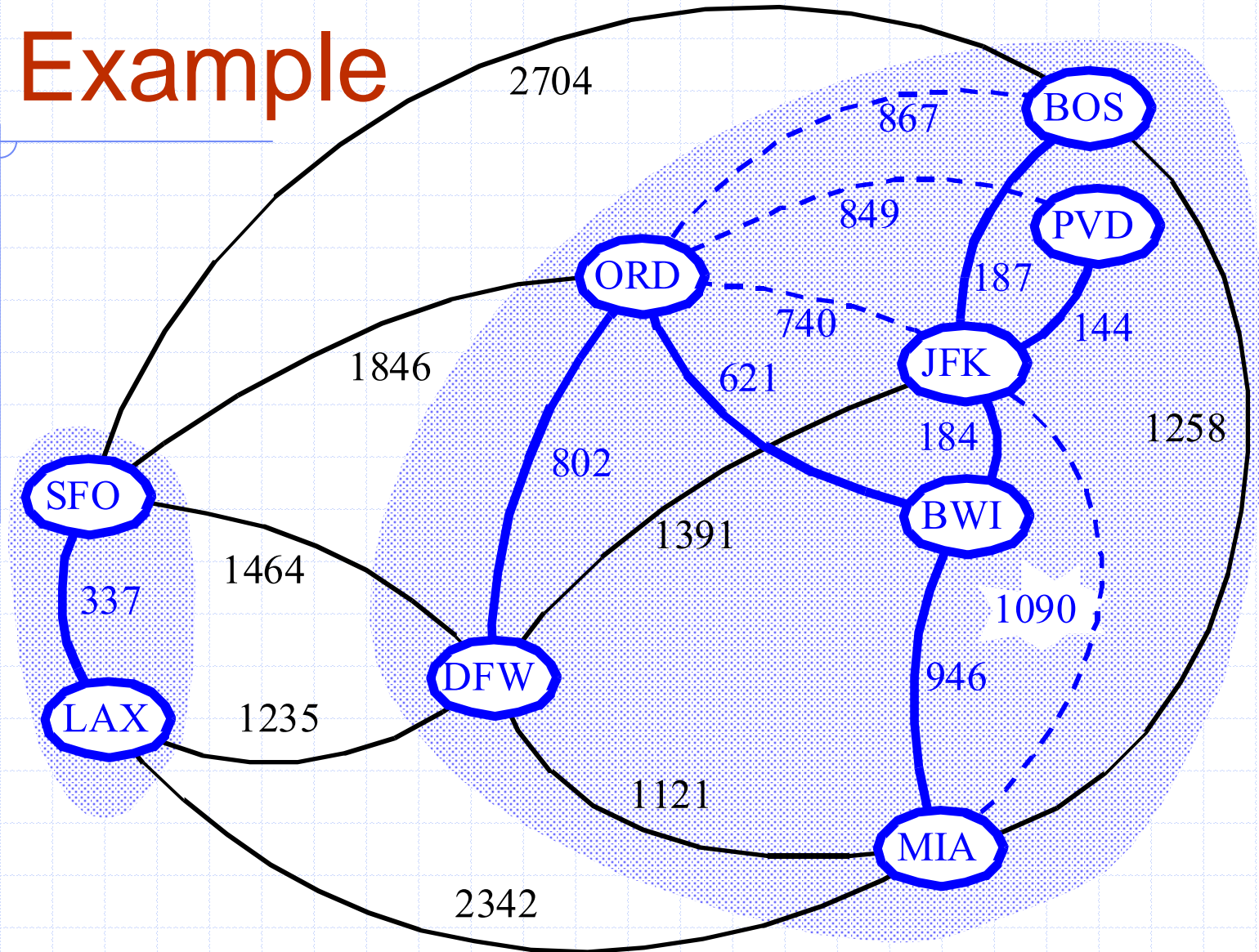
# Example



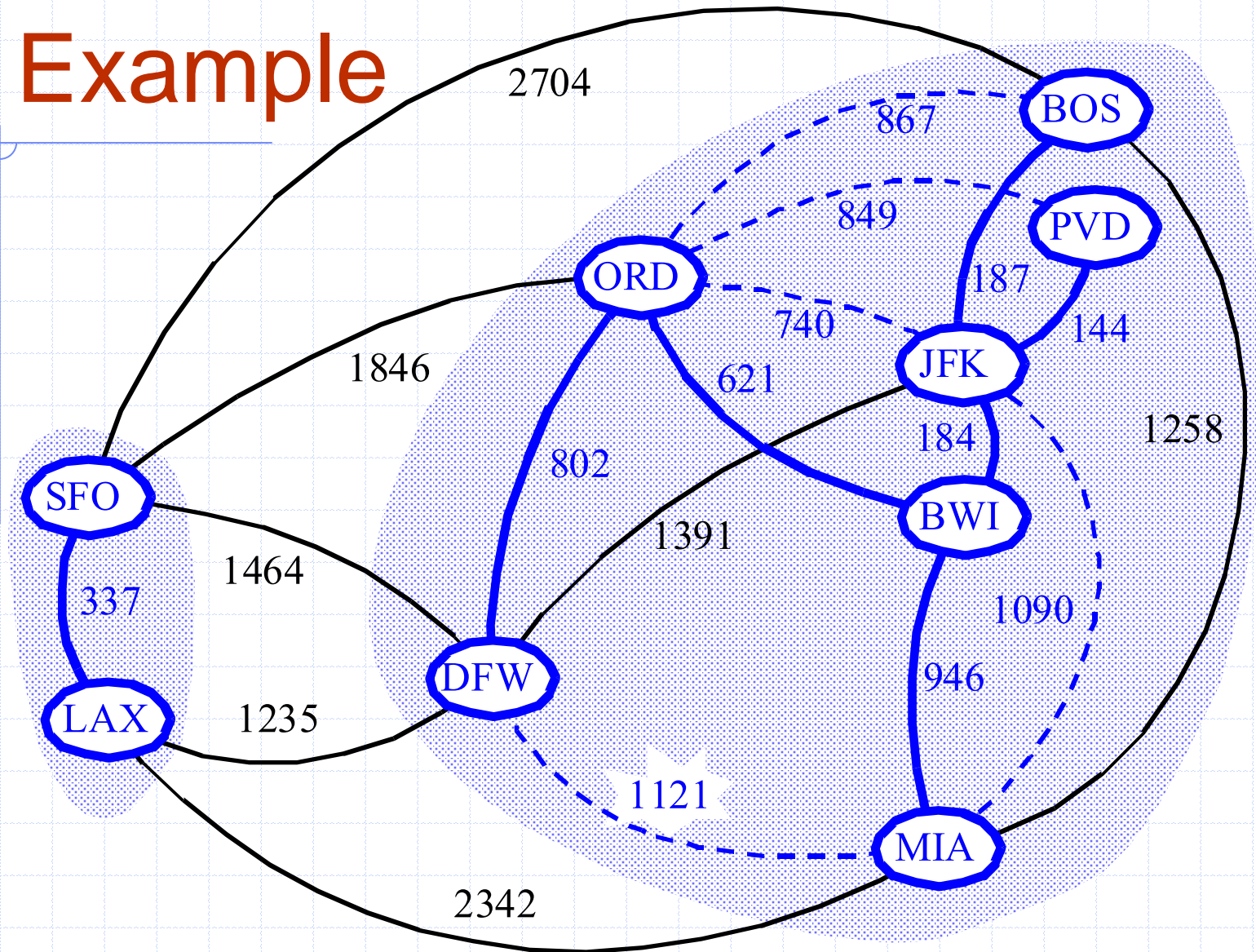
# Example



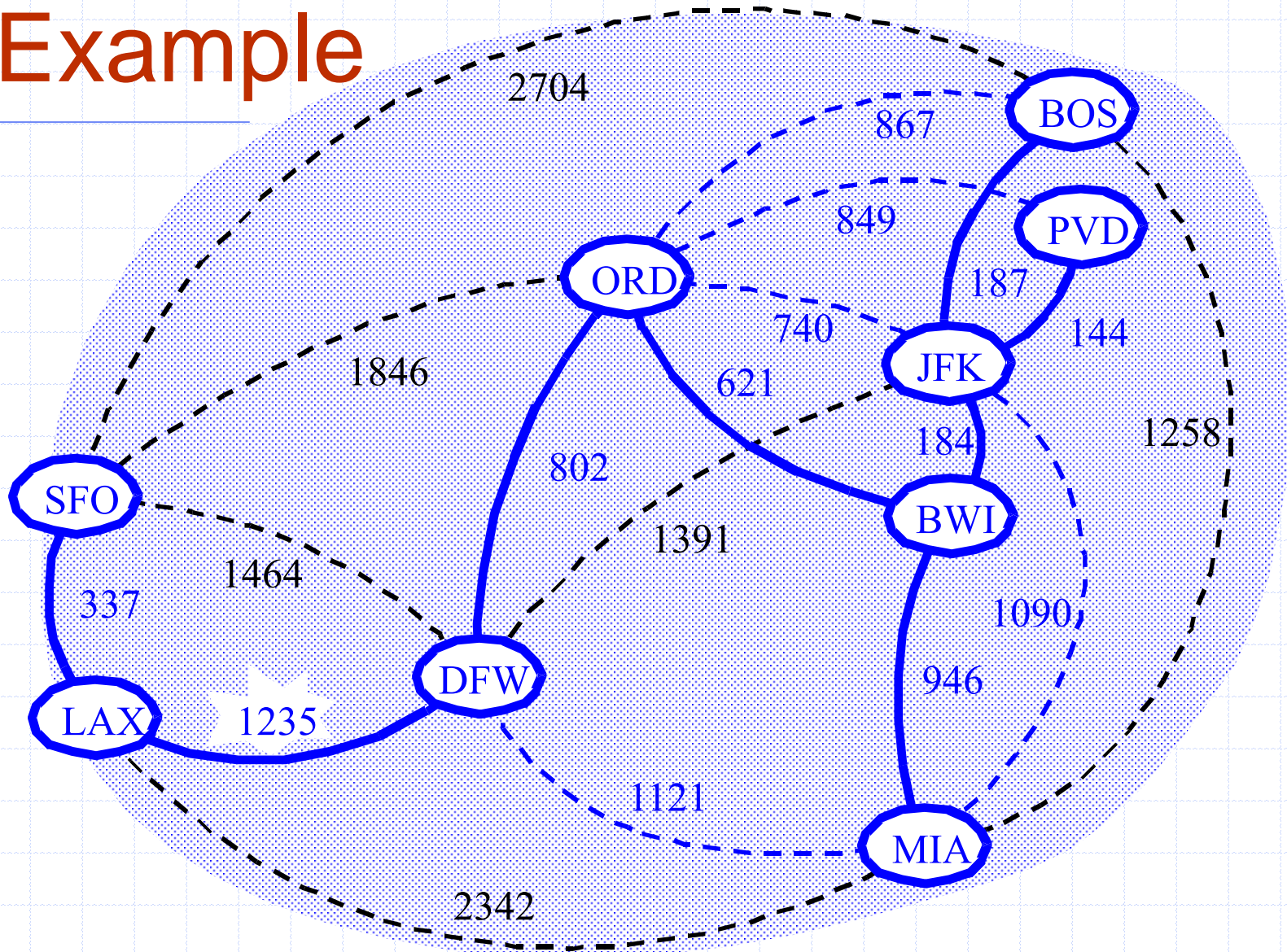
# Example



# Example



# Example



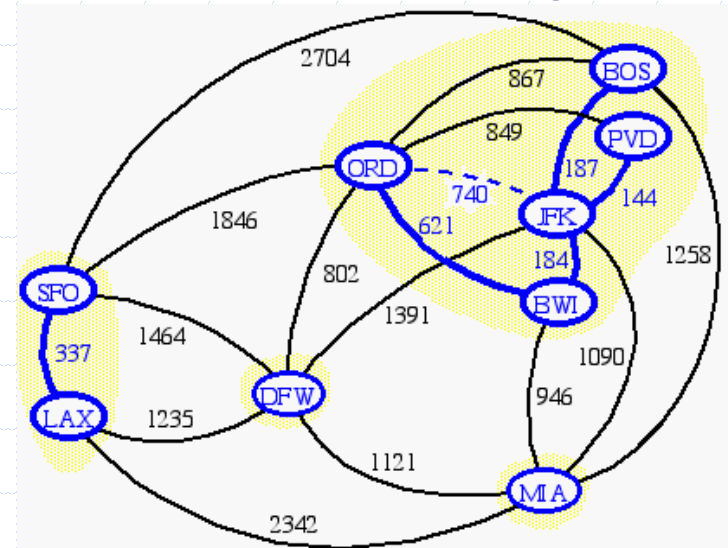
# Kruskal's Algorithm (1956) (High Level)

- ◆ A priority queue stores the edges outside the cloud
  - Key: weight
  - Element: edge
- ◆ At the end of the algorithm
  - We are left with one cloud that encompasses the MST
  - A tree  $T$  which is our MST

```
Algorithm KruskalMST( $G$ )  
  for each vertex  $v$  in  $G$  do  
    define a  $Cloud(v) \leftarrow \{v\}$   
   $Q \leftarrow$  new heap-based priority queue.  
  for all  $e \in G.edges()$   
     $Q.insert(weight(e), e)$   
   $T \leftarrow \emptyset$   
  while  $T$  has fewer than  $n-1$  edges do  
     $e \leftarrow Q.removeMin()$   
     $(u, v) \leftarrow G.endVertices(e)$   
    if  $Cloud(v) \neq Cloud(u)$  then  
      Add edge  $e$  to  $T$   
      Merge  $Cloud(v)$  and  $Cloud(u)$   
  return  $T$ 
```

# Data Structure for Kruskal Algorithm

- ◆ The algorithm maintains a forest of trees/clouds
- ◆ An edge is accepted if it connects distinct trees/clouds
- ◆ We need a data structure that maintains **clouds**, e.g., an array of disjoint sets of vertices (clouds), with the operations:
  - **getCloud**(u): returns the index of the cloud containing u (stored at the vertex as an integer attribute)
  - **mergeClouds**(u,v): merge the smaller cloud into the larger cloud to minimize time complexity





# Representation of a Cloud

- ◆ Each cloud/partition is stored in a sequence
- ◆ Each vertex has a reference back to the cloud containing it, which is an integer index into the **Clouds** array
  - operation **getCloud**(u) takes  $O(1)$  time, and returns the index of the cloud containing u.
  - in operation **mergeClouds**(u,v), we move the elements of the smaller cloud to the larger cloud and update the cloud attribute of vertices in the smaller cloud
  - the time for operation **mergeClouds**(u,v) is  $\text{MIN}(n_u, n_v)$ , where  $n_u$  and  $n_v$  are the sizes of the sequences storing u and v
- ◆ For each edge we do two **getCloud** operations
- ◆ We insert  $n-1$  edges into the MST, so we do at most  $O(n)$  merges
- ◆ Each element is processed (copied into a different cloud) at most  $\log n$  times (Why?)
  - Because whenever a vertex is processed in a union, it goes into a set of size at least double (Why?)
- ◆  $O(\log m)$  is  $O(\log n)$  (Why?)
  - Because  $m \leq n(n-1)/2 < n^2$

# Partition-Based Implementation

Uses an array to keep track of clouds and the cloud # is the index into the array of clouds.

**Algorithm Kruskal( $G$ ):**

**Input:** A weighted, simple, connected graph  $G$ .

**Output:** A MST  $T$  for  $G$  ( $T$  is a sequence containing the edges in the MST).

$Clouds \leftarrow$  new array of size  $n$

$c \leftarrow 0$

**for all**  $v \in G.vertices()$  **do**

$setCloudNum(v, c)$       // all start out in separate clouds

$Clouds[c] \leftarrow \{v\}$       // sequence containing  $v$  is stored in  $Clouds$  array

$c \leftarrow c + 1$

$Q \leftarrow$  new heap-based priority queue

**for all**  $e \in G.edges()$  **do**

$Q.insert(weight(e), e)$       //  $O(m \log m)$  which is  $O(m \log n)$

$T \leftarrow$  new Sequence      // edges in the MST

**while**  $T.size() < n-1$  **do**

$e \leftarrow Q.removeMin()$

$(u, v) \leftarrow G.endVertices(e)$

**if**  $getCloudNum(u) \neq getCloudNum(v)$  **then**

$T.insertLast(e)$

$mergeClouds(u, v, Clouds)$  {  $O(n \log n)$  since a vertex is merged  $O(\log n)$  times }

**return**  $T$

Running time:  
 $O((n+m)\log n)$

# Merge Two Clouds in time $O(\text{size of smaller cloud})$

**Algorithm** *mergeClouds*( $u, v, \text{Clouds}$ )

**Input:** Vertices  $u$  and  $v$  and array *Clouds* containing all clouds that still exist (or are non-empty).

**Output:** merges the two clouds containing  $u$  and  $v$

$cu \leftarrow \text{getCloudNum}(u)$

$cv \leftarrow \text{getCloudNum}(v)$

**if** *Clouds*[ $cu$ ].size()  $\geq$  *Clouds*[ $cv$ ].size() **then**

$\text{larger} \leftarrow \text{Clouds}[cu]$ ;  $\text{smaller} \leftarrow \text{Clouds}[cv]$

$\text{newCloud} \leftarrow cu$       // cloud # of larger

**else**

$\text{larger} \leftarrow \text{Clouds}[cv]$ ;  $\text{smaller} \leftarrow \text{Clouds}[cu]$

$\text{newCloud} \leftarrow cv$       // cloud # of larger

**while**  $\text{smaller.size()} > 0$  **do** // move vertices from smaller to larger cloud

$p \leftarrow \text{smaller.first}()$

$v \leftarrow \text{smaller.remove}(p)$       // smaller cloud will be empty

$\text{setCloudNum}(v, \text{newCloud})$       // set to cloud # of larger cloud

$\text{larger.insertLast}(v)$

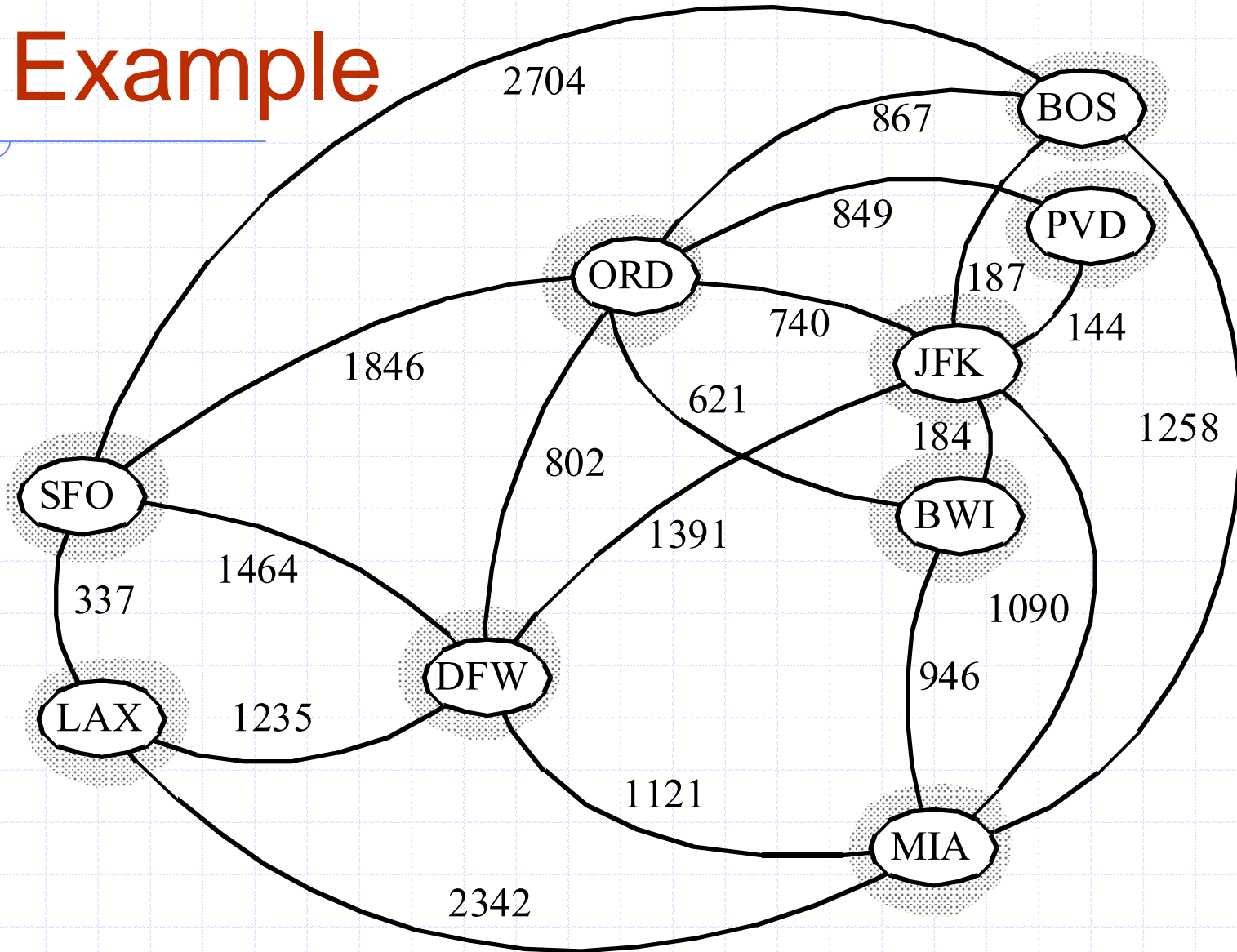
# Main Point

3. Kruskal's minimum spanning tree algorithm first finds the shortest edge, then tests it for safety; if it passes, it becomes part of the proposed solution.

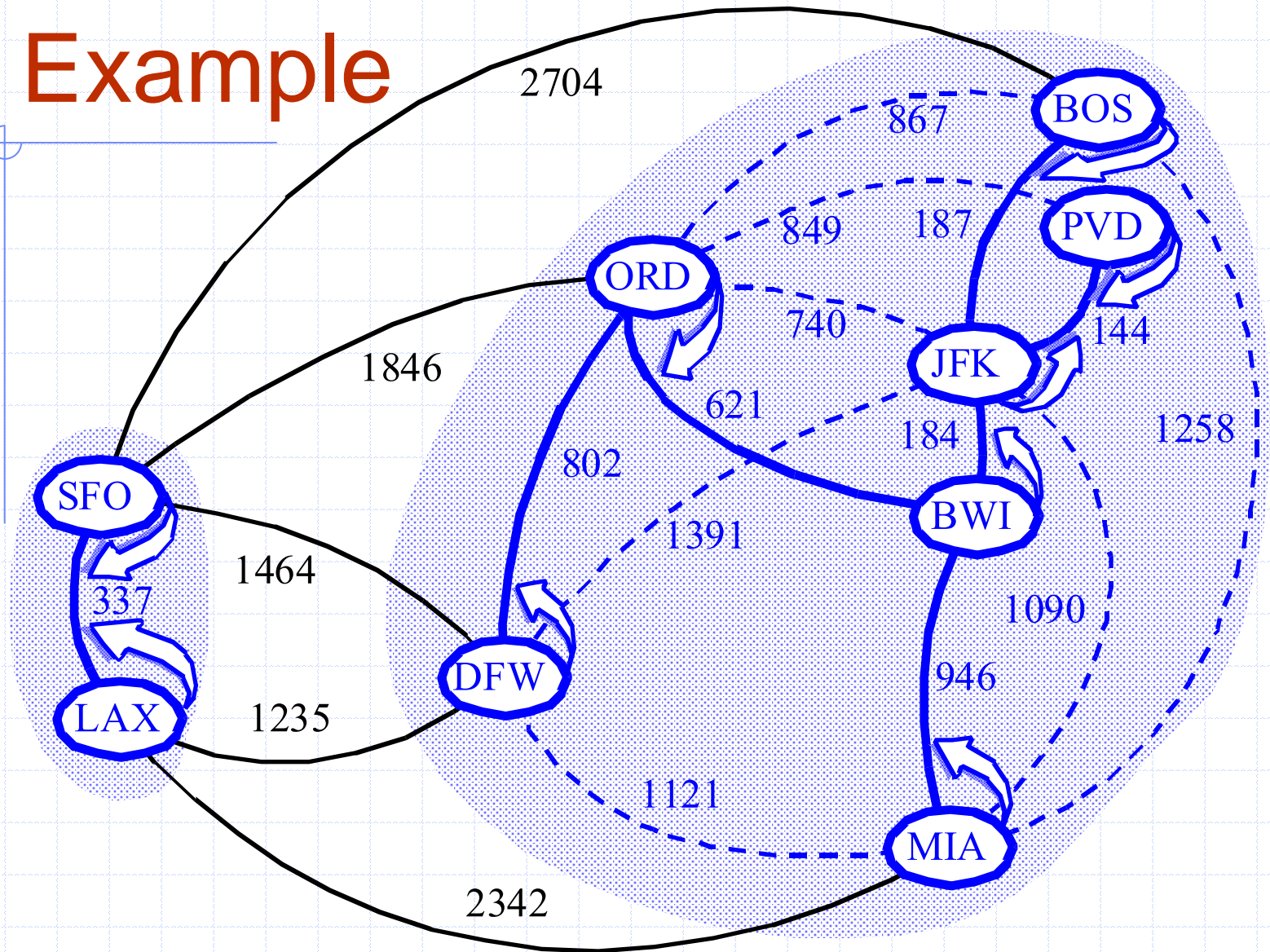
The fruits of an action cannot be determined precisely on the level of waking consciousness; it can only be determined from the level of infinite correlation where everything is interconnected. Thus only by establishing our awareness in pure awareness is right action possible.

# Baruvka's Algorithm (1926)

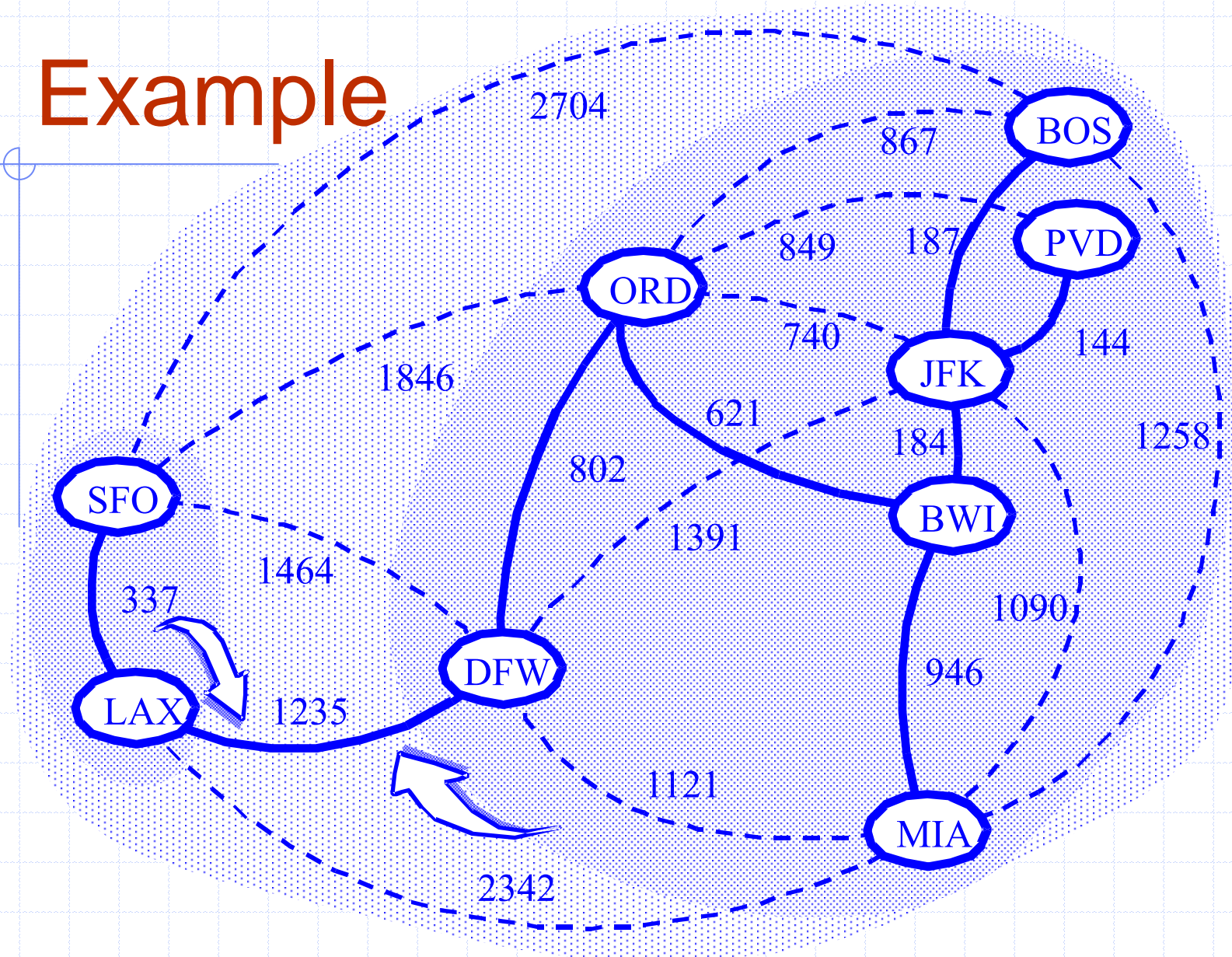
# Baruvka Example



# Example



# Example





# Baruvka's Algorithm

- ◆ Like Kruskal's Algorithm, Baruvka's algorithm grows many “clouds” at once.

## Algorithm *BaruvkaMST(G)*

```
 $T \leftarrow V$  {just the vertices of  $G$ , no edges,  $n$  connected components}  
while  $T$  has fewer than  $n-1$  edges do { $T$  is not yet an MST}  
  for each connected component  $C$  in  $T$  do  
    Find edge  $e$  with smallest-weight edge from  $C$  to  
      another component in  $T$ .  
    if  $e$  is not already in  $T$  then  
      Add edge  $e$  to  $T$   
return  $T$ 
```

- ◆ Each iteration of the while-loop halves the number of connected components in  $T$ .

# Baruvka's Algorithm (more details)

Algorithm *BaruvkaMST(G)*

for each  $e \in G.edges()$  do {label edges *NOT\_IN\_MST*}  
    set *MSTLabel*( $e$ , *NOT\_IN\_MST*) {no edges in MST}

$numEdges \leftarrow 0$

while  $numEdges < n-1$  do

    labelVerticesOfEachComponent( $G$ ) {BFS}

    insertSmallest-WeightEdgeOutOfComponents( $G$ )

return  $G$

# Required functionality

- ◆ Does not use a priority queue or locators!
- ◆ Does not use union-find data structures!
- ◆ Maintains a forest  $T$  subject to edge insertion
  - Can be supported in  $O(1)$  time using labels on edges in MST
- ◆ Step 1: Mark vertices with number of the component to which they belong
  - Traverse forest  $T$  to identify connected components
    - ◆  $O(1)$  time to label each vertex
    - ◆ Requires extra instance variable for each vertex
  - Takes  $O(n+m)$  time using a DFS or a BFS each time through the while-loop
- ◆ Step 2: Find a smallest-weight edge in  $E$  incident on each component  $C$  in  $T$  (insert into MST)
  - Scan edges to find the minimum weight edge incident on one of the vertices in  $C$  and incident on another vertex not in  $C$
  - Takes  $O(m)$  each time through the for-loop

# Analysis of Baruvka's Algorithm

- ◆ While-loop: each iteration (at worst) halves the number of connected components in  $T$ 
  - Thus executed  $\log n$  times
- ◆ Identifying connected components (in for-loop)
  - Vertices are labelled with component number
  - DFS or BFS of  $T$  runs in  $O(m+n)$  time (Skip edges of  $G$  that are not in  $T$ )
- ◆ For each component  $C$  find smallest edge connecting  $C$  to a different component of  $T$ 
  - Scan edges in  $G$
  - $O(m)$  time
- ◆ The running time is  $O(m \log n)$ .

# Output of the three MST algorithms is different

- ◆ Prim-Jarnik Algorithm (1957, 1930)
  - The edges in the MST are stored at the vertices by `setParent(e)`
- ◆ Kruskal's Algorithm (1956)
  - The MST edges are returned in a Sequence T
- ◆ Baruvka's Algorithm (1926)
  - The MST edges are labelled `IN_MST`

# After labeling each vertex with its component number (HW13)

## Algorithm *DFS*(*G*)

**Input** graph *G*

**Output** the edges of *G* are labeled as discovery edges and back edges

```
initResult(G)
for all u ∈ G.vertices()
    setLabel(u, UNEXPLORED)
    postVertexInit(G, u)
for all e ∈ G.edges()
    setLabel(e, UNEXPLORED)
    postEdgeInit(G, e)
for all v ∈ G.vertices()
    if getLabel(v) = UNEXPLORED
        preComponentVisit(G, v)
        DFScomponent(G, v)
        postComponentVisit(G, v)

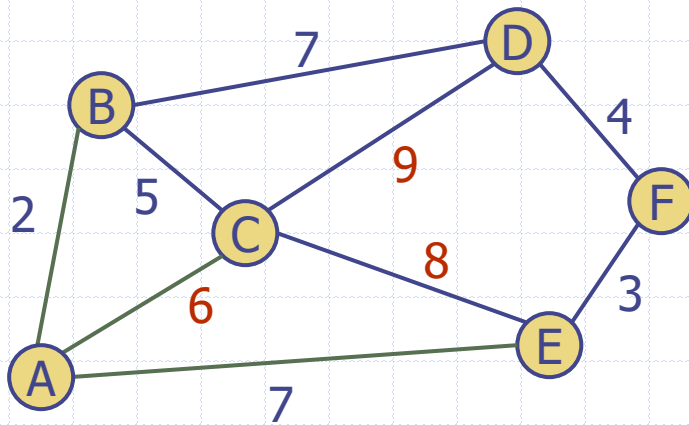
result(G)
```

## Algorithm *DFScomponent*(*G*, *v*)

```
setLabel(v, VISITED)
startVertexVisit(G, v)
for all e ∈ G.incidentEdges(v)
    preEdgeVisit(G, v, e, w)
    if getLabel(e) = UNEXPLORED
        w ← opposite(v, e)
        edgeVisit(G, v, e, w)
        if getLabel(w) = UNEXPLORED
            setLabel(e, DISCOVERY)
            preDiscoveryVisit(G, v, e, w)
            DFScomponent(G, w)
            postDiscoveryVisit(G, v, e, w)
        else
            setLabel(e, BACK)
            backEdgeVisit(G, v, e, w)
    finishVertexVisit(G, v)
```

HW14: Insert into *T* the smallest-weight edge going out from each component

# Cycle Property



By the Cycle Property, AC, CD, and CE cannot be in a MST.  
What about AE and BD?

Let's run the algorithms to verify this.

# Lower Bound on MST Computation

- ◆ There are randomized algorithms that compute MST's in expected linear time
- ◆ Linear time seems to be the lower bound
- ◆ Unknown whether there is a deterministic algorithm that runs in linear time (open question)



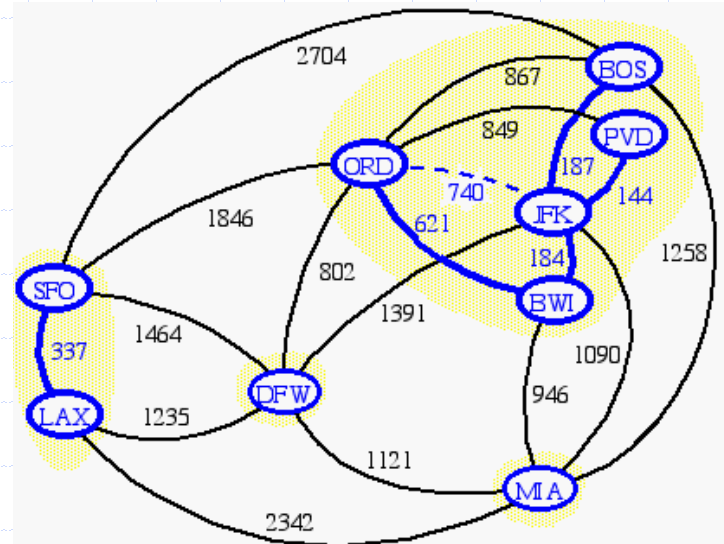
# Connecting the Parts of Knowledge with the Wholeness of Knowledge

1. Finding the minimum spanning tree can be done by an exhaustive search of all possible spanning trees, then choosing the one with minimum weight (but that takes at least exponential time).
2. To devise a greedy strategy, we identify a set of candidate choices, determine a selection procedure, and consider whether there is a feasibility problem. Then we have to prove that the strategy works.

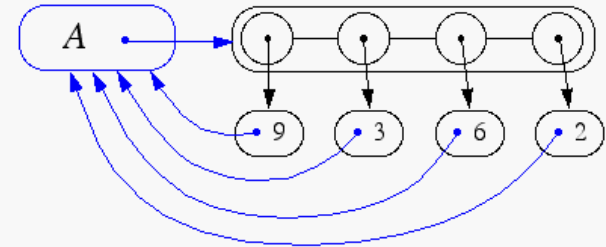
3. **Transcendental Consciousness** is the home of all the laws of nature, the source of all algorithms.
4. **Impulses within Transcendental Consciousness**: The natural laws within this unbounded field are the algorithms of nature governing all the activities of the universe.
5. **Wholeness moving within itself**: In Unity Consciousness, we perceive the spanning tree of natural law and appreciate the unity of all creation.

# Union-Find Data Structure for Kruskal Algorithm

- ◆ The algorithm maintains a forest of trees
- ◆ An edge is accepted if it connects distinct trees
- ◆ We need a data structure that maintains a **partition**, i.e., a collection of disjoint sets, with the operations:
  - find**(u): return the set storing u
  - union**(u,v): replace the sets storing u and v with their union



# Representation of a Partition



- ◆ Each set is stored in a sequence
- ◆ Each element has a reference back to the set
  - operation **find**(u) takes  $O(1)$  time, and returns the set of which u is a member.
  - in operation **union**(u,v), we move the elements of the smaller set to the sequence of the larger set and update their references
  - the time for operation **union**(u,v) is  $\min(n_u, n_v)$ , where  $n_u$  and  $n_v$  are the sizes of the sets storing u and v
- ◆ For each edge we do two finds
- ◆ We insert  $n-1$  edges into the MST, so we do at most  $O(n)$  unions
- ◆ Each element is processed (copied into a different cloud) at most  $\log n$  times (Why?)
  - Because whenever a vertex is processed in a union, it goes into a set of size at least double (Why?)
- ◆  $O(\log m)$  is  $O(\log n)$  (Why?)
  - Because  $m \leq n(n-1)/2 < n^2$

# Partition-Based Implementation

Performs cloud merges as unions and tests as finds.

**Algorithm Kruskal( $G$ ):**

**Input:** A weighted, simple, connected graph  $G$ .

**Output:** An MST  $T$  for  $G$ .

**for all**  $v \in G.vertices()$  **do**

    insert vertex  $v$  into  $T$

    define a *Cloud*( $v$ )  $\leftarrow \{v\}$

$Q \leftarrow$  new heap-based priority queue

**for all**  $e \in G.edges()$  **do**

$Q.insert(weight(e), e)$

{ $O(m \log m)$  which is  $O(m \log n)$ }

$T \leftarrow$  new empty tree

**while**  $T.numEdges() < n-1$  **do**

$e \leftarrow Q.removeMin()$

$(u, v) \leftarrow G.endVertices(e)$

**if**  $P.find(u) \neq P.find(v)$  **then**

        insert edge  $e$  into  $T$

$P.union(u, v)$

{ $O(n \log n)$  since a vertex is merged  $O(\log n)$  times}

**return**  $T$

Running time:  
 $O((n+m)\log n)$