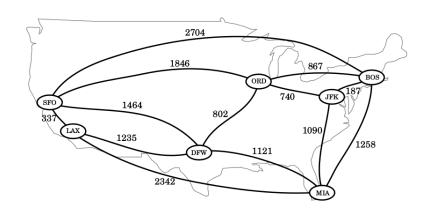
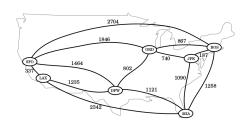
- The BFS search strategy can be used to find a shortest path from some starting vertex v to every other vertex in a connected graph
- This approach is useful when each edge is as good as any other, in particular it fails when a weighted graph has to be taken into account
- A weighted graph is a graph that has a numeric label w(e) associated with each edge e (the weight of that edge)
- For e = (u, v) we have w(u, v) = w(e)





Possible paths from JFK to LAX:

Path	Weights	Total
JFK - ORD - SFO - LAX	740 + 1846 + 337	2923
JFK - ORD - DFW - LAX	740 + 802 + 1235	2777
JFK - MIA - LAX	1090 + 2342	3432
JFK - MIA - DFW - LAX	1090 + 1121 + 1235	3446
JFK - BOS - ORD - SFO - LAX	187 + 867 + 1846 + 337	3237
JFK - BOS - ORD - DFW - LAX	187 + 867 + 802 + 1235	3091

The weight (or length) of a path P in a weighted graph G is the sum of the weights of the edges of P.

If
$$P = ((v_0, v_1), (v_1, v_2), \dots, (v_{k-1}, v_k))$$
 then

$$w(P) = \sum_{i=0}^{k-1} w(v_i, v_{i+1})$$

The distance from a vertex u to a vertex v in G, denoted d(u, v) is the length of a minimum-length path (shortest path) from u to v (if such a path exists)

Shortest Paths Dijkstra's Algorithm

- A class of algorithms solve the problem of finding a shortest path from some vertex s to each other vertex in a weighted graph G
- One of these, Dijkstra's algorithm, applies the greedy method: a given problem is solved by repeatedly selecting the best choice from among those available at each iteration
- The main idea is to perform a weighted BFS search starting from the source vertex s and creating a "cloud" of vertices, each entering the cloud in order of their distance from s
- Then, in each iteration, the next vertex chosen is the vertex outside the cloud closest to s
- The algorithm terminates when no more vertices are outside the cloud (or when those outside the cloud are not connected to those within the cloud)

Edge Relaxation

- Let us define a label D[v] for each vertex v in V
- These labels will always store the length of the best path obtained so far from s to v
- Initially, D[s] = 0 and $D[v] = \infty$ for each $v \neq s$
- We define the set C (the "cloud" of vertices) as the empty set
- At each iteration, a vertex u is selected not in C with smallest D[u] label, and u is put in C

Edge Relaxation

- Once a new vertex u is pulled into C, the label D[v] of each vertex adjacent to u and outside of C is updated, to reflect the fact that may be a new and better way to reach v via u
- This update is known as relaxation, because it takes an old estimate and checks if it can be improved to get closer to its value

More precisely:

```
if D[u] + w(u,v) < D[v] then
D[v] = D[u] + w(u,v)</pre>
```

Edge Relaxation

```
Algorithm ShortestPath(G,s):
 D[s] = 0
 D[v] = infinite for each vertex v != s
 Let a priority queue Q contain all the vertices
 of G using the D labels as keys
 while Q is not empty do
   u = value returned by Q.remove_min()
    for each vertex v adjacent to u such that v is in Q
       do
      if D[u] + w(u,v) < D[v] then
       D[v] = D[u] + w(u,v)
        Change to D[v] the key of vertex v in Q
  return the label D[v] of each vertex v
```

Computational Cost

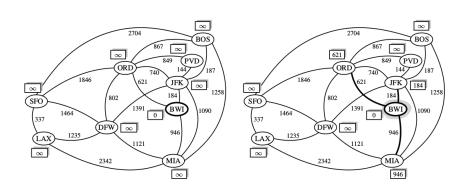
```
Let a priority queue Q contain all the vertices of G using the D labels as keys

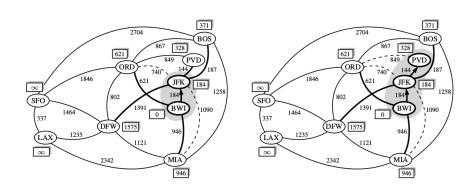
while Q is not empty do
    u = value returned by Q.remove_min()
    for each vertex v adjacent to u such that v is in Q do
    if D[u] + w(u,v) < D[v] then
        D[v] = D[u] + w(u,v)
        Change to D[v] the key of vertex v in Q

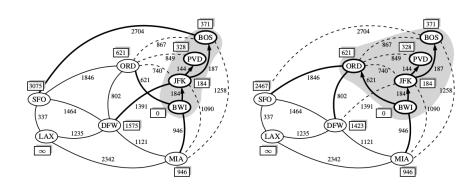
return the label D[v] of each vertex v
```

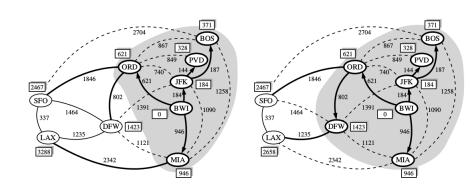
Computational Cost

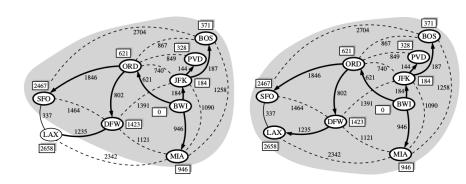
- The nested for loop runs in O(m) time
- The outer while loop executes in O(n) time
- The update of Q requires O(m) time
- If Q is a priority queue, each of the above operations run in O(logn), therefore the overall running time is O((n+m)logn), that is (approx) $O(n^2logn)$





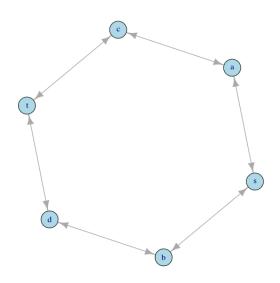


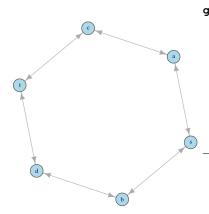




Dijkstra's Algorithm

Shortest Path Tree





```
def dijkstra(graph, src, dest, visited=[], distances={},
   predecessors={}):
    if src not in graph:
        raise TypeError('the root of the shortest path
            tree cannot be found in the graph')
    if dest not in graph:
        raise TypeError('the target of the shortest path
             cannot be found in the graph')
    if src == dest:
        path=[]
        pred=dest
        while pred != None:
            path.append(pred)
            pred=predecessors.get (pred, None)
        print('shortest path: '+str(path)+" cost="+str(
            distances[dest]))
    else :
        if not visited:
            distances[src]=0
```

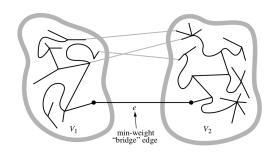
```
for neighbor in graph[src] :
  if neighbor not in visited:
  new_distance = distances[src] + graph[src][neighbor]
  if new distance < distances.get(neighbor,float('inf'))</pre>
    distances[neighbor] = new_distance
    predecessors[neighbor] = src
visited.append(src)
unvisited={}
for k in graph:
  if k not in visited:
    unvisited[k] = distances.get(k, float('inf'))
x=min(unvisited, key=unvisited.get)
dijkstra(graph, x, dest, visited, distances, predecessors)
```

- All the computers in an office must be connected using the least amount of cable (wireless is not admitted)
- This problem can be modeled using an undirected weighted graph G whose vertices represent the computers, and whose edges represent all the possible pairs (u, v) of computers
- The weight w(u, v) of the edge (u, v) is equal to the amount of cable needed to connect computer u to computer v
- The problem is finding a tree that contains all the vertices of G and has the minimum total weight over all such trees

Given an undirected weighted graph G, we are interested in finding a tree T that contains all the vertices in G and minimizes the sum

$$w(T) = \sum_{(u,v)inT} w(u,v)$$

Such a tree that contains every vertex of a connected graph G is said to be a **spanning tree**, and the problem of computing a spanning tree T with smallest total weight is known as the **minimum spanning tree** (MST) problem



Proposition

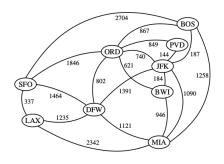
Let G be a weighted connected graph, and let V_1 and V_2 be a partition of the vertices of G into two disjoint nonempty sets. Let e be an edge in G with minimum weight from among those with one endpoint in V_1 and the other V_2 . There is a minimum spanning tree T that has e as one of its edges.

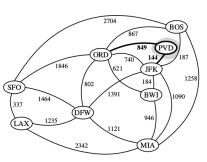
- In this algorithm, a minimum spanning tree is grown starting from a single cluster starting from some "root" vertex s
- The main idea is similar to that of Djkstra's algorithm: a cloud of vertices is defined that grows at each iteration
- At each iteration, a minimum-weight edge is chosen e=(u,v) which connects a vertex u in the cloud C to a vertex v outside the cloud C
- The vertex v is brought into the cloud C and the process is repeated until a spanning tree is formed
- As in Dijkstra'a algorithm, a label D[v] is maintained for each vertex outside the cloud C, so that D[v] stores the weight of the minimum observed edge for joining v to the cloud C

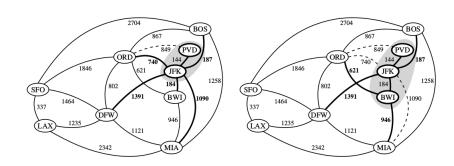
```
Algorithm Prim Jarnik(G):
  Input: An undirected, weighted, connected graph G
         with n vertices and m edges
 Output: A minimum spanning tree T for G
 Pick any vertex s of G
 D[s] = 0
  for each vertex v != s do
   D[v] = infinite
  Initialize T = 0
  Initialize a priority queue Q with an entry (D[v], (v,
      None)) for each vertex v, where D[v] is the key
     in the priority queue, and (v, None) is the
     associated value
 while Q is not empty do
    (u, e) = value returned by Q.remove_min()
   Connect vertex u to T using edge e
    for each edge e' = (u, v) such that v is in Q do
      if w(u,v) < D[v] do
```

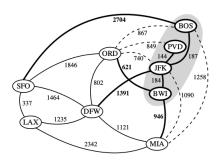
Analysis of the Prim-Jarnik Algorithm

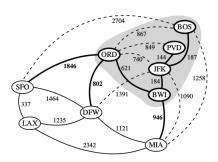
- n insertions are performed into Q
- Later, *n* extract-min are performed
- A total of *m* priorities are updates
- With a priority queue, each operation runs in O(logn), and the overall time for the algorithms is O((n+m)logn), which is O(mlogn) for a connected graph
- By using an unsorted list, the running time will be $O(n^2)$

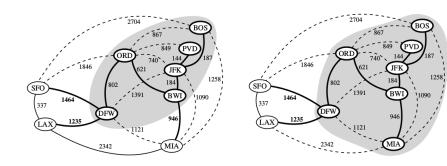


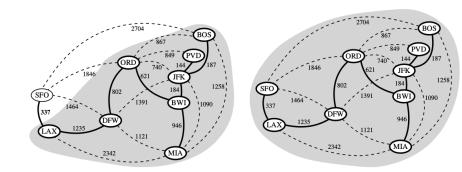












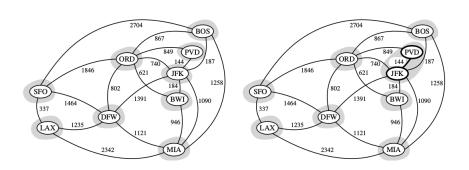
```
def MST_PrimJarnik(q):
 d = \{\}
 tree = []
 pq = AdaptableHeapPriorityQueue()
 pqlocator = {}
  for v in g.vertices():
    if len(d) == 0:
      d[v] = 0
    else.
      d[v] = float('inf')
    pqlocator[v] = pq.add(d[v], (v, None))
(to be continued)
```

```
while not pq.is_empty():
  key,value = pq.remove_min()
  u,edge = value
  del pglocator[u]
  if edge is not None:
    tree.append(edge)
  for link in q.incident_edges(u):
    v = link.opposite(u)
    if v in pqlocator:
      wgt = link.element()
      if wqt < d[v]:
        d[v] = wqt
        pq.update(pqlocator[v], d[v], (v, link))
return tree
```

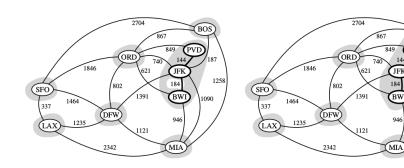
- Kruskal's algorithm maintains a forest of clusters, repeatedly merging pairs of clusters until a single cluster spans the graph
- Initially, each vertex is by itself a singleton cluster
- The algorithm considers each edge in turn, ordered by increasing weight
- If an edge e connects two different clusters, then e is added to the set of edges of the MST, and the clusters connected by e are merged into a single cluster
- On the contrary, if e connects two vertices that are already in the same cluster, then e is discarded
- The algorithms ends when enough edges have been added to form a spanning tree

```
Algorithm Kruskal(G):
  Input: A simple connected weighted graph G
         with n vertices and m edges
  Output: A minimum spanning tree T for G
  for each vertex v in G do
    Define an elementary cluster C(v) = \{v\}
    Initialize a priority queue Q to contain all edges
        in G, using the weights as keys
    T = 0
    while T has fewer than n-1 edges do
      (u, v) = value returned by Q.remove_min()
      Let C(u) be the cluster containing u, and let C(v)
          be the clustre containing v
      if C(u) != C(v) then
        Add edge (u,v) to T
        Merge C(u) and C(v) into one cluster
    return tree T
```

- The correctness of Kruskal's algorithm is based upon the crucial fact about minimum spanning trees from Proposition
- Each time that Kruskal's algorithm adds an edge (u, v) to the MST T, we can define a partitioning of the set of vertices V by letting V_1 be the cluster containing v and letting V_2 contain the rest of the vertices in V
- ullet This defines a disjoint partitioning of the vertices of V
- Moreover, since we are extracting edges from Q in order by their weights, e must be a minimum-weight edge with one vertex in V_1 and the other in V_2 .
- Therefore, Kruskal's algorithm always adds a valid MST edge

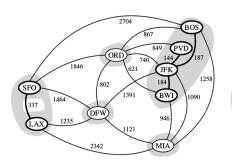


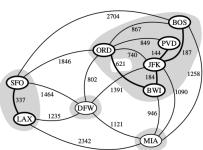
Kruskal's Algorithm

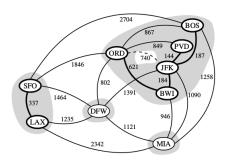


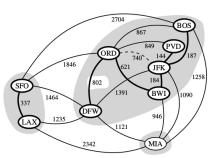
1258

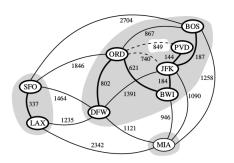
1090

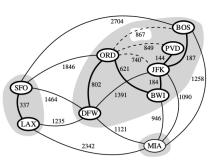


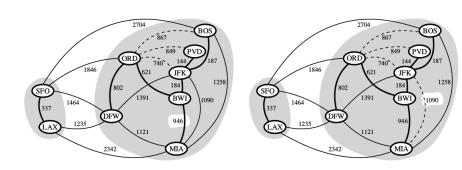


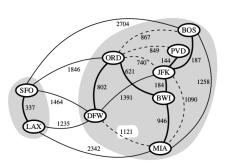


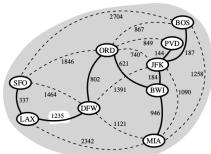












Analysis of Kruskal's Algorithm

- The ordering of edges can be implemented in O(mlogn) (by using a priority queue)
- To implement Kruskal's algorithm, we must be able to find the cluster for vertices u and v that are endpoints of an edge e, to test whether those two clusters are distinct, and if so, to merge those two cluster into one
- The management of **disjoint partitions** can be performed in O(m + nlogn)
- For a connected graph we have $m \ge n-1$ and therefore the dominant term is O(mlogn), which is the running time of Kruskal's algorithm

Kruskal's Algorithm

```
def MST_Kruskal(g):
  tree = []
  pq = HeapPriorityQueue()
  forest = Partition()
  position = {}
  for v in q.vertices():
    position[v] = forest.make_group(v)
  for e in q.edges():
    pq.add(e.element(), e)
  size = q.vertex_count()
  while len(tree) != size - 1 and not pq.is_empty():
    weight,edge = pq.remove_min()
    u, v = edge.endpoints()
    a = forest.find(position[u])
    b = forest.find(position[v])
    if a != b:
```

345 / 347

Disjoint Partitions

- A partition data structure manages an universe of elements that are organized into disjoint sets (an element belongs to one and only one of these sets)
- For clarity reason, the clusters of a partition are referred to as groups
- To differentiate between one group and another, we assume that at any point in time, each group has a designated entry called leader of the group

Disjoint Partitions

Method	Functionality
make_group(x)	Create a singleton group containing new element x and return the position storing x
union(p, q)	Merge the groups containing positions p and q
find(p)	Return the position of the leader of the group containing position p