

# Week 3 Greedy algorithms

A greedy algorithm is an algorithm that follows the problem solving approach of making a locally optimal choice at each stage with the hope of finding a global optimum

**Common Problems can be solved using greedy approach:**

- Interval scheduling/partitioning
- Scheduling to minimize lateness
- **Shortest path**
- **Minimum spanning trees**

**How to design greedy algorithm:**

Step1: Understand problem

Step2: Start with simple algorithm

Step3: Test it correctness:

- Try some simple examples to get feel for algorithm
- If none of them break algorithm, see if there's underlying structural property we can use to prove correctness

Step4: Better understanding algorithm, and improve

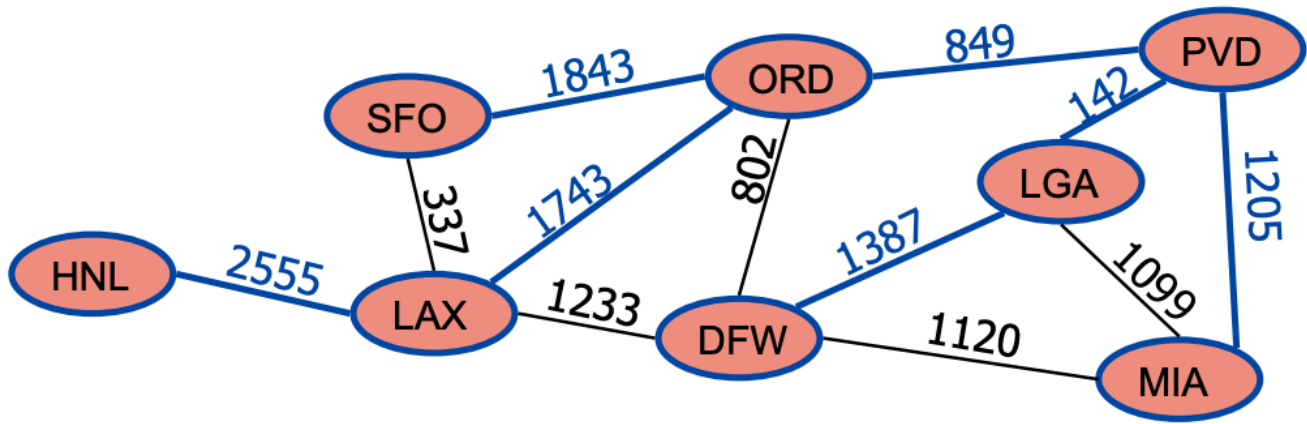
## Shortest path

**Problems:** Given an edge weighted graph and two vertices  $u$  and  $v$ , we want to find a path of minimum total weight between  $u$  and  $v$ , where the weight of a path is the sum of the weights of its edges.

**Applications:** Internet packet routing, flight reservations and driving directions

**Property:**

1. A subpath of a shortest path is itself a shortest path
2. There is a tree of shortest paths from a start vertex to all the other vertices (shortest path tree)



## Dijkstra's Algorithm 🤖

Maintain a set of explored nodes  $S$  for which we have determined the shortest path distance  $D(u)$  from  $s$  to  $u$ . ( $s$  is the starting vertex)

- Input:
  - Graph  $G = (V, E)$
  - Edges with non-negative weights
  - Start vertex  $s$
- Output:
  - Distance from  $s$  to all  $v \in V$
  - Shortest path tree rooted at  $s$
- Initialize  $S = \{s\}$ ,  $D(s) = 0$ ,  $D[v] = \text{infinite}$  for all  $v$  (vertices) in  $V - s$
- Repeatedly choose unexplored node  $v$  which minimizes:

$$D[v] = \min\{ D[u] + w(l_e) \}, \text{ for each } u \in S,$$

- $l_e$  is a single edge (neighbor edge) connected between  $v$  and  $u$
- $w(l_e)$  — the weight of edge  $l_e$
- $v$  can be any vertices not explored before (not in  $S$ )

## Pseudocode

```
def Dijkstra(G = (V, E), w, s):
    # initialize algorithm
    D = [] # shortest Distance list from vertex s to other vertices
    parent = {} # Represent the rooted tree generated by Dijkstra
    for v in V do
        D[v] = sys.maxint #set D[v] to infinity
    D[s] = 0
```

```

Q <- new priority queue for { (v, D[v]) : v in V}

#iteratively add vertices to S
while Q is not empty do
  u <- Q.remove_min()
  for z in G.neighbors(u) do
    if (D[u] + w[u,z]) < D[z] then
      D[z] <- D[u] + w[u, z]
      Q.update_priority(z, D[z])
      parent[z] <- u # set the z.parent as vertex u
return D, parent

```

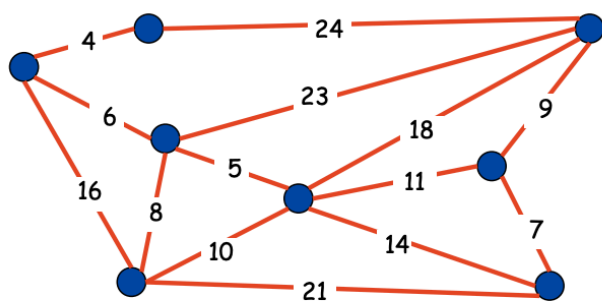
## Time Complexity

Assume the graph is connected:  $m \geq n - 1$ , the algorithm spends  $O(m)$  time on everything except PQ operations

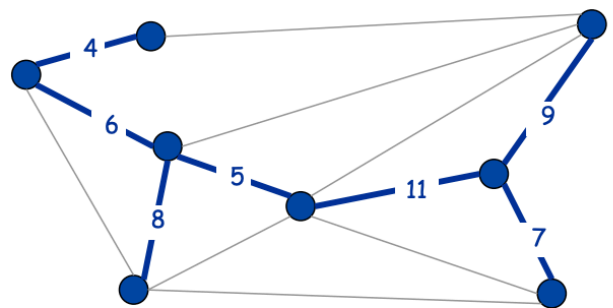
- But using normal **Heap PQ**, Disjksstar runs in  $O(m * \log n)$  time
  - Remove\_min() in each operation cost  $O(\log n)$ , and iterate  $m$  times
  - Using Fibonacci heaps, instead we get  $O(m + n \log n)$

## Minimum Spanning Tree

Given a connected graph  $G = (V, E)$  with real-valued edge weights  $c_e$ , an MST is a subset of the edges  $T \subseteq E$  such that  $T$  is a spanning tree whose sum of edge weights is minimized.



$G = (V, E)$

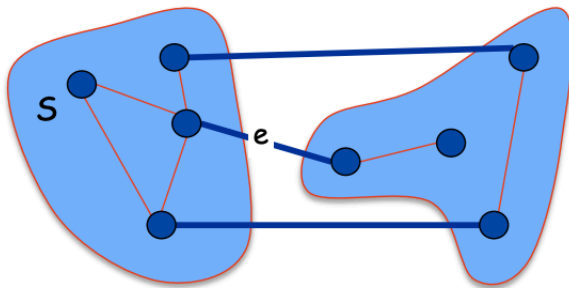


$T$  with  $\sum_{e \in T} c_e = 50$

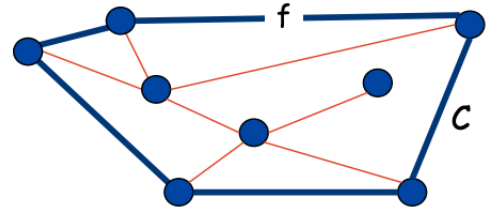
## MST Properties

**Assumption:** All edge costs  $c_e$  are distinct

- **Cut properties:** Let  $S$  be any subset of nodes, and let  $e$  be the min cost edge with exactly one endpoint in  $S$ . Then the MST contains  $e$
- **Cycle Property:** Let  $C$  be any cycle, and let  $f$  be the max cost edge belonging to  $C$ . Then the MST does not contain  $f$ .



$e$  is in the MST



$f$  is not in the MST

## Terminology

Cutset — A **Cut** is a subset of nodes  $S$ . The corresponding **cutset**  $D$  is the subset of edges with exactly one endpoint in  $S$ .

## Find the Minimum Spanning Tree (MST):

Prim's Algorithm

Cut property — Let  $S$  be any subset of nodes, and let  $e$  be the min cost edge with exactly one endpoint in  $S$ . Then the MST contains  $e$

So in Prim's Algorithm, everytime we add an edge, we follow cut property.

```
def prim(G = (V, E), c):
    u <- arbitrary vertex in V
    S <- { u } #let vertex u to be our starting vertex
    T <- NULL # That's our initialization of MST

    while len(S) < len(V) do
        # u in S and v not in S
        (u, v) <- min(S.cutset()) # min cost edge with exactly one endpoint in S
        add (u, v) to T # v never discover before, add it to T
        add v to S
    return T
```

# Implementation

```
# Sample Graph Strature
class Vertex:
    def __init__(self, value):
        self.value = value;
        self.adjacent_edges = [];

    def getValue(self):
        return self.value;

    def __repr__(self):
        return self.value;

    def __str__(self):
        return self.value;

    def addEdge(self, edge):
        self.adjacent_edges.append(edge);

    def getAdjacentEdges(self):
        return self.adjacent_edges;

class Edge:
    def __init__(self, name, weight):
        self.name = name;
        self.weight = weight;
        self.neighborVertices = []

    def __str__(self):
        return self.name + " " + str(self.weight);

    def __repr__(self):
        return self.name + " " + str(self.weight);

    # For compare x < y
    def __lt__(self, other):
        return self.weight < other.weight

    def addVertex(self, vertex: Vertex):
        if len(self.neighborVertices) == 2:
            print("Already connected with two vertices");
            return;
```

```

        self.neighborVertices.append(vertex);

def get_Neighbor_Vertices(self):
    return self.neighborVertices;

def get_weight(self):
    return self.weight;

class Adjacency_list_graph:
    def __init__(self):
        self.vertices = [];
        self.edges = []

    def addVertex(self, vertex: Vertex):
        self.vertices.append(vertex)

    def addVertices(self, list):
        self.vertices.extend(list);

    def addEdge(self, edge: Edge):
        self.edges.append(edge);
        for v in self.vertices:
            if v in edge.get_Neighbor_Vertices():
                v.addEdge(edge);

    def addEdges(self, edgeList):
        self.edges.extend(edgeList);
        for edge in edgeList:
            for v in self.vertices:
                if v in edge.get_Neighbor_Vertices():
                    v.addEdge(edge);

    def getVertices(self):
        return self.vertices;

    def getEdges(self):
        return self.edges;

```

```

from queue import PriorityQueue

```

```

def prim(G: Adjacency_list_graph):
    V = G.getVertices()
    discovered_Edge = PriorityQueue()
    spanningTree = []
    S = [V[0]]
    for edge in V[0].getAdjacentEdges():
        discovered_Edge.put((edge.get_weight(), edge))

    while len(S) < len(V):
        # discovered_Edge.sort(key = sortWeight);
        print([edge for edge in discovered_Edge.queue]);

        weight, local_min_edge = discovered_Edge.get()
        end_point1 = local_min_edge.get_Neighbor_Vertices()[0]
        end_point2 = local_min_edge.get_Neighbor_Vertices()[1]
        print("Add light edge: {} -- {}".format(end_point1, end_point2));
        print("");
        spanningTree.append((end_point1, end_point2))

        if end_point1 not in S:
            nexVertex = end_point1;
        else:
            nexVertex = end_point2;

        S.append(nexVertex)
        for e in nexVertex.getAdjacentEdges():
            if (e.get_weight(), e) not in discovered_Edge.queue and e !=
local_min_edge:
                discovered_Edge.put((e.get_weight(), e))
    return spanningTree

```

## Time Complexity — Similar analysis to Dijkstra's Algorithm

Using heap implementation:  $O(m \log n)$

Using Fibonacci heap:  $O(m + n \log n)$

## Kruskal's Algorithm

Consider edges in ascending order of weight.

**Case1:** If adding  $e$  to  $T$  creates a cycle, discard  $e$  according to cycle property.

**Case2:** Otherwise, insert  $e = (u, v)$  into  $T$  according to cut property where  $S$  = set of nodes in  $u$ 's connected component.

### Kruskal's Algorithm: Time complexity

Sorting edges takes  $O(m \log m)$  time

We need to be able to test if adding a new edge creates a cycle, in which case we skip the edge

- Option1: Run DFS in each iteration to see if the number of connect components stays the same. This leads to  $O(m * n)$  time for the main loop

### Union Find ADT

Union find — data structure defined on a ground set of elements  $A$

Used to keep track of an evolving partition of  $A$

Supported operations:

- $make\_sets(A)$  — make  $|A|$  singleton sets with elements in  $A$
- $find(a)$  — returns an id for the set element  $a$  belongs to
- $union(a, b)$  — union the sets elements  $a$  and  $b$  belong to

### Implementation

```
def Kruskal(V, E, c):
    sort E in increasing c-value
    answer = []
    comp = make_sets(V)
    for (u, v) in E do
        # check if u and v are already connected in comp
        if comp.find(u) != comp.find(v) then
            answer.append((u, v))
            comp.union(u, v)
    return answer
```

Sets are represented with a lists. An array points to the set each element belongs to:



- *make\_sets(A)* Create and initialize the array, take  $O(n)$
- *find(a)* is a simple lookup in the array, take  $O(1)$
- *union(a, b)* Add elements in u's set to v's set, take  $O(n)$

Kruskal's Algorithm would run in  $O(n^2)$  time

However, in **Better union-find implementation**:

Sets are represented with a lists. An array points to the set each element belongs to. Keep track of cardinality of each set. When taking the union of two sets change the smallest.

Kruskal's algorithm would be **optimized in  $O(m \log n)$  time**