# Faculty of Information Technology, Monash University

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# FIT2004: Algorithms and Data Structures

Week 5: Greedy (Graph) Algorithms

# **Overview**

Divide and conquer (W 1-3)

Greedy algorithms (W 4-5) Dynamic programming (W 6-7)

Network flow (W 8-9) structures (W 10-11)

- Last Lecture
  - Introduction to Graphs
  - Graph Traversal Algorithms
- Today's Lecture
  - Dijkstra's Algorithm
  - Prim's Algorithm
  - Kruskal's Algorithm

FIT2004: Seminar 5 - Greedy (Graph) Algorithms

# **Greedy Algorithm**

- Greedy paradigm for algorithms: make a locally optimal choice at each stage and commit to it.
- Often it is easy to come up with ideas to solve a problem using a greedy strategy, but in many cases the proposed greedy algorithm does not find the overall optimal solution for the problem.
- Correctness proofs are very important.
- o Today we will take a look at three very important greedy algorithms:
  - Dijkstra's algorithm solves the shortest path problem in graphs with non-negative weights.
  - Prim's and Kruskal's algorithms present two different ways to find a minimum spanning tree of a graph (an important problem for which there are actually multiple distinct greedy strategies that work fine).

## **Outline**

- 1. Dijkstra's Algorithm
- 2. Prim's Algorithm
- 3. Kruskal's Algorithm

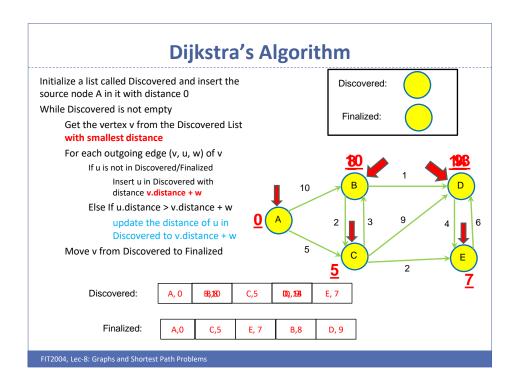


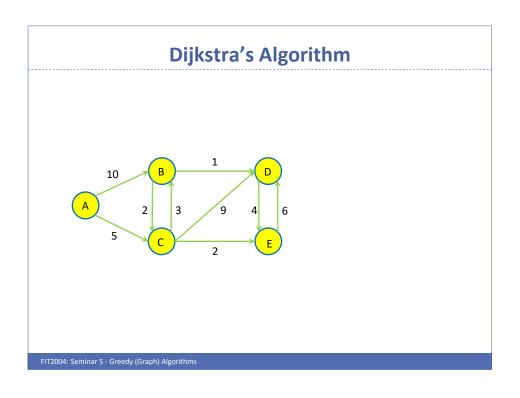
Edsger W. Dijkstra Turing Award 1972

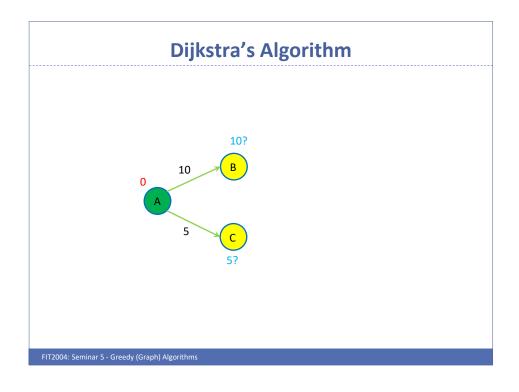
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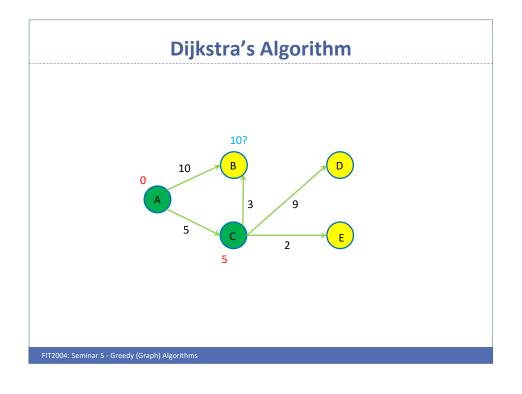
# Dijkstra's Algorithm

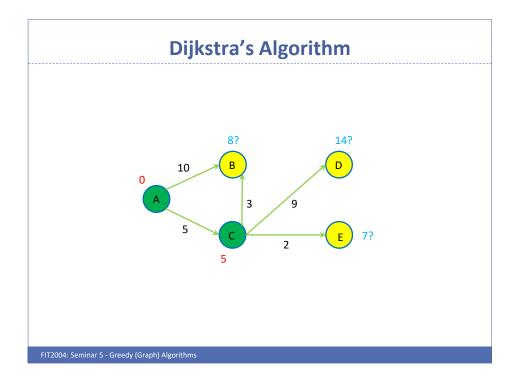
- Solves the single source, all targets shortest path problem on graphs with non-negative weights. Closely related to BFS.
- Keeps track of a set S of nodes whose distance to the source has already been determined.
- Initially S only contains the source node, and it grows by one element at a time until finding all nodes that are reachable from the source node.
- At each iteration, from all nodes that are one edge away from
   S, add the node with the smallest distance to the source to S.
- By doing so, all the nodes will be added to S in the increasing order of distance to the source.

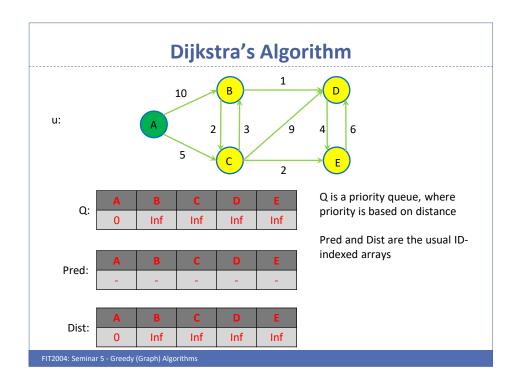


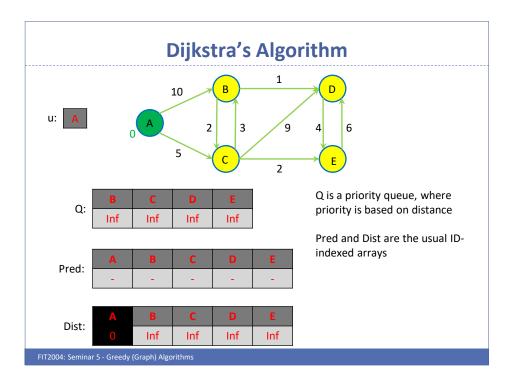


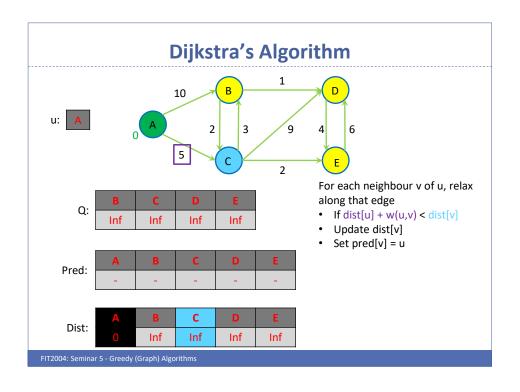


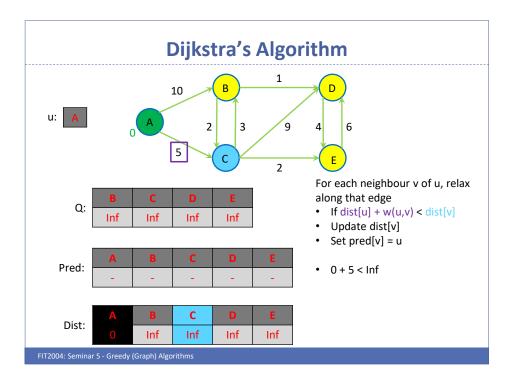


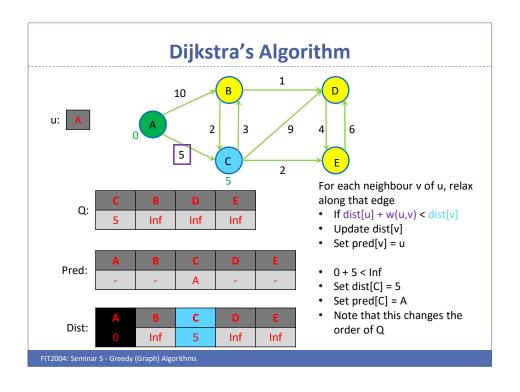


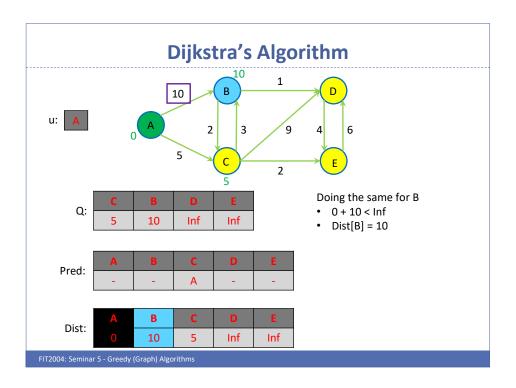


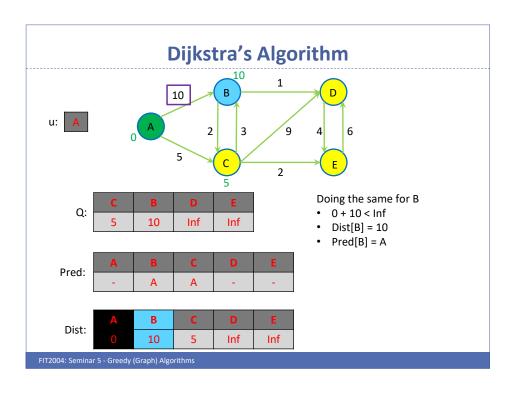


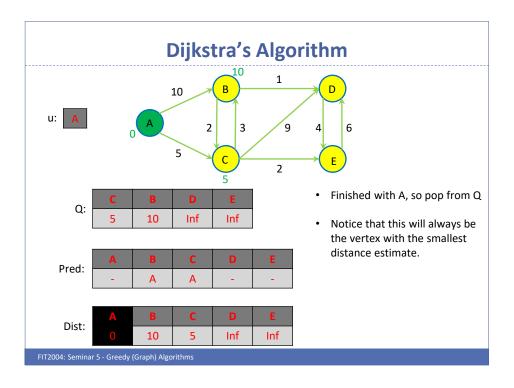


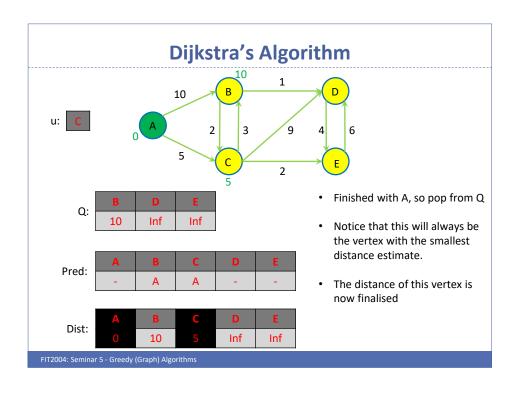


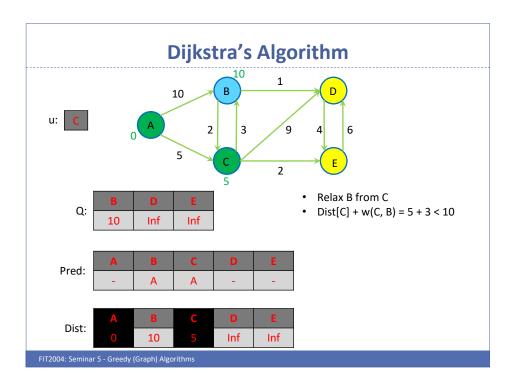


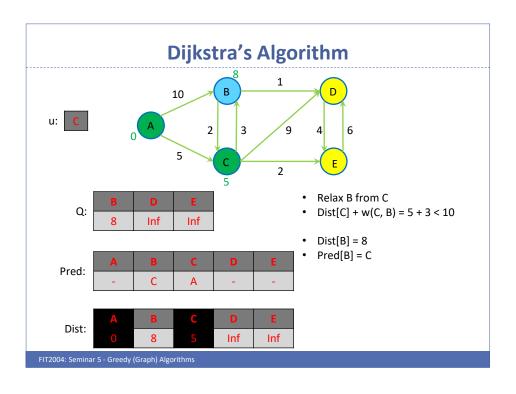


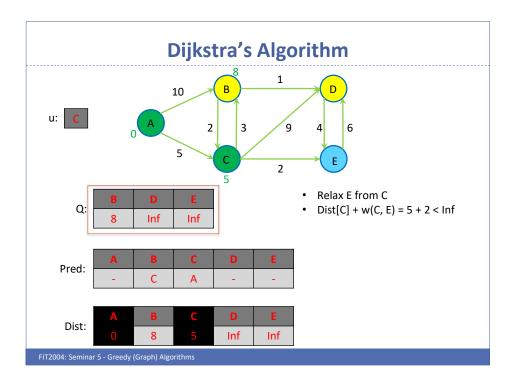


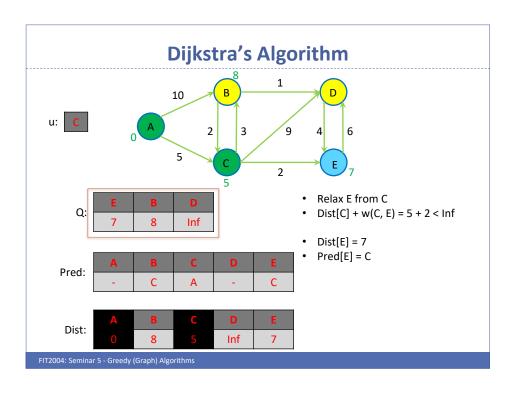


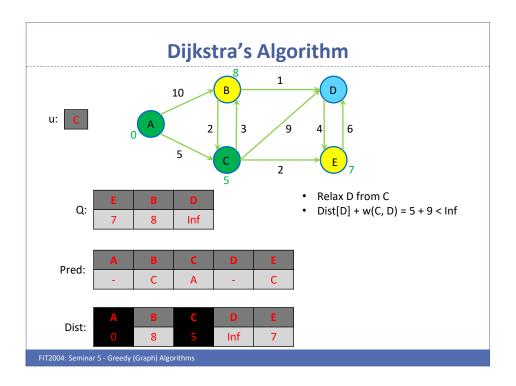


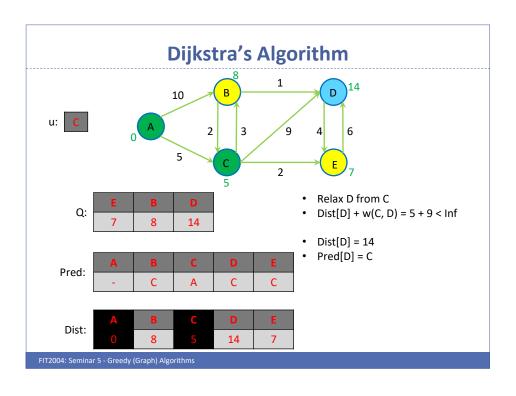


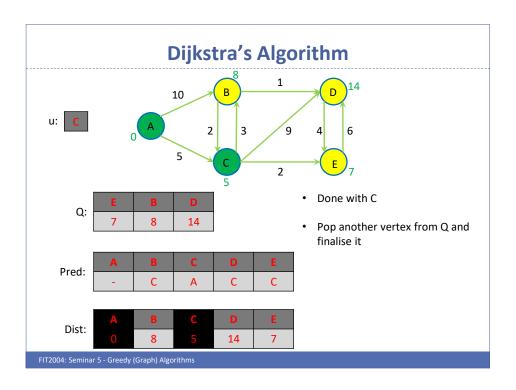


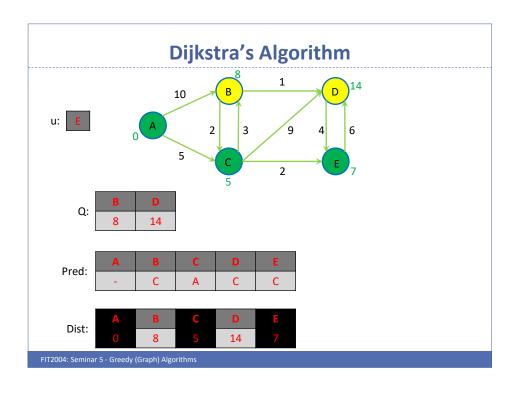


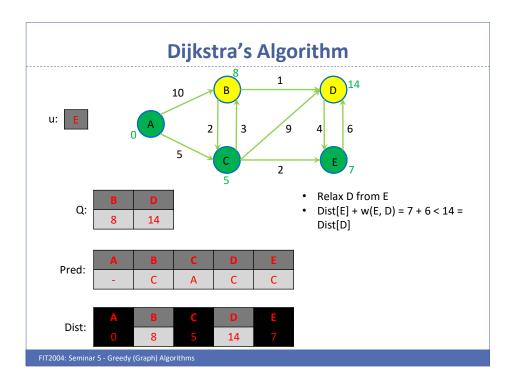


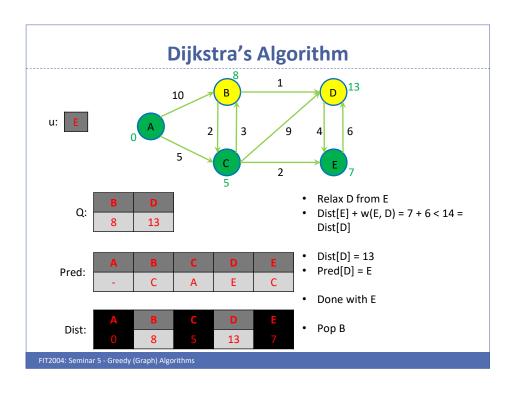


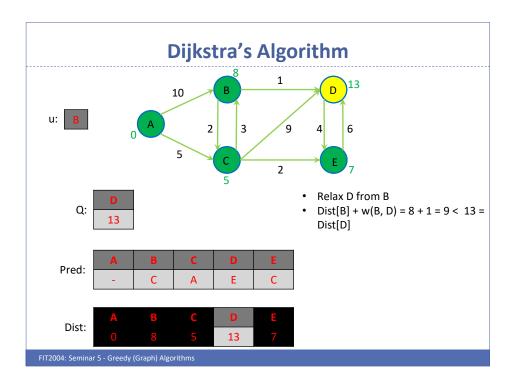


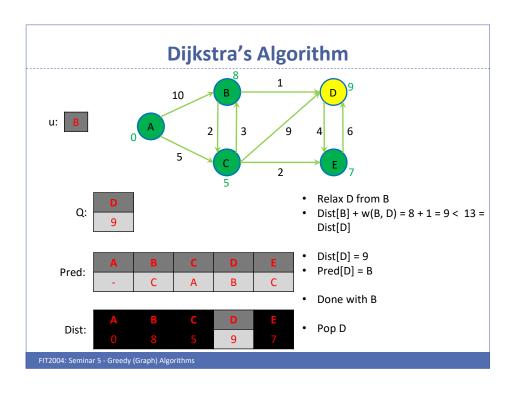


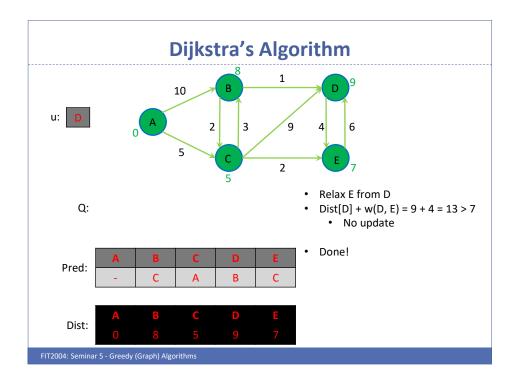


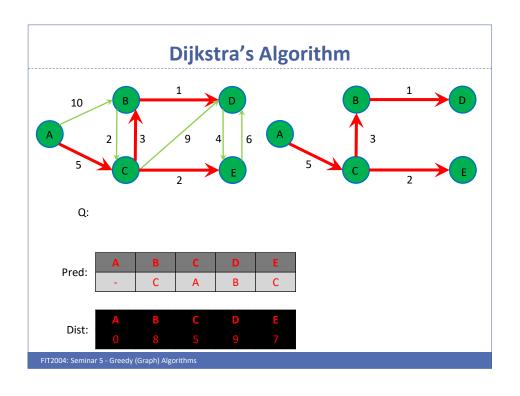












# Dijkstra's Algorithm

```
Algorithm 38 Dijkstra's algorithm
 1: function DIJKSTRA(G = (V, E), s)
        dist[1..n] = \infty
        pred[1..n] = 0
 3:
        dist[s] = 0
        Q = \overrightarrow{\text{priority\_queue}}(V[1..n], \text{key}(v) = dist[v])
        while Q is not empty do
           u = Q.pop_min()
           for each edge e that is adjacent to u do
               // Priority queue keys must be updated if relax improves a distance estimate!
               RELAX(e)
       return dist[1..n], pred[1..n]
Algorithm 37 Edge relaxation
 1: function RELAX(e = (u, v))
       if dist[v] > dist[u] + w(u, v) then
```

EIT2004: Saminar E Grandy (Granh) Algarithm

dist[v] = dist[u] + w(u, v)pred[v] = u

4:

# Dijkstra's Algorithm

```
Algorithm 38 Dijkstra's algorithm
 1: function DIJKSTRA(G = (V, E), s)
       dist[1..n] = \infty
                                                                                                         While loop: \Theta(V)
        pred[1..n] = 0
        dist[s] = 0
        Q = \text{priority\_queue}(V[1..n], \text{key}(v) = \text{dist}[v])
        while Q is not empty do
           u = Q.pop min()
                                                                                                            For loop: \Theta(E)
           for each edge e that is adjacent to u do
               // Priority queue keys must be updated if relax improves a distance estimate!
 9:
10:
               RELAX(e)
       \textbf{return} \ \textit{dist}[1..n], \textit{pred}[1..n]
```

### Time Complexity:

- $\circ$  While loop executes  $\theta(V)$  times
  - Find the vertex with smallest distance and <u>delete</u> it: depends on priority queue implementation (Q.extract\_min)
- Each edge visited once  $\rightarrow \theta(E)$ 
  - Updating the priority queue: depends on implementation (Q.update)
  - Relaxation is  $\theta(1)$  since we can find distances and compare them in  $\theta(1)$
- Total cost: θ(E\*Q.update+ V\*Q.extract\_min)

# Dijkstra's Algorithm using min-heap

Total cost: θ(E\*Q.update+ V\*Q.extract\_min)

### Required additional data structure:

- Create an array called Vertices.
- Vertices[i] will record the location of i-th vertex in the min-heap

### Updating the distance of a vertex v in min-heap in $\theta(\log V)$

- o Find the location in the queue (heap) in  $\theta(1)$  using Vertices
- Now do the normal heap-up operation in  $\theta(\log V)$ 
  - For each swap performed between two vertices x and y during the upHeap
    - Update Vertices[x] and Vertices[y] to record their updated locations in the min-heap

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# Time Complexity of Dijkstra's Algorithm

- $\circ$  While loop executes  $\theta(V)$  times
  - Extract the vertex with smallest distance using min heap:  $\theta(\log V)$
- Each edge visited once  $\rightarrow \theta(E)$ 
  - Updating the priority queue using min heap:  $\theta(\log V)$
  - Relaxation is  $\theta(1)$  since we can find distances and compare them in  $\theta(1)$
- Total cost: θ(E\*Q.update+ V\*Q.extract\_min)
- o Total cost:  $\theta(E \log V + V \log V)$ 
  - o Which term dominates?

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- o If the graph is connected, then E ≥ V-1
  - E dominates V
- Total cost:  $\theta(E \log V)$

Algorithm 38 Dijkstra's algorithm

- 1: **function** DIJKSTRA(G = (V, E), s) 2:  $dist[1..n] = \infty$
- pred[1..n] = 0
   dist[s] = 0
- 5:  $Q = \text{priority\_queue}(V[1..n], \text{key}(v) = \text{dist}[v])$ 6: **while** Q **is not empty do**
- 6: while Q is not empty do 7:  $u = Q.pop\_min()$
- 8: **for each** edge *e* that is adjacent to *u* **do**9: // Priority queue keys must be updated
  10: RELAX(*e*)
- 11: **return** dist[1..n], pred[1..n]
- Otal Cost. O(L log V)

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# **Proof of Correctness**

Claim: For every vertex v which has been removed from the queue, dist[v] is correct.

- Notation:
  - V is the set of vertices
  - o Q is the set of vertices in the queue
  - S = V / Q = the set of vertices that have been removed from the queue

### **Base Case**

 dist[s] is initialised to 0, which is the shortest distance from s to s (since there are no negative weights).

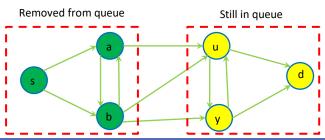
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# **Proof of Correctness**

Claim: For every vertex v which has been removed from the queue, dist[v] is correct.

### Inductive Step:

- Assume that the claim holds for all vertices which have been removed from the queue, in other words it holds for all vertices in the set S.
- o Let u be the next vertex which is removed from the queue.
- We will show that dist[u] is correct.



# **Proof of Correctness**

Claim: For every vertex v which has been removed from the queue, dist[v] is correct.

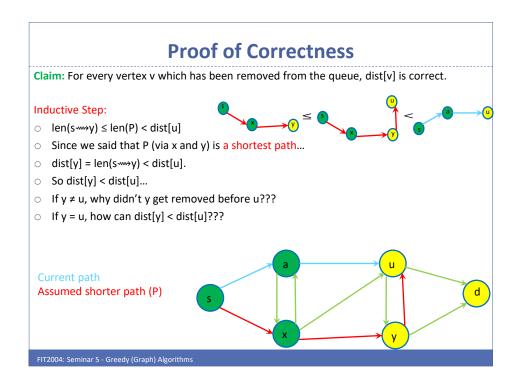
### **Inductive Step:**

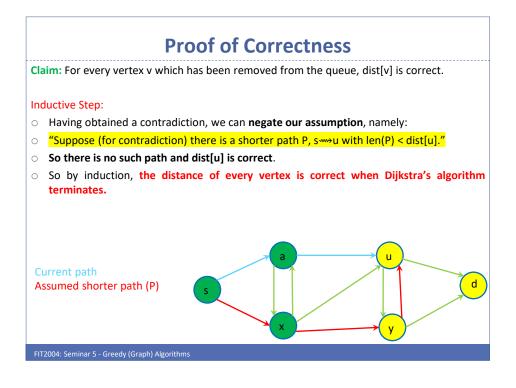
- Suppose (for contradiction) there is a shortest path P, s→u with len(P) < dist[u].
- O Let x be the furthest vertex on P which is in S (i.e. has been finalised).
- By the inductive hypothesis, dist[x] is correct (since it is in S).

Current path
Assumed shorter path (P)

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# Proof of Correctness Claim: For every vertex v which has been removed from the queue, dist[v] is correct. Inductive Step: By the inductive hypothesis, dist[x] is correct (since it is in S). Let y be the next vertex on P after x. By assumption len(P) < dist[u]. Edge weights are non-negative. len(s→→y) ≤ len(P) < dist[u]. Current path Assumed shorter path (P) Assumed shorter path (P)





# **Outline**

- 1. Dijkstra's Algorithm
- 2. Prim's Algorithm
- 3. Kruskal's Algorithm

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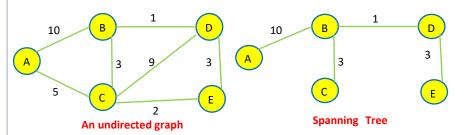
# What is a Spanning Tree

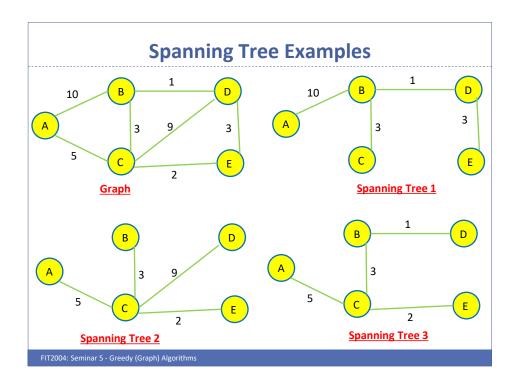
### Tree

A tree is a connected undirected graph with no cycles in it.

### Spanning Tree:

A spanning tree of a general undirected weighted graph G is a tree that spans G
 (i.e., a tree that includes every vertex of G) and is a subgraph of G (i.e., every
 edge in the spanning tree belongs to G).





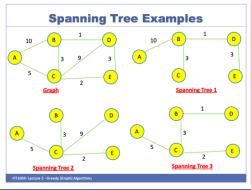


Is it true that a spanning tree of a connected graph G is a maximal set of edges of G that contains no cycles?

Yes

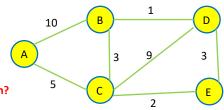
Is it true that a spanning tree of a connected graph G is a minimal set of edges that connect all vertices?

Yes

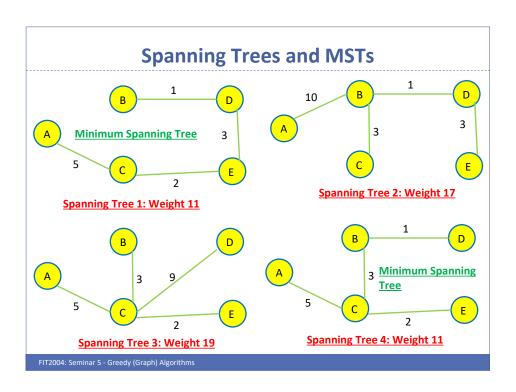


# **Minimum Spanning Tree (MST)**

- Weight of a spanning tree is the sum of the weights of the edges in the tree.
- A minimum spanning tree of a weighted general graph G is a tree that spans G, whose weight is minimum over all possible spanning trees for this graph.
- There may be more than one minimum spanning tree for a graph G (e.g., two or more spanning trees with the same minimum weight).



What is the weight of the MST in this graph? How many MSTs are in this graph?



# **MST Algorithms**

- Let T denote the MST we are constructing, initialised to be empty.
- An edge e is said to be safe if {T U e} is a subset of some MST

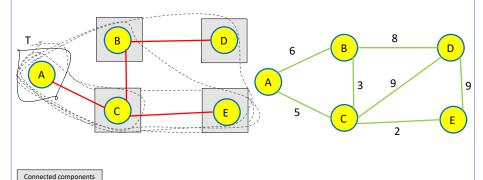
### General Strategy:

- ➤ T = null
- while T can be grown safely:
- find an edge e=<x,y> along which T is safe to grow
- $T = \{T\} \text{ union } \{\langle x,y\rangle\}$
- return T
- We will study two **greedy** algorithms that follow this strategy.

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# Prim's Algorithm: Overview

- Start by picking any vertex v to be the source of T.
- While T does not contain <u>all</u> vertices in the graph:
  - Find shortest edge e that goes from T to a different connected component.
  - Add e to the tree T.



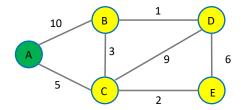
# MST Algorithms: Prim's Algorithm

Prim's Algorithm (small modification of Dijkstra's Algorithm)

- We start with no edges in T, so there are V connected components in T (each with a single node) and select one node as the source.
- At each iteration, we add to T the shortest edge that links the connected component containing the source to a different connected component.
- In each iteration, the number of connected components decreases by 1 and the number of nodes in the connected component containing the source increases by 1.
- Cycles are avoided as no edges linking two nodes that are already in the same connected component are ever added to T.
- Time complexity:  $\theta(E \log V)$

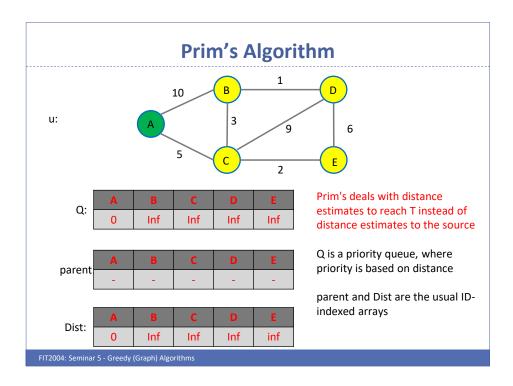
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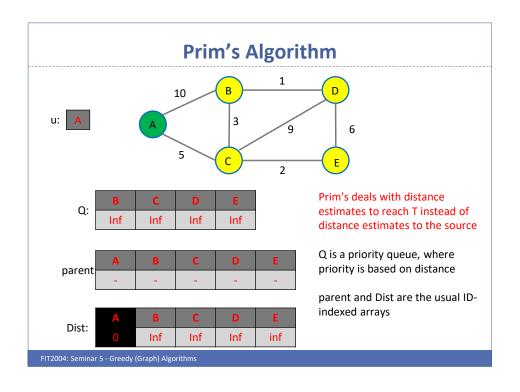
# **Prim's Algorithm**

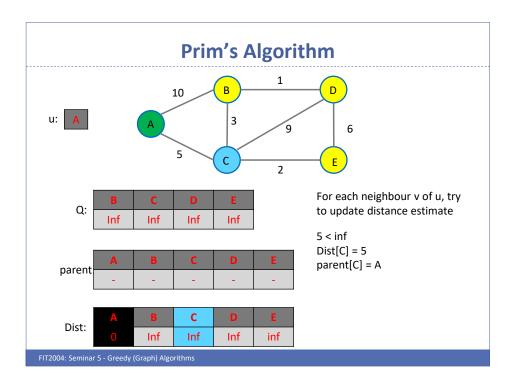


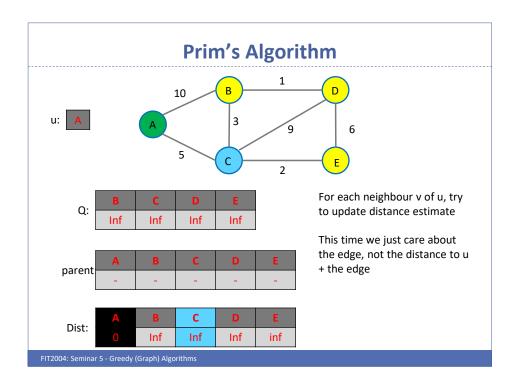


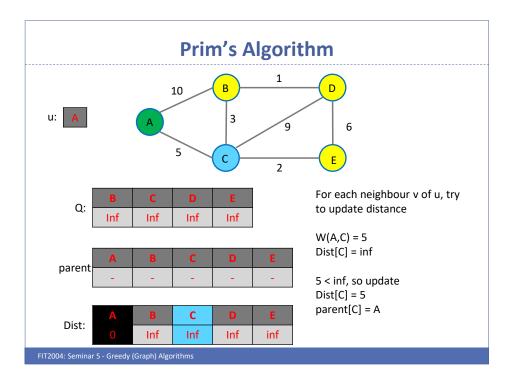
Robert C. Prim

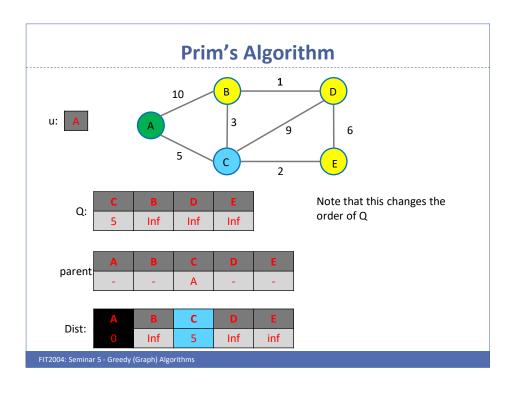


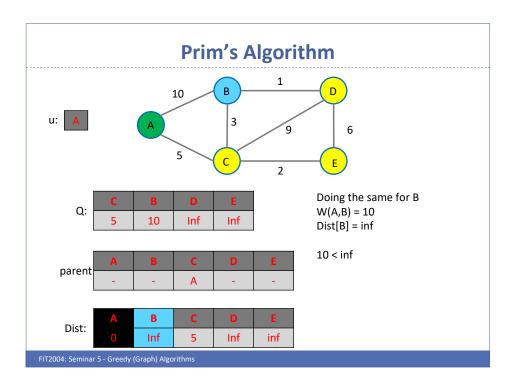


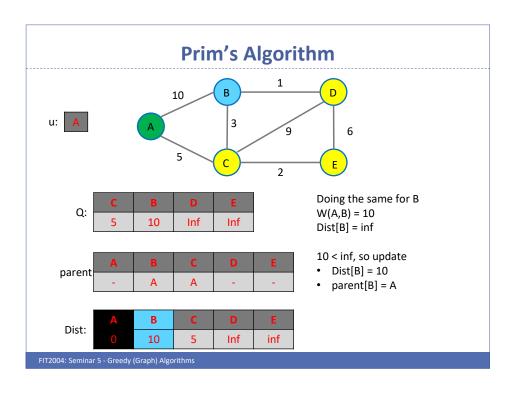


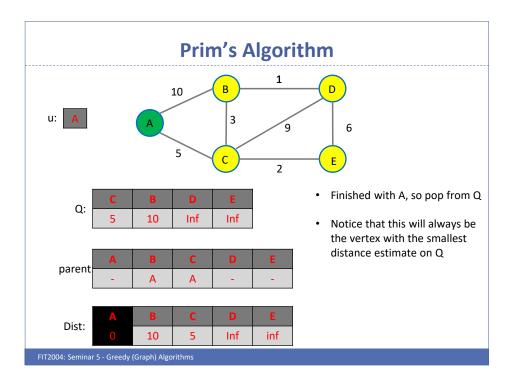


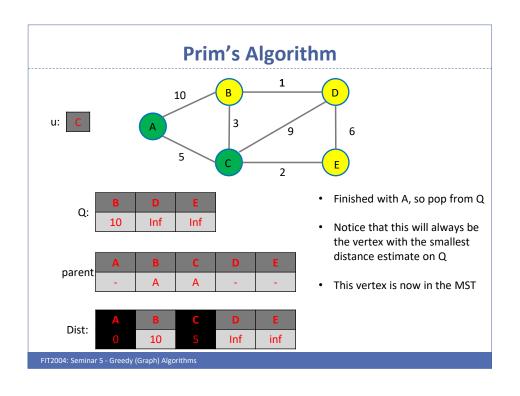


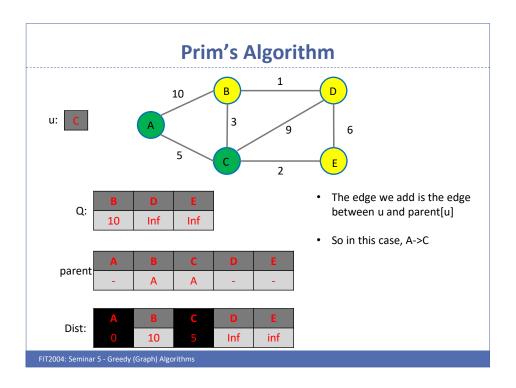


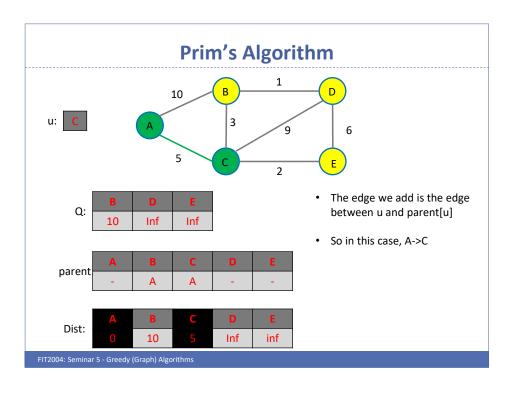


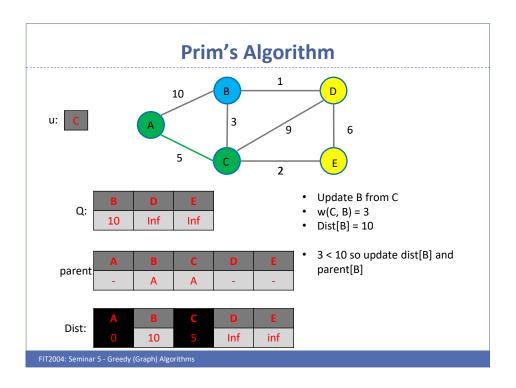


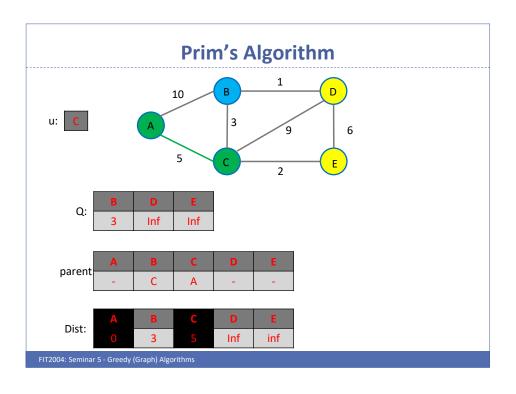


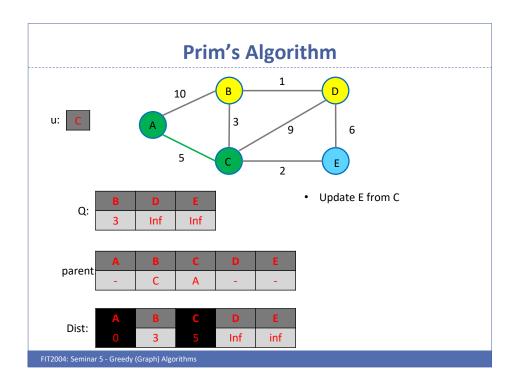


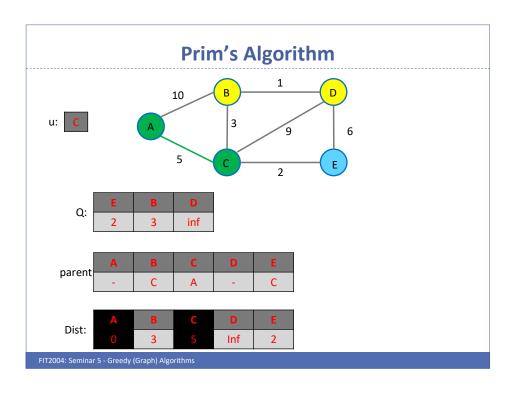


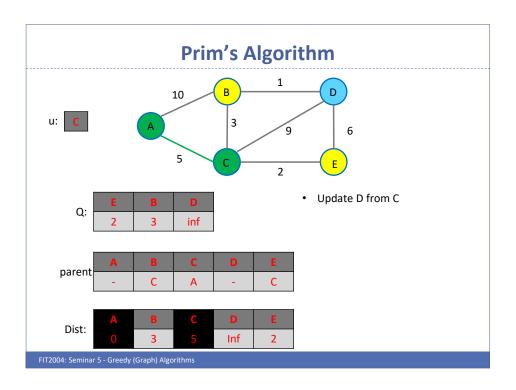


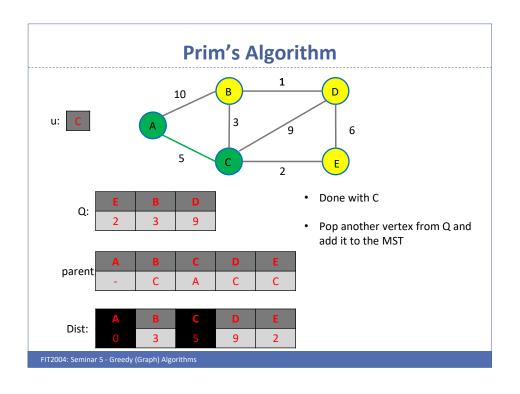


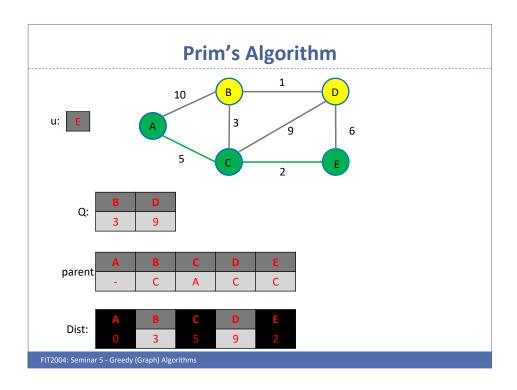


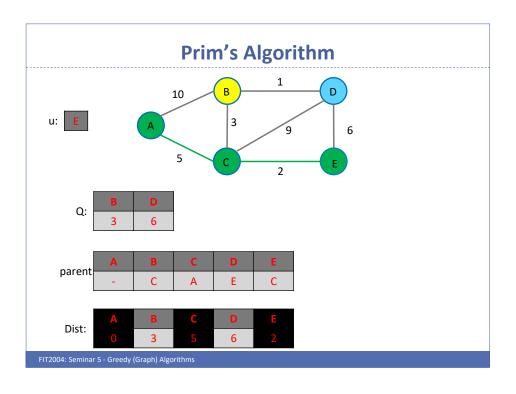


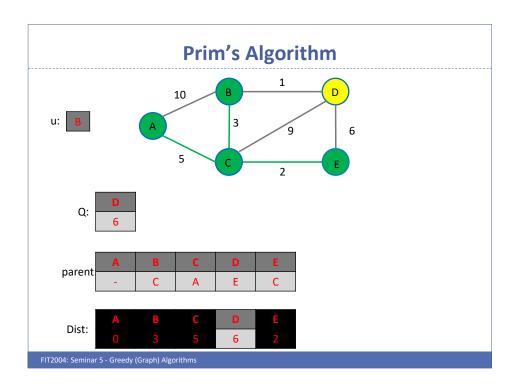


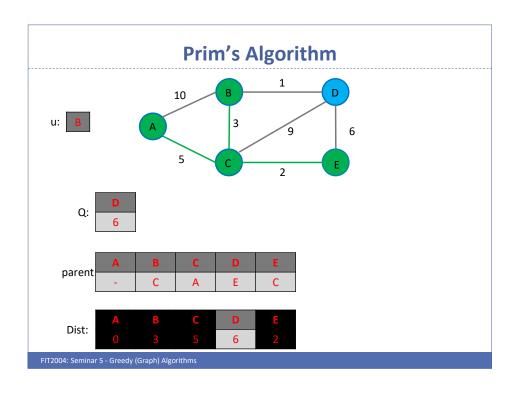


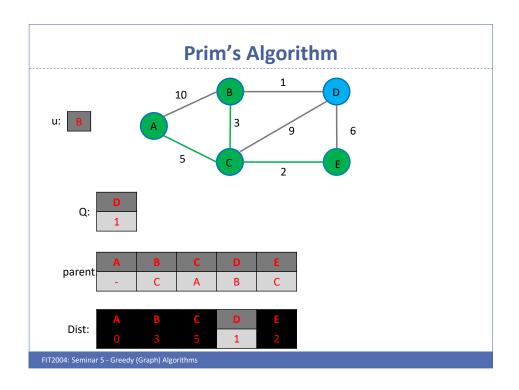


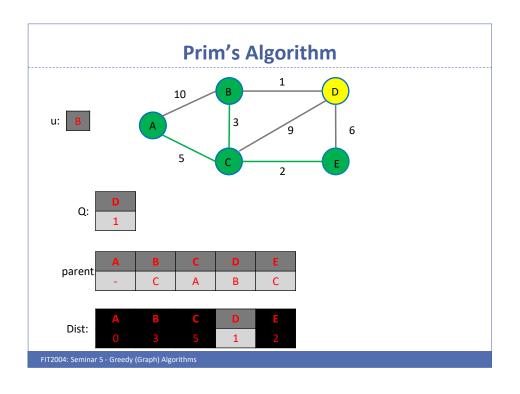


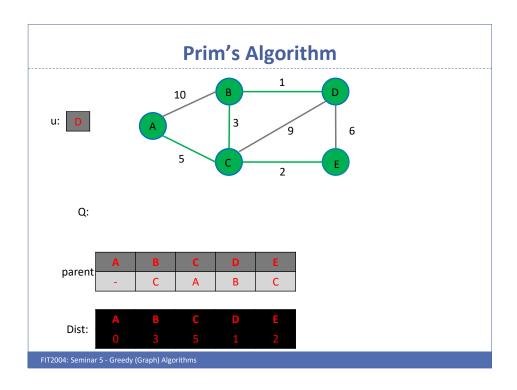


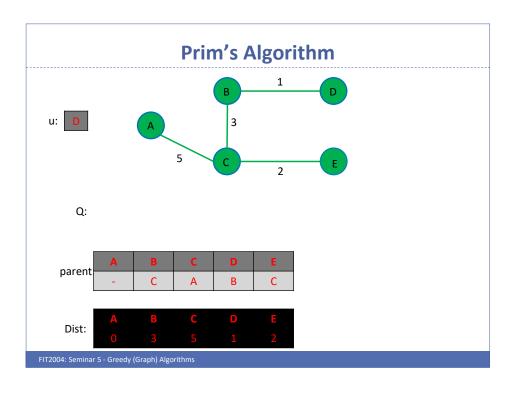












## **Prim's Algorithm**

#### Algorithm 69 Prim's algorithm

```
1: function PRIM(G = (V, E), r)
       dist[1..n] = \infty
       parent[1..n] = null
       T = (\{r\}, \emptyset)
 4:
       dist[r] = 0
 5:
       Q = \text{priority\_queue}(V[1..n], \text{key}(v) = dist[v])
 6:
       while Q is not empty do
 7:
           u = Q.pop_min()
 8:
           T.add\_vertex(u)
 9:
           T.add\_edge(parent[u], u)
10:
           for each edge e = (u, v) adjacent to u do
11:
               if not v \in T and dist[v] > w(u, v) then
12:
                  // Remember to update the key of v in the priority queue!
                  dist[v] = w(u, v)
14:
                  parent[v] = u
15:
       return T
16:
```

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## **Prim's Algorithm: Complexity**

It is very similar to Dijkstra's Algorithm and its complexity is the same as Dijkstra's Algorithm

o  $\theta(V \log V + E \log V)$  if min-heap is used

Since the input graph G is connected, E  $\geq$  V-1. Hence, the complexity can be simplified to  $\theta(E \log V)$ .

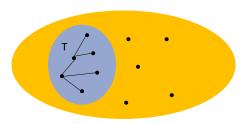
#INV: Every iteration of Prim's algorithm, the current set of selected edges in T is a subset of some minimum spanning tree of G

#### **Base Case:**

o The invariant is true initially when T is empty

#### Inductive step:

- We want to show that, if T is a subset of some MST at the start of some iteration, it is still a subset of some MST at the start of the next iteration
- Assume T is a subset of some MST M



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## **Prim's Algorithm: Correctness**

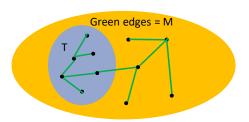
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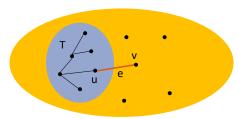
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- o If e is in M, then T U (e) is a subset of M, which is an MST, so the invariant holds



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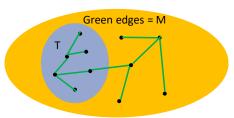
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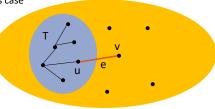
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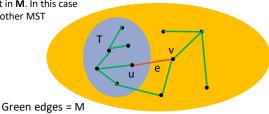
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#### Inductive step:

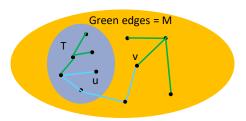
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#### Inductive step:

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- o Since M is a tree, there is exactly one path from **u** to **v** in **M** (shown in blue)
- u and v are not connected in T (since v is not in T). Consider the first edge on the blue path which is not contained in T (call this edge x).

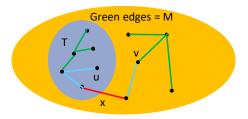


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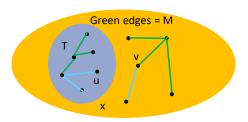
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- $\circ\quad$  One vertex of this edge is in  ${\bf T},$  the other is not.
- o Removing this edge would disconnect M



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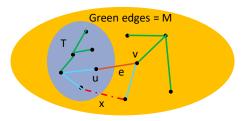


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## **Prim's Algorithm: Correctness**

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- u and v are not connected in T (since v is not in T). Consider the first edge on the blue path which is not contained in T (call this edge x).
- Removing this edge would disconnect M
- Adding the edge e=(u,v) would form a new spanning tree, M'
- Since Prim's algorithm always selects the lightest edge incident to T, we know that w(e) ≤ w(x)
- So the weight of M' is no greater than the weight of M, therefore choosing e is correct



### **Outline**

- 1. Dijkstra's Algorithm
- 2. Prim's Algorithm
- 3. Kruskal's Algorithm

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## MST Algorithms: Kruskal's Algorithm

#### Kruskal's Algorithm

- We start with no edges in T, so there are V connected components in T (each with a single node).
- We process edges in ascending order of edge weights, and add an edge e to T if this addition does not create a cycle.
- Each time that one edge e is added to T, e is the smallest edge that links two distinct connected components of T.
- Cycles are avoided as no edges linking two nodes that are already in the same connected component are ever added to T. At any point of the execution T is a forest (i.e., a set of trees)
- Time complexity:  $\theta$ (E log V).

## Kruskal's Algorithm

#### Kruskals(G(V, E))

- Sort the edges in ascending order of weights
- Let T be a graph with V as its vertices, and no edges
- For each edge (u, v) in ascending order
  - If adding (u, v) does not create a cycle in T
    - > Add (u, v) to T
- Return T

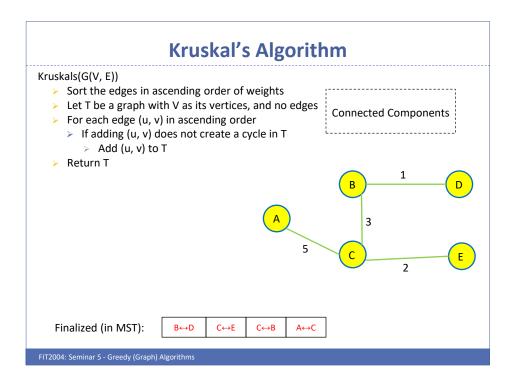


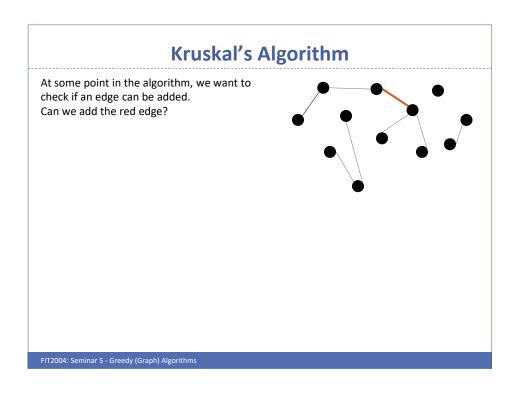
Joseph Kruskal

1 В D 10 9 4

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#### Kruskal's Algorithm Kruskals(G(V, E)) Sort the edges in ascending order of weights Let T be a graph with V as its vertices, and no edges **Connected Components** For each edge (u, v) in ascending order If adding (u, v) does not create a cycle in T > Add (u, v) to T Return T 1 10 How to determine if the edge will create a cycle??? 9 B↔D, 1 C↔E, 2 C↔B, 3 E↔D, 4 A↔C, 5 A↔B, 10 Sorted Edges: Finalized (in MST): C↔E C↔B A↔C FIT2004: Seminar 5 - Greedy (Graph) Algorithms





## Kruskal's Algorithm

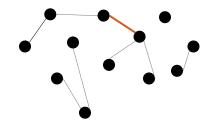
At some point in the algorithm, we want to check if an edge can be added.
Can we add the red edge?

An edge can be added if it does not create a cycle.

OR

An edge can be added if it is not between two nodes which belong to the same connected component.

Importantly, adding an edge between two distinct connected components "merges" those connected components into one.

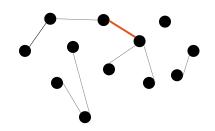


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## Kruskal's Algorithm

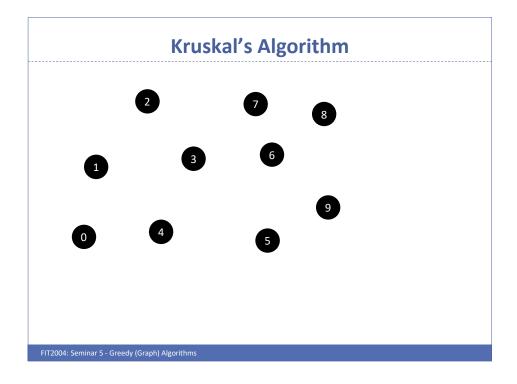
An edge can be added if it is NOT between two nodes which belong to the same connected component.

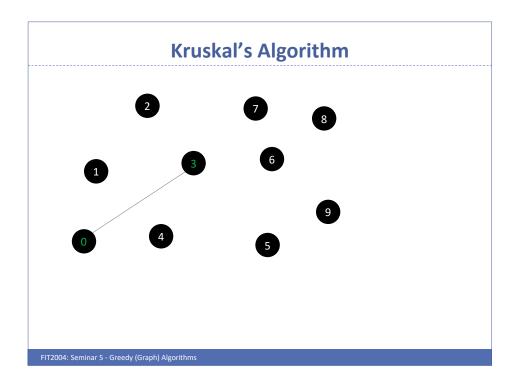
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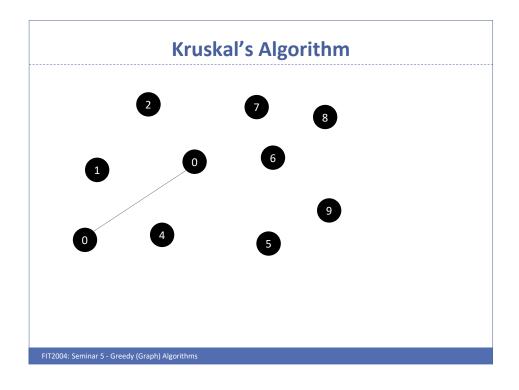


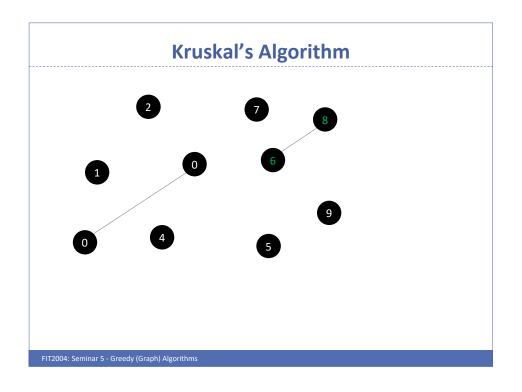
Remember that in Kruskal's algorithm, each time that one edge e is added to T, e is the smallest edge that links two distinct connected components of T.

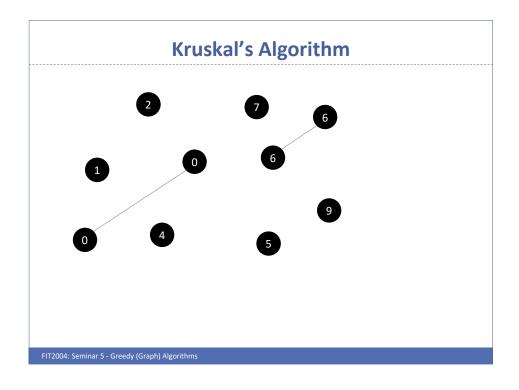
Intuitively, Kruskal's algorithm runs for V-1 iterations, and in each iteration it merges the two distinct connected components that have the smallest distance between themselves.

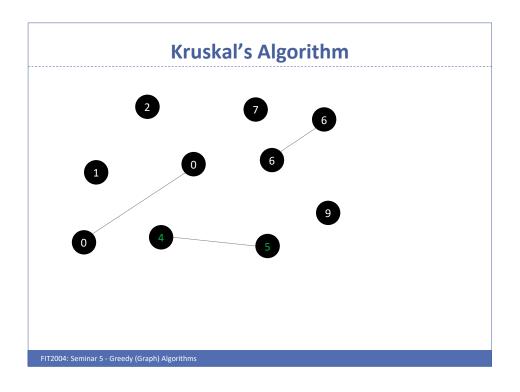


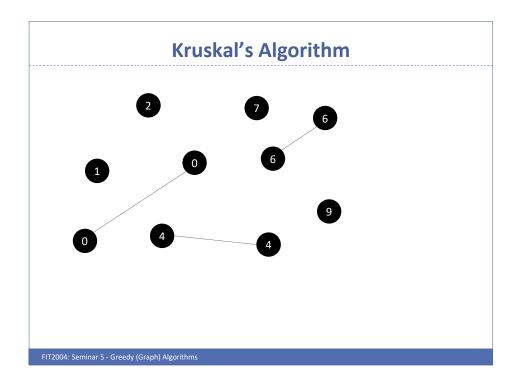


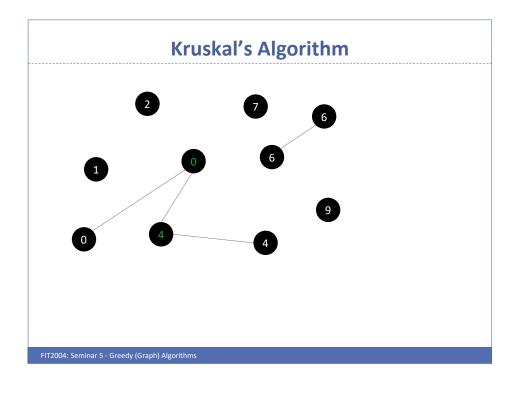


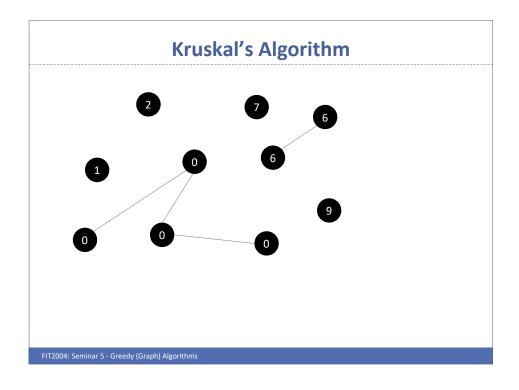


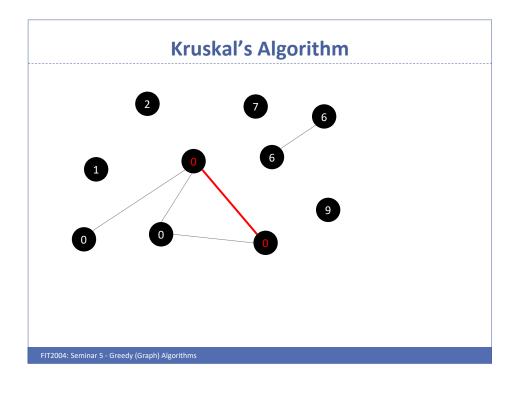












## Kruskal's Algorithm

#### Algorithm 41 Kruskal's algorithm

```
1: function KRUSKAL(G = (V, E))
      sort(E, key((u, v)) = w(u, v))
                                                      // Sort edges in ascending order of weight
3:
      forest = UnionFind.initialise(n)
      T = (V, \emptyset)
4:
5:
      for each edge (u, v) in E do
         if forest.FIND(u) \neq forest.FIND(v) then
                                                        // Ignore edges that would create a cycle
6:
7:
             forest.UNION(u, v)
             T.add_edge(u, v)
8:
9:
      return T
```

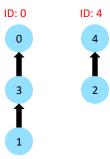
How can we quickly find the set a vertex belongs to, and also union two sets?

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### **Union-Find Data Structure**

- Define two operations: find(u) and union(u,v)
- Find(u): Given a vertex u, return its set ID
- Union(u,v): Given two vertices u and v, if they have different set IDs, union the
  two sets they belong to (and update all the set IDs of the vertices in one of the
  sets)
- We need both find(u) and union(u,v) to be fast.
  - What would be the time complexity of find(u) and union(u,v) if we store the set ID of each vertex in an array?
    - $\circ$  **Find** would be θ(1), but **union** would be θ(V) in the worst case, thus not good enough...

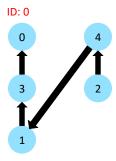
- We want to union faster
- o Linked lists are fast to union (i.e. append one linked list to another)



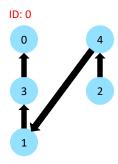
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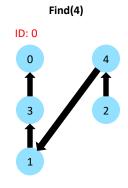
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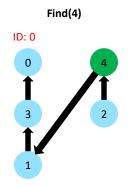
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### **Union-Find Data Structure**

- Linked lists are fast to union (i.e. append one linked list to another)
- O How do we find, with linked lists?
- Heads know their set ID
- Traverse to the head to find the ID of an element in the list
- O This is  $\theta$ (size of the linked list)



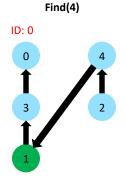
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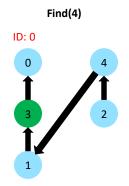
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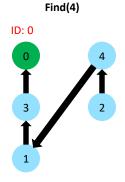
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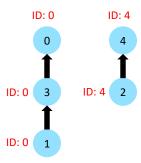
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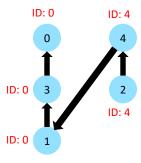
- Alternatively, every node could know its ID
- Now find is  $\theta(1)$
- o But Union is now slower, we have to update all the IDs



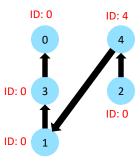
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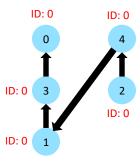
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## **Union-Find Data Structure**

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- O Where are we?
- We want find and union operations to be fast
- Linked lists are an improvement over arrays, since they stop us looking at items which are not relevant to the union we are currently doing
- Linked lists allows  $\theta(1)$  union
- We can't make find in  $\theta(1)$  since we have to store the ID at every node which makes union slow (we have to change all the IDs)
- Solution: Change from linked list to linked tree

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### **Union-Find Data Structure**

**Operations:** 



1

1 2

3

1

**1** 5

Vertex ID	0	1	2	3	4	5
Parent	0	1	2	3	4	5

**Operations:** Union(0,1)

Find(0) = 0 Find(1) = 1







**1** 3

4

**1** 5

Vertex ID	0	1	2	3	4	5
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## **Union-Find Data Structure**

**Operations:** Union(0,1)



1

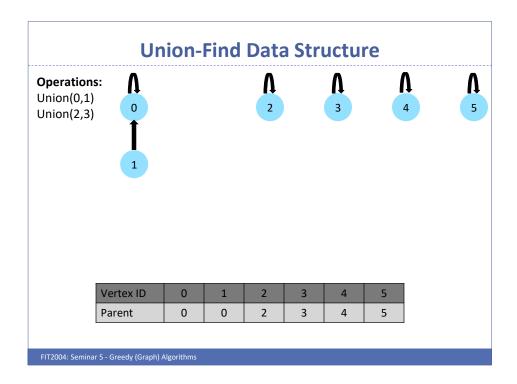
**1** 3

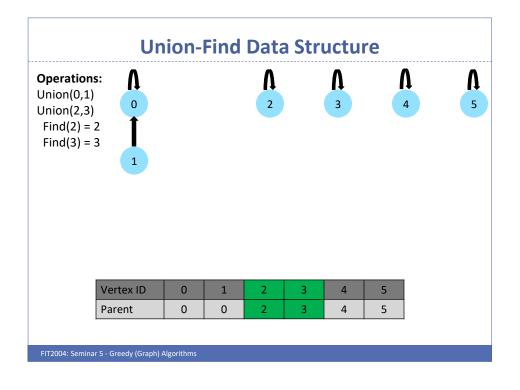
1

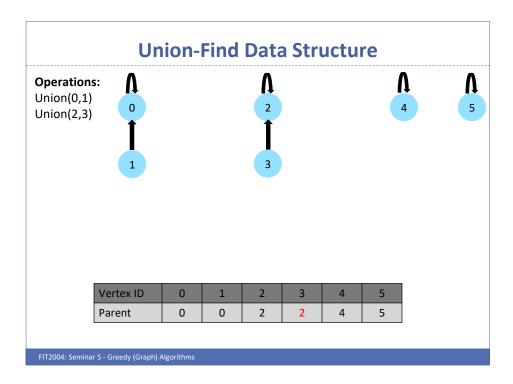
**1** 5

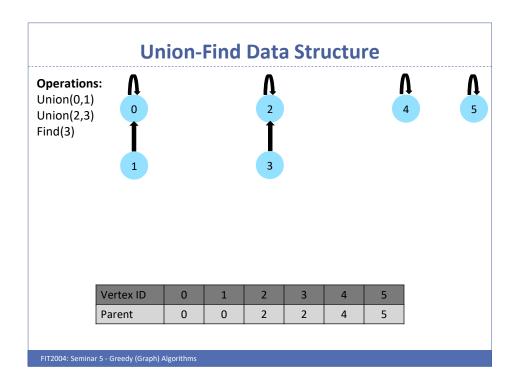
 Vertex ID
 0
 1
 2
 3
 4
 5

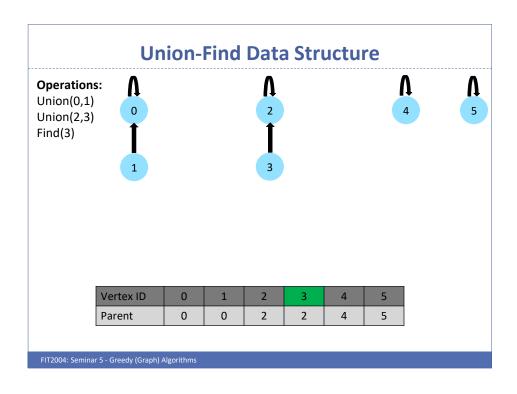
 Parent
 0
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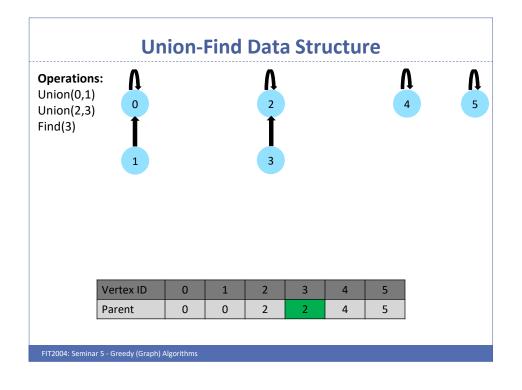


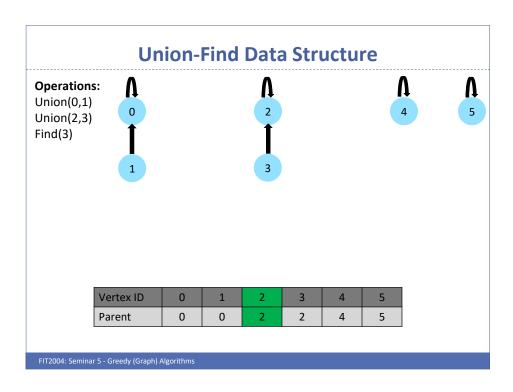


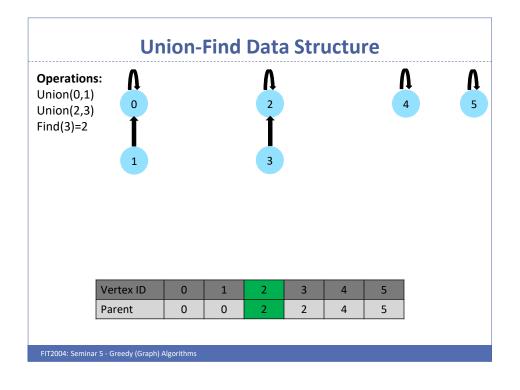


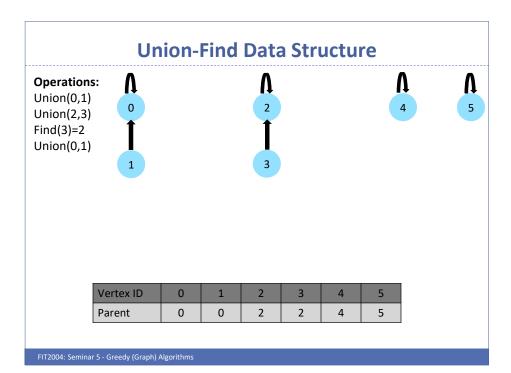


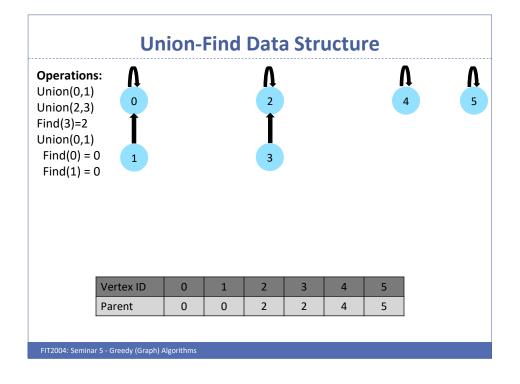


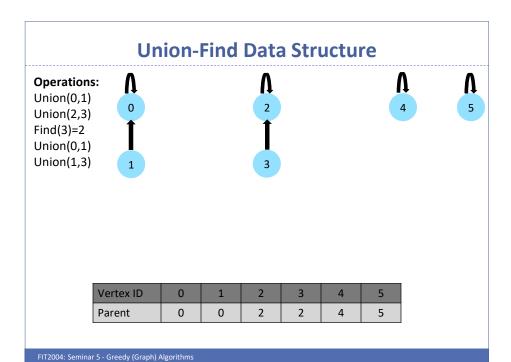


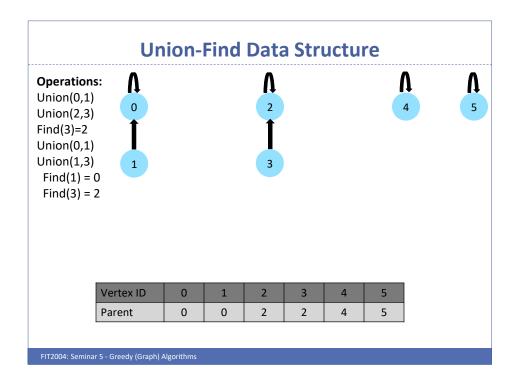




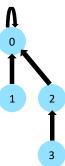








Operations: Union(0,1)Union(2,3)Find(3)=2Union(0,1) Union(1,3) Find(1) = 0Find(3) = 2



V	
4	



Vertex ID 4 5 Parent 0 2 4 5

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### **Union-Find Data Structure**

#### **Operations:**

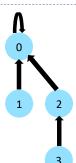
Union(0,1)

Union(2,3)Find(3)=2

Union(0,1)Union(1,3)

Find(1) = 0

Find(3) = 2



Λ

Find: traverse parent pointers until you find a vertex who is its own parent (i.e. a root). That vertex ID is the set ID

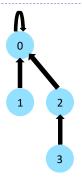
Union(u,v): If find(u) != find(v), set parent[find(u)] = find(v) (or vice versa)

What is the complexity of union and find? Union is  $\theta$ (find). Find could in theory be  $\theta$ (V), but if we can keep the heights of the trees low, then it will be at most  $\theta$ (max height)

Vertex ID	0	1	2	3	4	5
Parent	0	0	0	2	4	5



Operations: Union(0,1) Union(2,3) Find(3)=2 Union(0,1) Union(1,3) Find(1) = 0 Find(3) = 2



4

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**Optimisation:** When we union, we have a choice of the new root.

What should we choose?

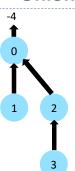
The set with more nodes!

Vertex ID	0	1	2	3	4	5
Parent	0	0	0	2	4	5

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### **Union-Find Data Structure**

Operations: Union(0,1) Union(2,3) Find(3)=2 Union(0,1) Union(1,3) Find(1) = 0 Find(3) = 2



-1 **1**  -1 **1** 5

**Optimisation:** When we union, we have a choice of the new root.

What should we choose?

The set with more nodes!

Union by size: When doing a union, add the sizes and update the parent value of the root appropriately

Vertex ID	0	1	2	3	4	5
Parent	-4	0	0	2	-1	-1

Parent array values are now either parents OR negative sizes

## Kruskal's Algorithm: Complexity

```
Algorithm 41 Kruskal's algorithm
  1: function KRUSKAL(G = (V, E))
         sort(E, key((u, v)) = w(u, v))
                                                             // Sort edges in ascending order of weight
         forest = UnionFind.initialise(n)
  3:
         T = (V, \emptyset)
   4:
         for each edge (u, v) in E do
  6:
             if forest.FIND(u) \neq forest.FIND(v) then
                                                               // Ignore edges that would create a cycle
                 forest. UNION(u, v)
  7:
                 T.add_edge(u, v)
  8:
  9:
         return T
                                         Time Complexity:
But this is not tight. A closer

 Sorting edges: θ(E log E)

look reveals that the total
                                            • E log E \leq E log V<sup>2</sup> = 2 E log V so θ(E log V)
cost of all UNION_SETS is
                                         o Initialization of union find: \theta(V)
\theta(V \log V).
                                         o For loop executes \theta(E) times

    SET_ID() takes θ(x) where x is height of the tree

                                               UNION SET() takes \theta(1) + 2 finds, so it takes \theta(x) where x is the height
                                               of the deeper of the two trees to be unioned (which could be at most V)
                                         Total cost: θ(EV)
```

## Complexity of UNION\_SETS

- We can show that, when using the union-by-size rule, the number of elements, N, in any tree is at least 2<sup>h</sup>, where h is the height of the tree.
- In other words, h ≤ log 2 N.
- Union needs 2 find operations + ⊕(1) effort.
- Find needs an effort equal to the height of the tree, which is ≤ log 2 N.
- So Union is Θ(log 2 N) where N is the number of nodes in the taller tree being unioned.
- We need to do V-1 unions in total.
- Each one has a worst case cost of ⊕(log V) (this is a significant overestimation, but its makes the maths easy).
- So the total cost of all unions is bounded by Θ(V log V).

## Kruskal's Algorithm: Complexity

```
Algorithm 41 Kruskal's algorithm
  1: function KRUSKAL(G = (V, E))
        sort(E, key((u, v)) = w(u, v))
                                                           // Sort edges in ascending order of weight
        forest = UnionFind.initialise(n)
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  4:
        for each edge (u, v) in E do
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                                                             // Ignore edges that would create a cycle
  6:
 7:
               forest. UNION(u, v)
                T.add_edge(u, v)
 8:
  9:
        return T
                                      Time Complexity:

 Sorting edges: θ(E log E)

                                         • E log E = E log V^2 = 2 E log V so \theta(E log V)
                                      o Initialization of union find: \theta(V)
                                      o For loop executes \theta(E) times
                                            The two finds take the same effort as the union, log(v)

 UNION takes θ(V log V) in total

                                      o Total cost: \theta(E log V + V log V) = \theta(E log V)
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```

## **Complexity of UNION\_SETS**

- We can improve the disjoint sets data structure significantly with 2 other optimisations:
  - Union by rank
  - Path compression
- These are discussed in the notes, but are not examinable.
- The complexity can be improved from V log V to Vα(V), where α denotes the inverse Ackermann function, an **extremely** slow growing function.
- Note:  $\alpha$ (any number which can be represented using the matter in the universe) < 5, so  $V\alpha(V)$  is effectively  $\theta(V)$ .

## **Kruskal's Algorithm: Correctness**

The proof of correctness for Kruskal's algorithm can be done following the same approach as Prim's one. Please try to write it down by yourself and check the details on the course notes.

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## **Comparison- Greedy Algorithms**

	Dijkstra	Prim's	Kruskal			
Purpose	Finds shortest path from a single source to all vertices	Finds Minimum Spanning Tree (MST)	Finds Minimum Spanning Tree (MST)			
Graph Type	Weighted (both directed & undirected)	Weighted, connected, undirected	Weighted, connected, undirected			
Approach	Greedy algorithm using priority queue	Greedy algorithm (like Dijkstra)	Greedy algorithm using sorting			
Process	Expands the closest vertex and updates distances	Expands the closest vertex and adds it to MST	Sorts edges & adds the smallest <b>non-cyclic</b> edge to MST			
Data Structures Used	Priority Queue (Min- Heap), Adjacency List/Matrix	Priority Queue (Min- Heap), Adjacency List/Matrix	Disjoint Set (Union- Find), Edge List			
Time Complexity	$O((V+E)\log V)$ with a priority queue (simplified to $O(E \log V)$ )	$O((V+E)\log V)$ with a priority queue (simplified to $O(E \log V)$ )	$O(V \log V)$ or $O(V\alpha V)$ where $\alpha$ is inverse Ackermann function			
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## Reading

- Course Notes: Chapter 6
- You can also check algorithms' textbooks for contents related to this lecture, e.g.:
  - o CLRS: Chapter 23, Section 24.3
  - KT: Chapter 4

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### **Concluding Remarks**

#### Take home message

- Dijkstra's algorithm is a greedy algorithm for determining shortest paths on graphs with non-negative weights.
- Prim's and Kruskal's algorithms are both greedy algorithms that correctly determine minimum spanning trees. While Prim's algorithm keeps growing the same connected component at each iteration; Kruskal's algorithm merges the two closest connected components.

#### Things to do (this list is not exhaustive)

- Make sure you understand:
  - The algorithms, and how to implement Union-Find data structure for Kruskal's algorithm.
  - How proofs of correctness for greedy algorithms work.

#### **Coming Up Next**

Dynamic Programming