COMMONWEALTH OF AUSTRALIA

Copyright Regulations 1969

WARNING

This material has been reproduced and communicated to you by or on behalf of the University of Sydney pursuant to Part VB of the Copyright Act 1968 (**the Act**). The material in this communication may be subject to copyright under the Act. Any further copying or communication of this material by you may be the subject of copyright protection under the Act.

Do not remove this notice.

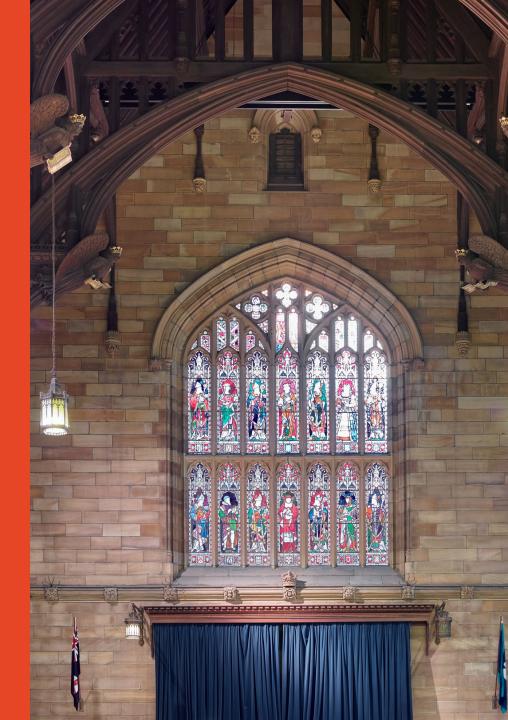
COMP2823

Lecture 8: Shortest Paths and Minimum Spanning Trees [GT 14.1-2, 15.1-3]

Joachim Gudmundsson School of Computer Science

Some content is taken from material provided by the textbook publisher Wiley.

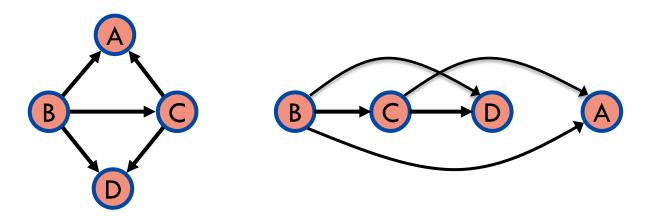




Last week

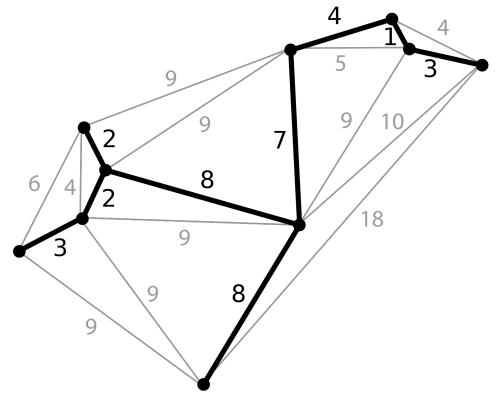
Introduced

- Unweighted graphs G=(V,E) directed and undirected
- Many new concepts: cycle, path, subgraphs, ...
- Graph ADT
- Graph traversal: BFS and DFS
 and some of its applications (cut set, topological sort, ...)



This week

- Weighted graphs G=(V,E) directed and undirected
- Shortest path using Dijkstra's algorithm
- Minimum spanning tree: Prim's and Kruskal's algorithm
- Briefly: Union find



Weighted Graphs

- In a weighted graph, each edge has an associated numerical value,
 called the weight of the edge
- Edge weights may represent, distances, costs, etc.

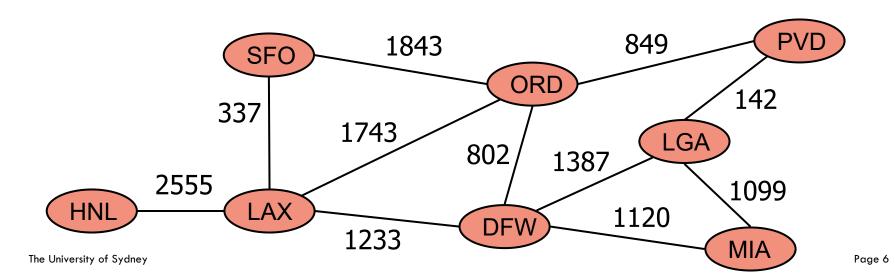
Example: In a road network, the weight of an edge represents the distance or time it takes to traverse the edge.



Weighted Graphs

- In a weighted graph, each edge has an associated numerical value,
 called the weight of the edge
- Edge weights may represent, distances, costs, etc.

Example: In a flight route graph, the weight of an edge represents the distance in miles between the endpoint airports

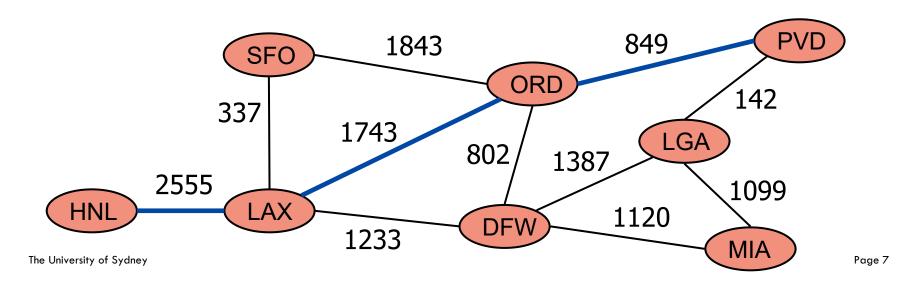


Shortest Paths

Given an edge weighted graph and two vertices \mathbf{u} and \mathbf{v} , we want to find a path of minimum total weight between \mathbf{u} and \mathbf{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.

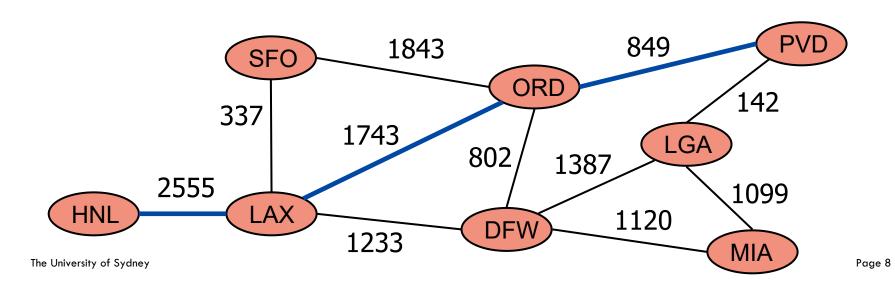
Example: Shortest path between Providence (PVD) and Honolulu (HNL)



Shortest Path Properties

Property: A subpath of a shortest path is itself a shortest path

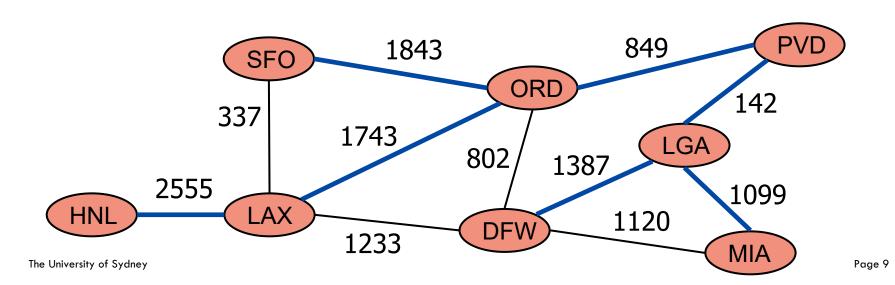
Example: Shortest path from Providence (PVD) to Honolulu (HNL) also contains a shortest path from Providence (PVD) to Los Angeles (LAX)



Shortest Path Properties

Property: There is a tree of shortest paths from a start vertex to all the other vertices (shortest path tree).

Example: Tree of shortest paths from Providence (PVD)



Input:

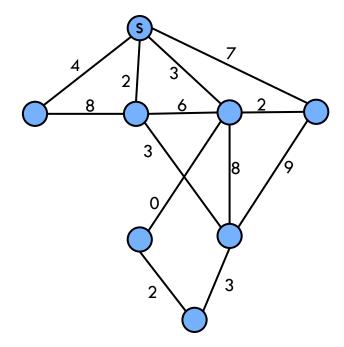
- Graph G = (V, E)
- Edges weights $w : E \rightarrow R_+$
- Start vertex s

Output:

- Distance from s to all v in V
- Shortest path tree rooted at s

Assumptions:

- G is connected and undirected
- edge weights are nonnegative



Input:

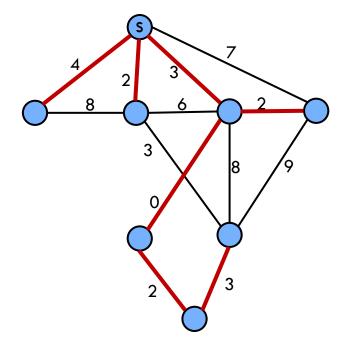
- Graph G = (V, E)
- Edges weights $w : E \rightarrow R_+$
- Start vertex s

Output:

- Distance from s to all v in V
- Shortest path tree rooted at s

Assumptions:

- G is connected and undirected
- edge weights are nonnegative



High level idea:

Keep track of a subset S of V s.t.

$$D[v] = dist_w(s, v)$$
 for all v in S

Maintain a distance estimate

$$D[v] \ge dist_w(s, v)$$
 for all v in V/S

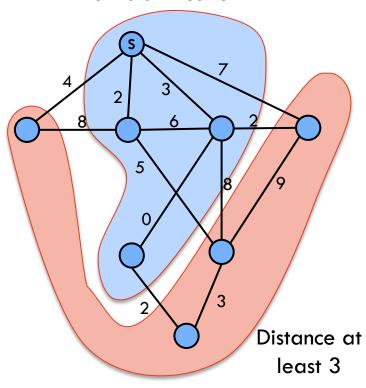
Initially:

- D[s] = 0
- D[v] = ∞ for all v in V s

In each iteration we:

- add to S vertex u in V \ S with smallest D[u]
- update D-values for vertices adjacent to u

Exact distances from s at most 3



High level idea:

Keep track of a subset S of V s.t.

$$D[v] = dist_w(s, v)$$
 for all v in S

Maintain a distance estimate

$$D[v] \ge dist_w(s, v)$$
 for all v in V/S

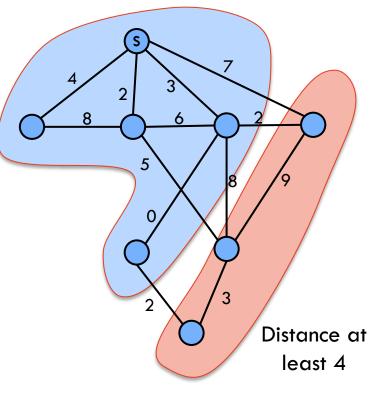
Initially:

- D[s] = 0
- D[v] = ∞ for all v in V s

In each iteration we:

- add to S vertex u in V \ S with smallest D[u]
- update D-values for vertices adjacent to u

Exact distances from s at most 4



High level idea:

Keep track of a subset S of V s.t.

$$D[v] = dist_w(s, v)$$
 for all v in S

Maintain a distance estimate

$$D[v] \ge dist_w(s, v)$$
 for all v in V/S

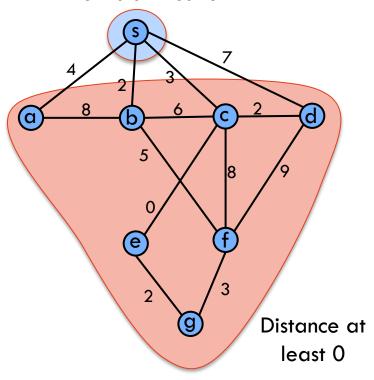
Initially:

- D[s] = 0
- D[v] = ∞ for all v in V s

In each iteration we:

- add to S vertex u in V \ S with smallest D[u]
- update D-values for vertices adjacent to u

Exact distances from s at most 0



High level idea:

- Keep track of a subset S of V s.t.
 - $D[v] = dist_w(s, v)$ for all v in S
- Maintain a distance estimate

$$D[v] \ge dist_w(s, v)$$
 for all v in V/S

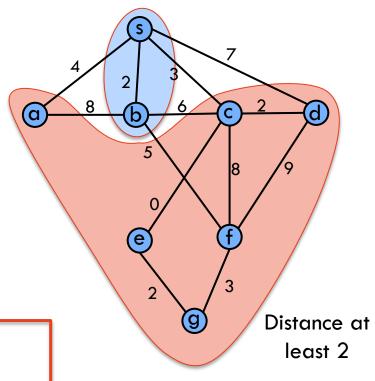
Initially:

- D[s] = 0
- D[v] = ∞ for all v in V s

In each iteration we:

- add to S vertex u in V \ S with smallest D[u]
- update D-values for vertices adjacent to u

Exact distances from s at most 2



High level idea:

- Keep track of a subset S of V s.t.
 - $D[v] = dist_w(s, v)$ for all v in S
- Maintain a distance estimate

$$D[v] \ge dist_w(s, v)$$
 for all v in V/S

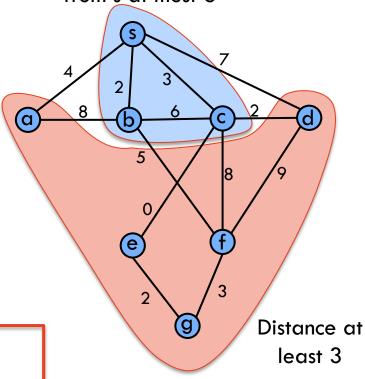
Initially:

- D[s] = 0
- D[v] = ∞ for all v in V s

In each iteration we:

- add to S vertex u in V \ S with smallest D[u]
- update D-values for vertices adjacent to u

Exact distances from s at most 3



High level idea:

Keep track of a subset S of V s.t.

$$D[v] = dist_w(s, v)$$
 for all v in S

Maintain a distance estimate

$$D[v] \ge dist_w(s, v)$$
 for all v in V/S

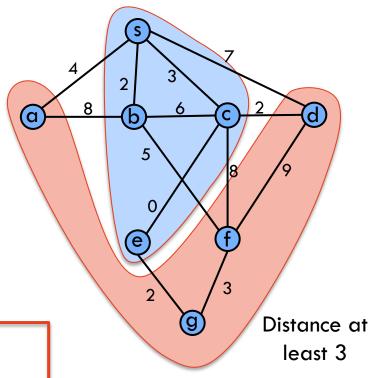
Initially:

- D[s] = 0
- D[v] = ∞ for all v in V s

In each iteration we:

- add to S vertex u in V \ S with smallest D[u]
- update D-values for vertices adjacent to u

Exact distances from s at most 3



High level idea:

Keep track of a subset S of V s.t.

$$D[v] = dist_w(s, v)$$
 for all v in S

Maintain a distance estimate

$$D[v] \ge dist_w(s, v)$$
 for all v in V/S

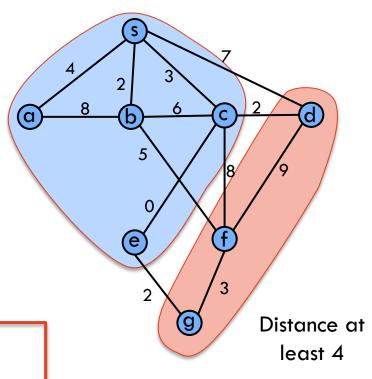
Initially:

- D[s] = 0
- D[v] = ∞ for all v in V s

In each iteration we:

- add to S vertex u in V \ S with smallest D[u]
- update D-values for vertices adjacent to u

Exact distances from s at most 4



High level idea:

Keep track of a subset S of V s.t.

$$D[v] = dist_w(s, v)$$
 for all v in S

Maintain a distance estimate

$$D[v] \ge dist_w(s, v)$$
 for all v in V/S

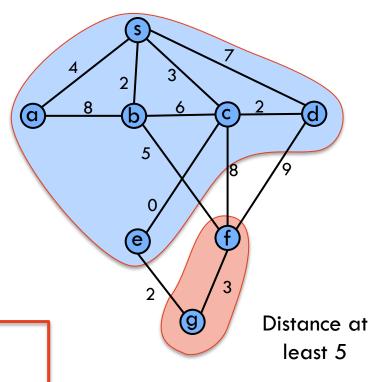
Initially:

- D[s] = 0
- D[v] = ∞ for all v in V s

In each iteration we:

- add to S vertex u in V \ S with smallest D[u]
- update D-values for vertices adjacent to u

Exact distances from s at most 5



High level idea:

Keep track of a subset S of V s.t.

$$D[v] = dist_w(s, v)$$
 for all v in S

Maintain a distance estimate

$$D[v] \ge dist_w(s, v)$$
 for all v in V/S

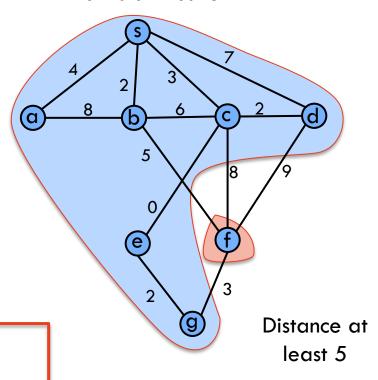
Initially:

- D[s] = 0
- D[v] = ∞ for all v in V s

In each iteration we:

- add to S vertex u in V \ S with smallest D[u]
- update D-values for vertices adjacent to u

Exact distances from s at most 5



High level idea:

Keep track of a subset S of V s.t.

$$D[v] = dist_w(s, v)$$
 for all v in S

Maintain a distance estimate

$$D[v] \ge dist_w(s, v)$$
 for all v in V/S

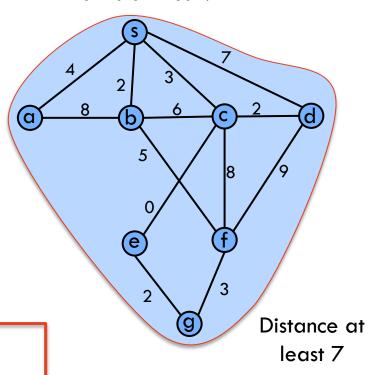
Initially:

- D[s] = 0
- D[v] = ∞ for all v in V s

In each iteration we:

- add to S vertex u in V \ S with smallest D[u]
- update D-values for vertices adjacent to u

Exact distances from s at most 7



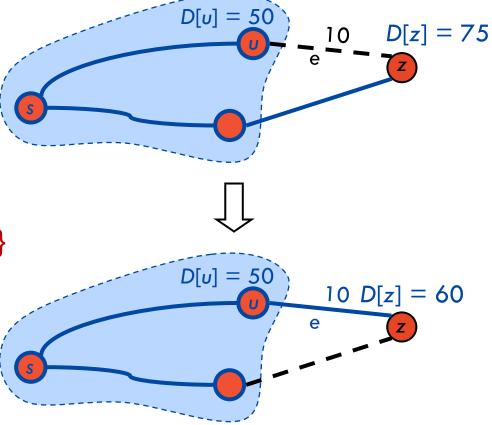
Edge Relaxation

Consider edge e = (u, z) such that:

- u is the last vertex added to \$
- z is not in \$

The relaxation of edge (u, z) updates D[z] as follows:

$$D[z] \leftarrow \min\{D[z], D[u] + w(u, z)\}$$



Dijkstra's Algorithm pseudocode

```
def Dijkstra(G, w, s):

# initialize algorithm

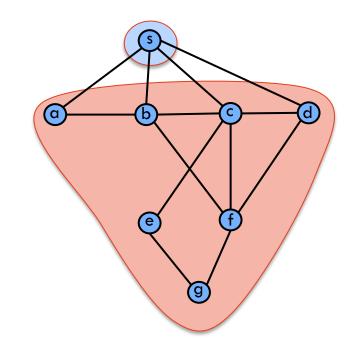
for v in V do

D[v] \leftarrow \infty

parent[v] \leftarrow \emptyset

D[s] \leftarrow 0

Q \leftarrow new priority queue for { (v, D[v]) : v in V }
```

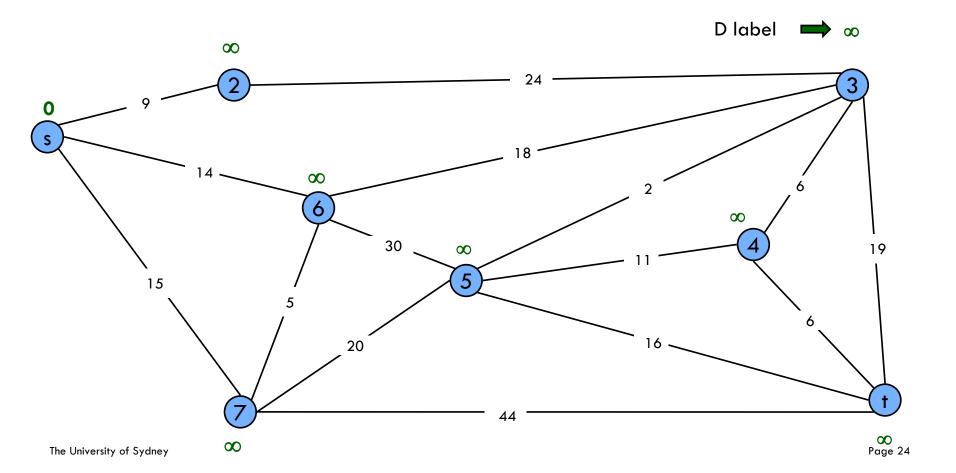


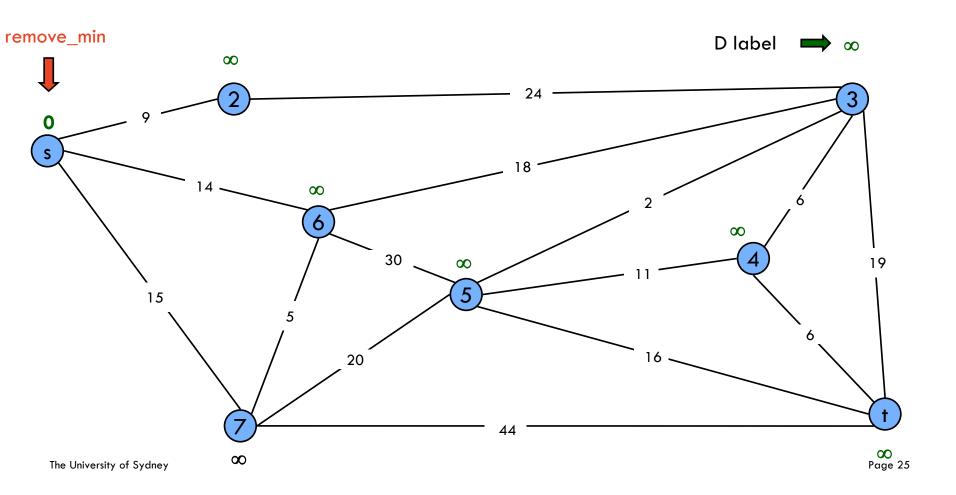
```
# iteratively add vertices to S
while Q is not empty do
  u ← Q.remove_min()
for z in G.neighbors(u) do
  if D[u] + w[u, z] < D[z] then
   D[z] ← D[u] + w[u, z]
   Q.update_priority(z, D[z])
   parent[z] ← u
return D, parent</pre>
```

Recall: heap-based priority queue (Lect 5)

Operation	Time
size, isEmpty	O(1)
min,	O(1)
insert	$O(\log n)$
removeMin	$O(\log n)$

+ update key O(log n)

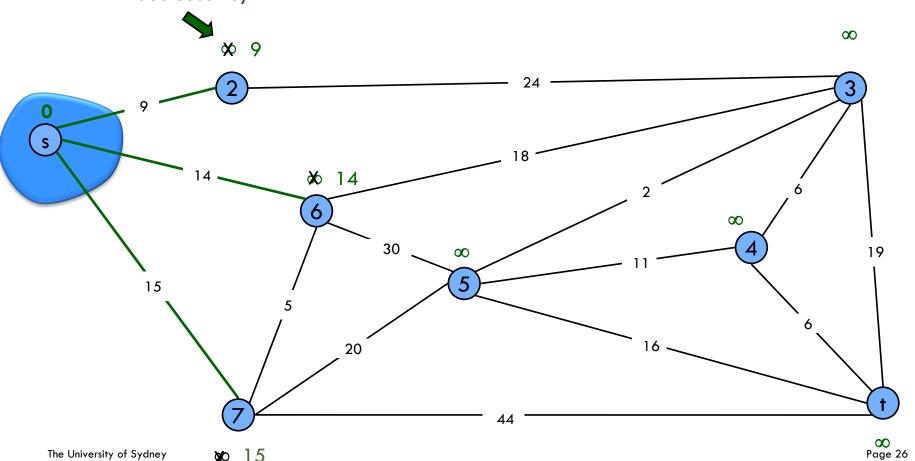


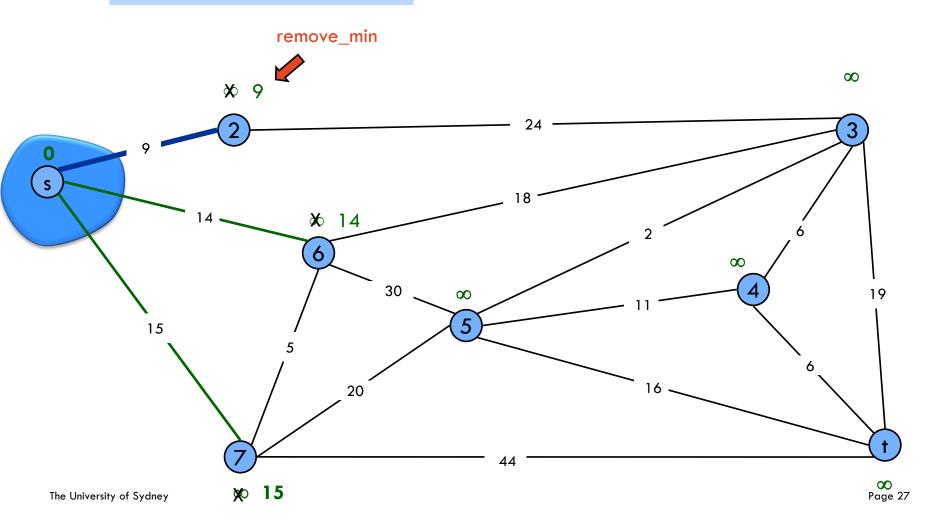


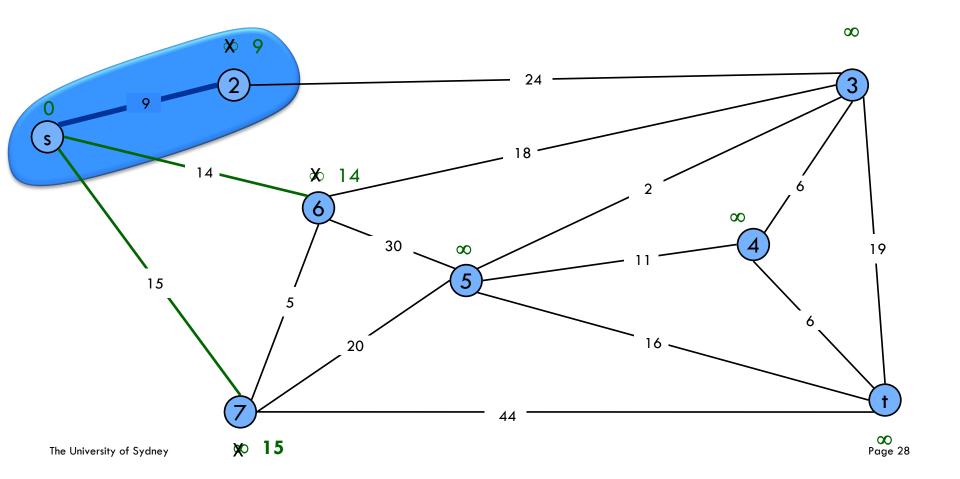
15

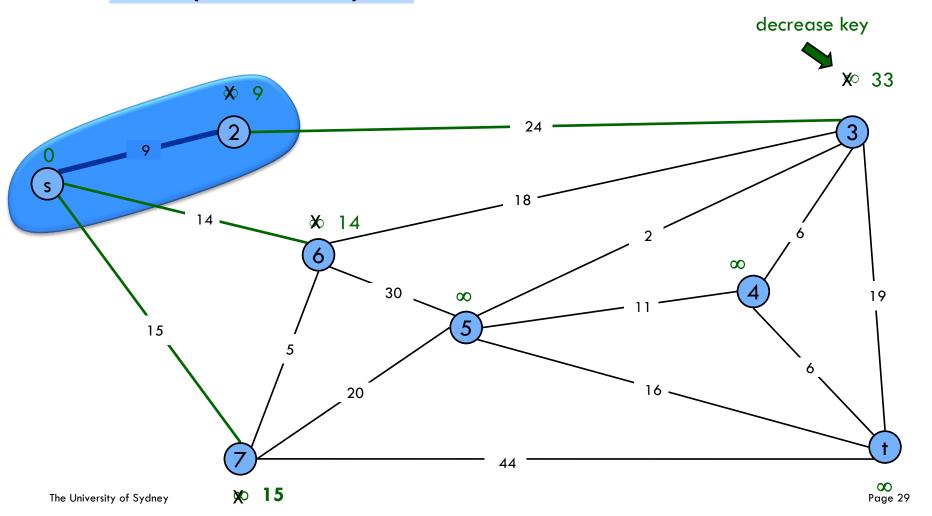
The University of Sydney

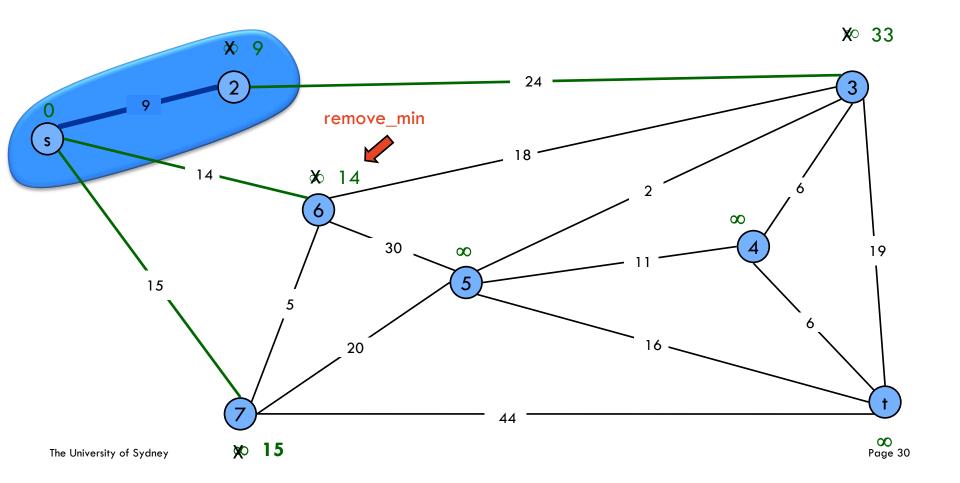
decrease key

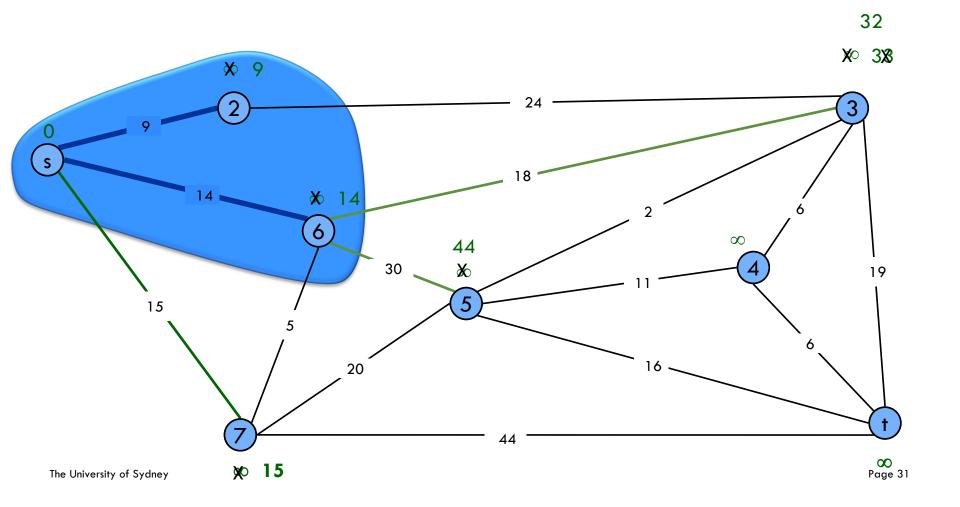


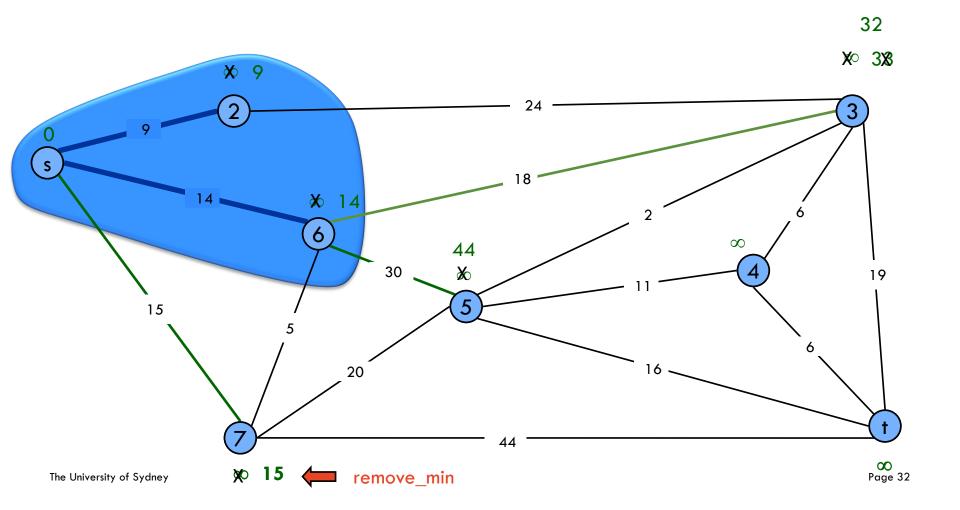


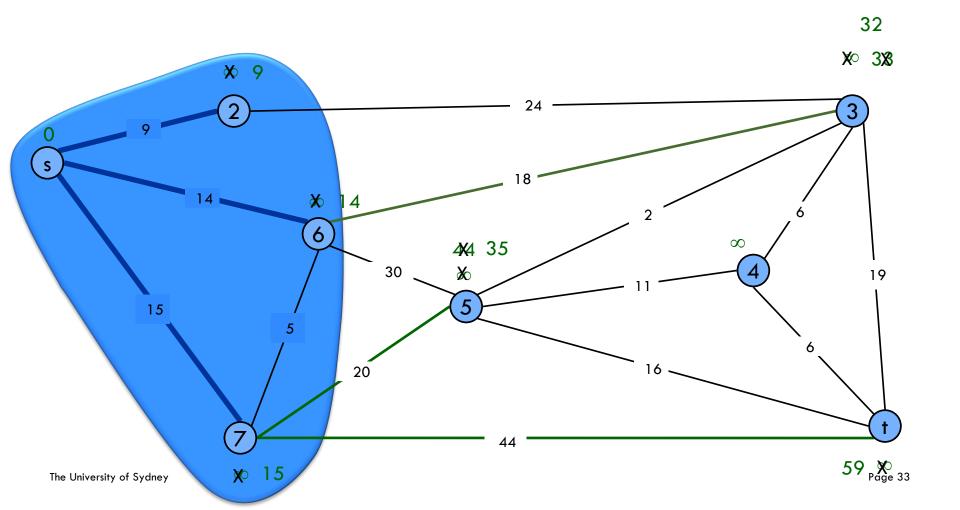


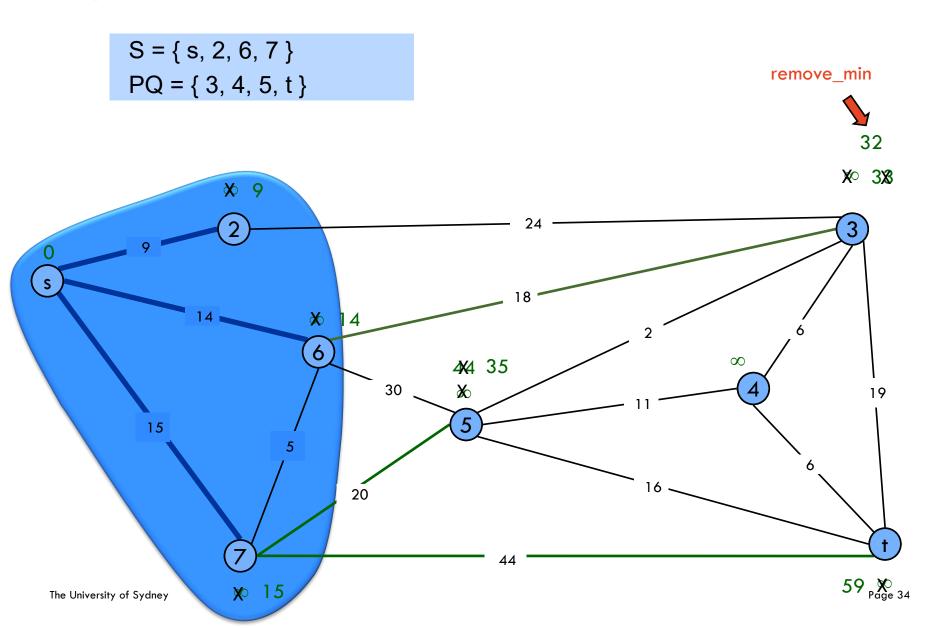


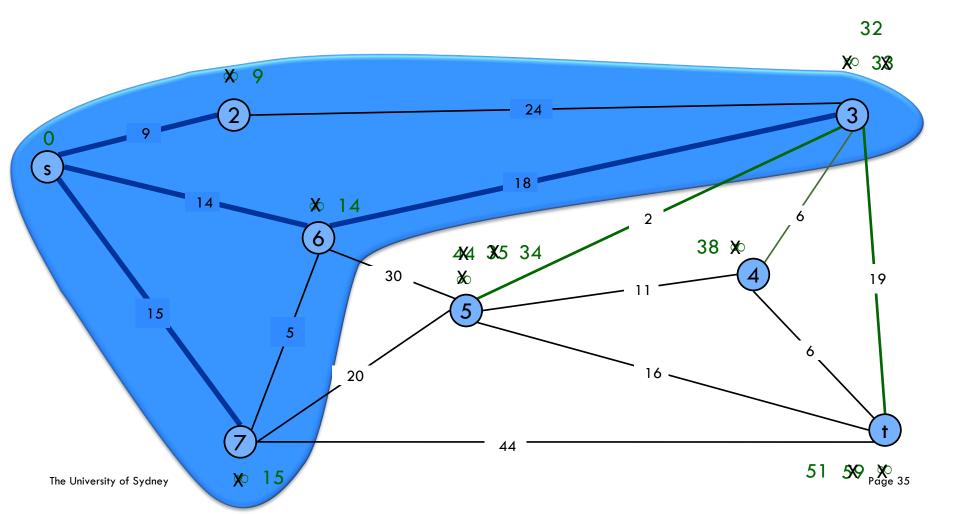


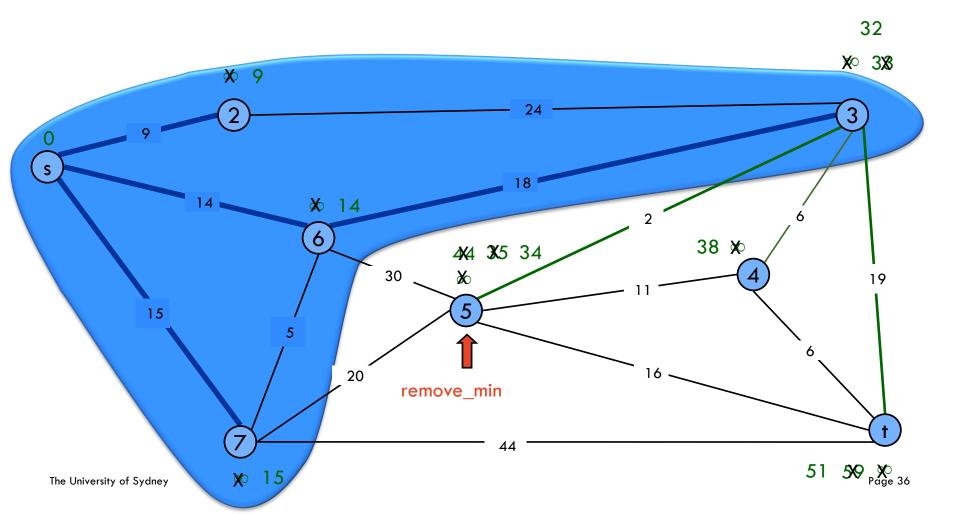


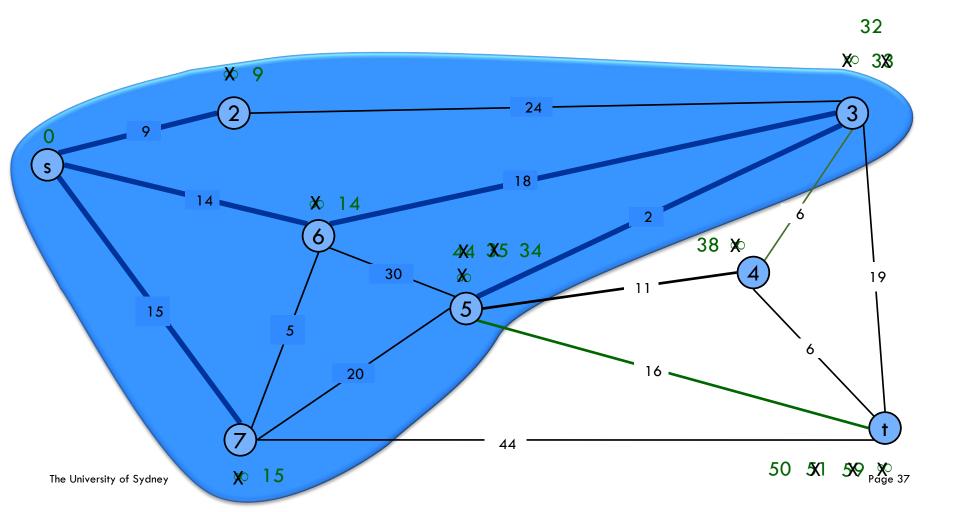


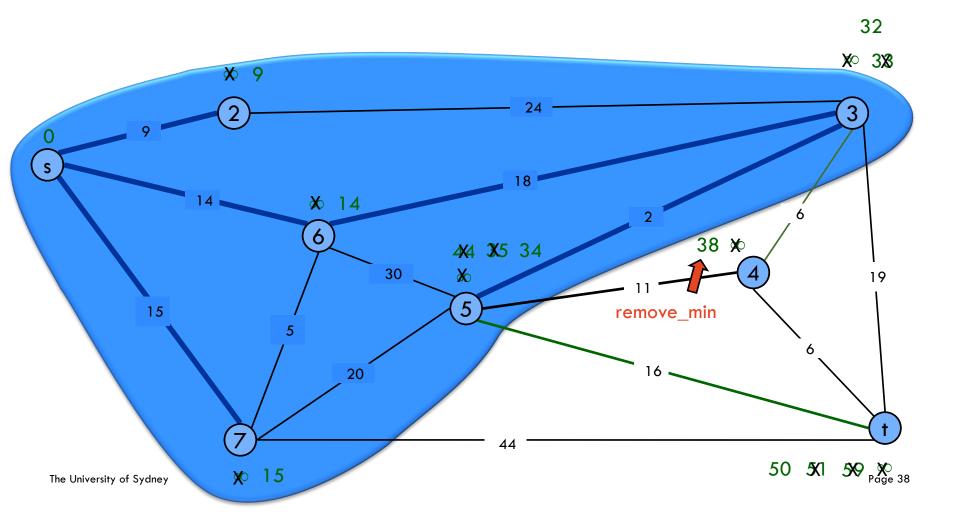


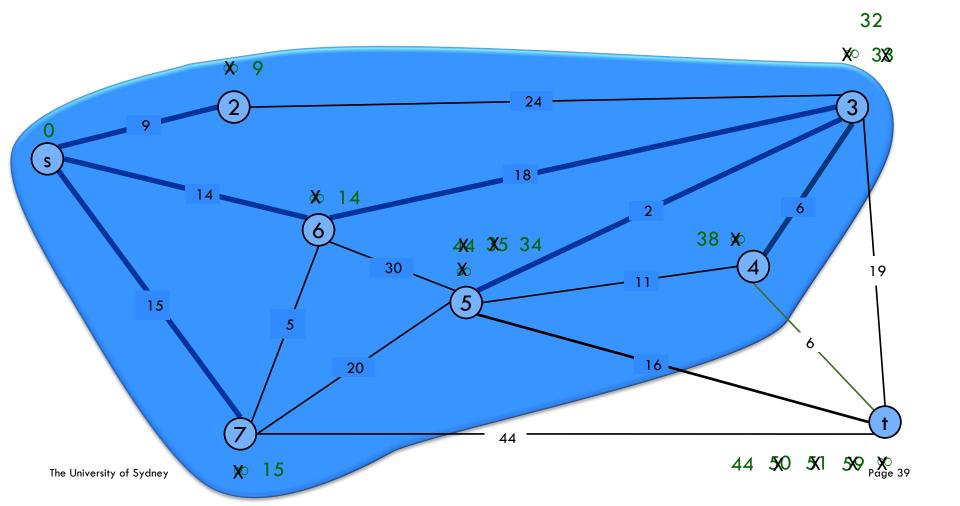


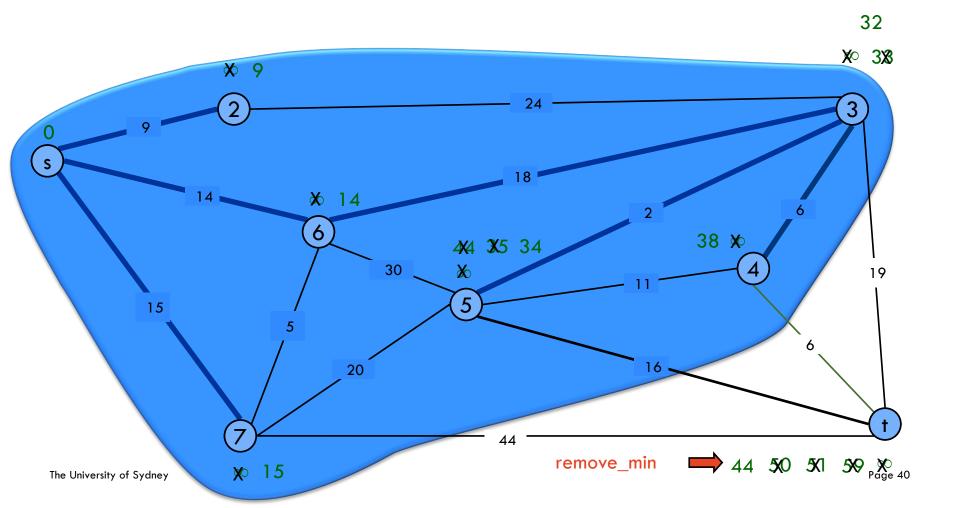






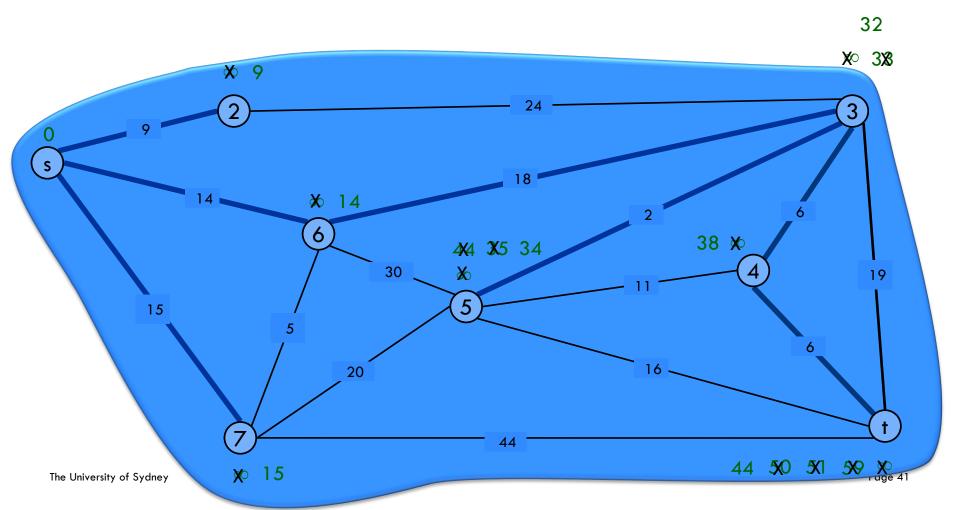






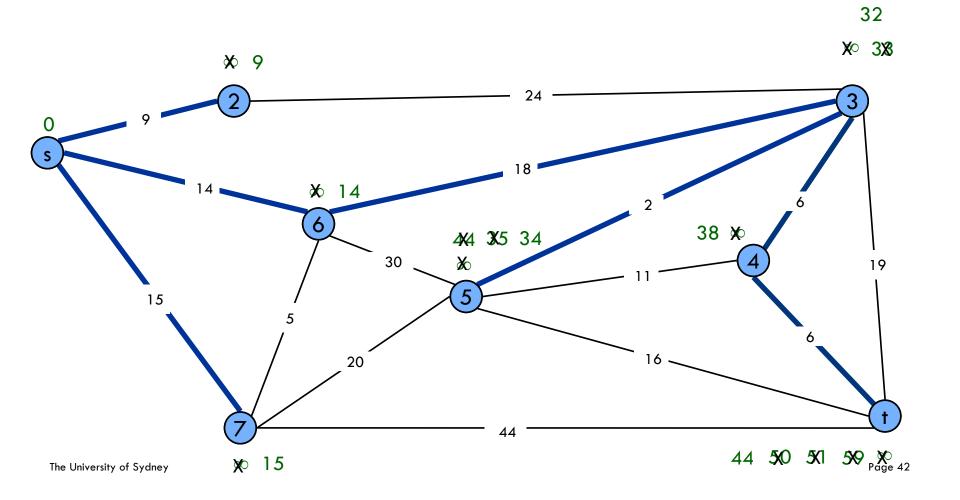
$$S = \{ s, 2, 3, 4, 5, 6, 7, t \}$$

PQ = { }



$$S = \{ s, 2, 3, 4, 5, 6, 7, t \}$$

PQ = { }



Dijkstra complexity analysis except PQ ops

```
def Dijkstra(G, w, s):
 # initialize algorithm
 for v in V do
  D[v] \leftarrow \infty
  parent[v] \leftarrow \emptyset
 D[s] \leftarrow 0
 Q \leftarrow \text{new priority queue for } \{ (v, D[v]) : v \text{ in } V \}
                                                               O(n) insert operations
 # iteratively add vertices to S
 while Q is not empty do
                                           O(n) remove_min operations
  u \leftarrow Q.remove min()
  for z in G.neighbors(u) do
   if D[u] + w[u, z] < D[z] then
                                          O(deg(u)) times for each u in V
    D[z] \leftarrow D[u] + w[u,z]
                                           plus update_priority work
    Q.update_priority(z, D[z])
                                           ⇒ O(m) PQ update operations (decrease_key)
    parent[z] \leftarrow u
 return D, parent
```

[Dijkstra 1956]

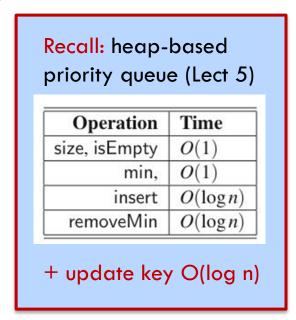
Dijkstra's Algorithm complexity analysis

Assuming the graph is connected (so $m \ge n-1$), the algorithm spends O(m) time on everything except PQ operations

Priority queue operation counts:

- insert: n
- decrease_key: m
- remove_min: n

Fact: Using a heap for PQ, Dijkstra runs in O(m log n) time



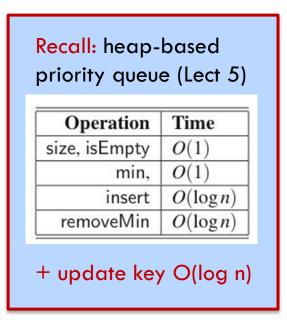
Dijkstra's Algorithm complexity analysis

Assuming the graph is connected (so $m \ge n-1$), the algorithm spends O(m) time on everything except PQ operations

Priority queue operation counts:

- insert: n
- decrease_key: m
- remove_min: n

Fact: Using a heap for PQ, Dijkstra runs in O(m log n) time



Fibonacci heaps is a PQ that can carry out decrease key in O(1) amortized time. Using that instead we get $O(m + n \log n)$ time.

Dijkstra's Algorithm Correctness

Invariant: For each $u \in S = V \setminus Q$, we have $D[u] = dist_w(s, u)$

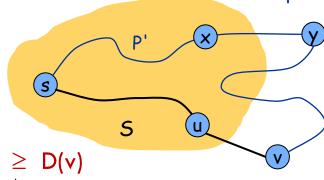
Proof: (by induction on |S|)

Base case: |S| = 1 is trivial since D[s] = 0

Inductive hypothesis: Assume true for $|S| = k \ge 1$.

Induction step:

- Assume v is next node added to S and u=parent[v]
- Assume the shortest s-u path plus (u, v) is an s-v path of length D[v]
- Consider any s-v path P. We'll see that it's no shorter than D[v].
- Let x-y be the first edge in P that leaves S,
 and let P' be the subpath from s to x.
- P is already too long as soon as it leaves S:



$$w(P) \ge w(P') + w(x,y) = D(x) + w(x,y) \ge D(y) \ge D(v)$$
 $\uparrow \qquad \qquad \uparrow \qquad \uparrow$

inductive Def of Dijkstra chose v

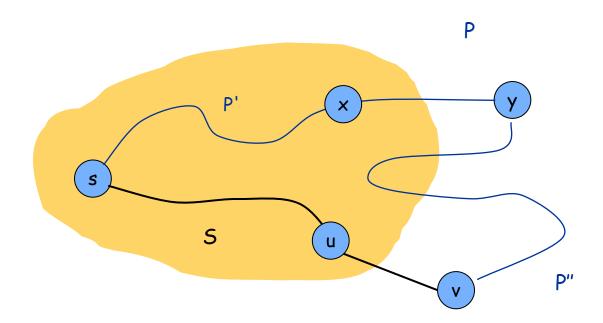
hypothesis

The University of Sydney

D(y) instead of y

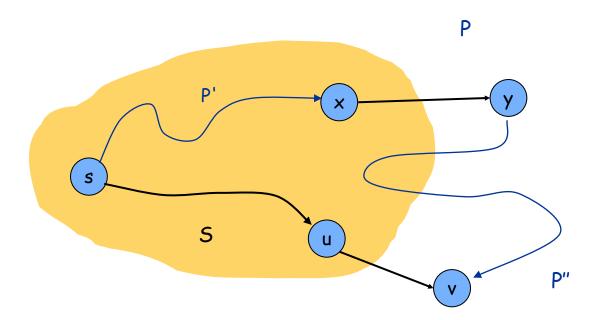
Warning: Dijkstra may not work for negative-weight edges

In the proof of correctness, even if D[v] is the smallest label, it may be that $dist_w(s, v) < D[v]$ if w(P'') < 0



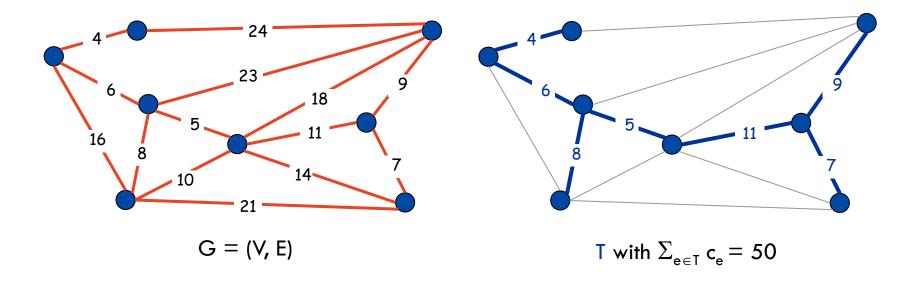
Dijkstra works for directed graph

In the proof of correctness, we need to use directed paths instead



Minimum Spanning Tree (MST)

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.



Applications

MST is fundamental problem with diverse applications.

Network design: Telephone, electrical, hydraulic, TV cable, computer, road

Approximation algorithms for NP-hard problems: traveling salesperson problem, Steiner tree

Indirect applications.

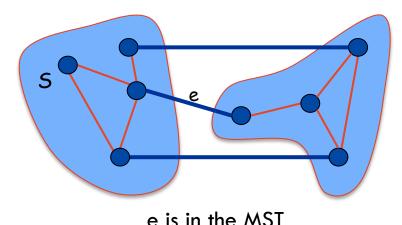
- max bottleneck paths
- LDPC codes for error correction
- image registration with Renyi entropy
- learning salient features for real-time face verification
- reducing data storage in sequencing amino acids in a protein

– ...

MST properties

Simplifying assumption. All edge costs c_e are distinct.

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.



MST properties

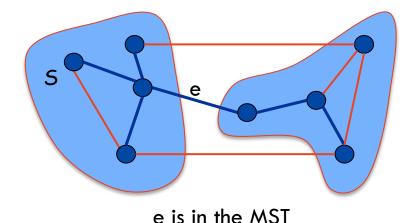
Simplifying assumption. All edge costs c_e are distinct.

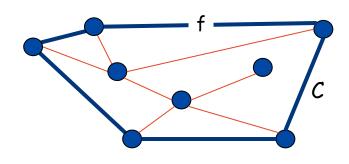
Cut property. Let S be any subset of nodes, and let e be the min cost edge with exactly one endpoint in S. Then the MST contains e.

[Let's try to prove it]

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.

[Obvious]

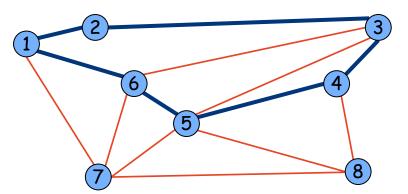




f is not in the MST

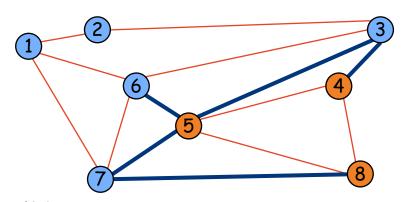
Cycles and Cuts

Cycle. Set of edges of the form a-b, b-c, c-d, ..., y-z, z-a.



Cycle C = 1-2, 2-3, 3-4, 4-5, 5-6, 6-1

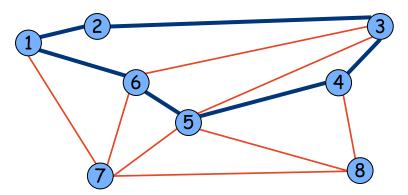
Cutset. A cut is a subset of nodes S. The corresponding cutset D is the subset of edges with exactly one endpoint in S.



Cut S = $\{4, 5, 8\}$ Cutset D = 5-6, 5-7, 3-4, 3-5, 7-8

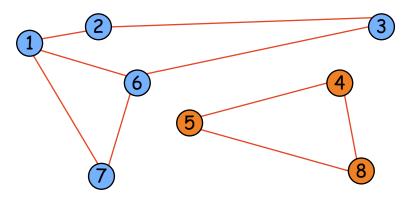
Cycles and Cuts

Cycle. Set of edges of the form a-b, b-c, c-d, ..., y-z, z-a.



Cycle C = 1-2, 2-3, 3-4, 4-5, 5-6, 6-1

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



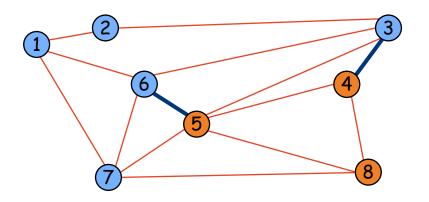
Cut S =
$$\{4, 5, 8\}$$

Cutset D = 5-6, 5-7, 3-4, 3-5, 7-8

Removal of the cutset separates the cut from the rest of the graph

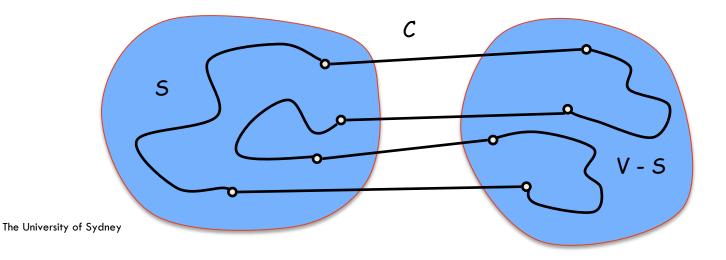
Cycle-Cut Intersection

Claim. A cycle and a cutset intersect in an even number of edges.

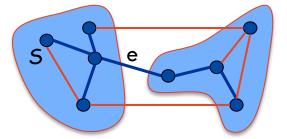


Cycle C = 1-2, 2-3, 3-4, 4-5, 5-6, 6-1 Cutset D = 3-4, 3-5, 5-6, 5-7, 7-8 Intersection = 3-4, 5-6

Proof:



Proving the Cut Property

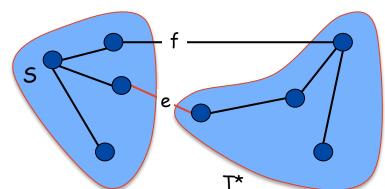


Simplifying assumption. All edge costs c_e are distinct.

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.

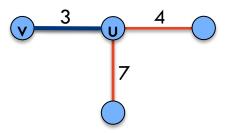
Proof: (exchange argument)

- Let T* be MST and suppose e does not belong to T*
- Adding e to T* creates a cycle C in T*
- Edge e is both in the cycle C and in the cutset D corresponding to S \Rightarrow there exists another edge, say f, that is in both C and D [cycle-cut inters.].
- $T' = T^* \cup \{e\} \{f\}$ is also a spanning tree.
- Since $c_e < c_f$, $cost(T') < cost(T^*)$.
- A contradiction, so e must belong in T*



Prim's Algorithm

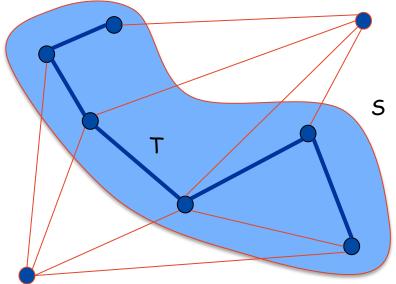
```
def prim(G, c):
    u ← arbitrary vertex in V
    S ← { u }
    T ← Ø
    while |S| < |V| do
        (u, v) ← min cost edge s.t. u in S and v not in S
        add (u, v) to T
        add v to S
    return T</pre>
```



Prim's Algorithm: Correctness

```
def prim(G, c):
    u ← arbitrary vertex in V
    S ← { u }
    T ← Ø
    while |S| < |V| do
        (u, v) ← min cost edge s.t. u in S and v not in S add (u, v) to T
        add v to S
    return T</pre>
```

Every time we add an edge we follow cut property! (add shortest cut edge between S and V/S)

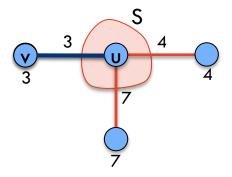


Implementation: Prim's Algorithm

```
def prim(G, c) {
 for v in V do
  d[v] \leftarrow \infty
   parent[v] \leftarrow \emptyset
 u ← arbitrary vertex in V
 d[u] \leftarrow 0
 Q \leftarrow \text{new } PQ \text{ with items } \{ (v, d[v]) \text{ for } v \text{ in } V \}
 S \leftarrow \emptyset
 while Q is not empty do
   u \leftarrow delete min element from Q
   add u to S
  for (u, v) incident to u do
    if v \notin S and c_{u,v} < d[v] then
      parent[v] \leftarrow u
      decrease priority d[v] to c_{u,v}
 return parent
```

Main idea: for every v in $V \setminus S$ we keep

- d[v] = distance to closest neighbor in S
- parent[v] = closest neighbor in S

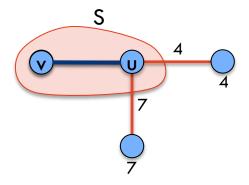


Implementation: Prim's Algorithm

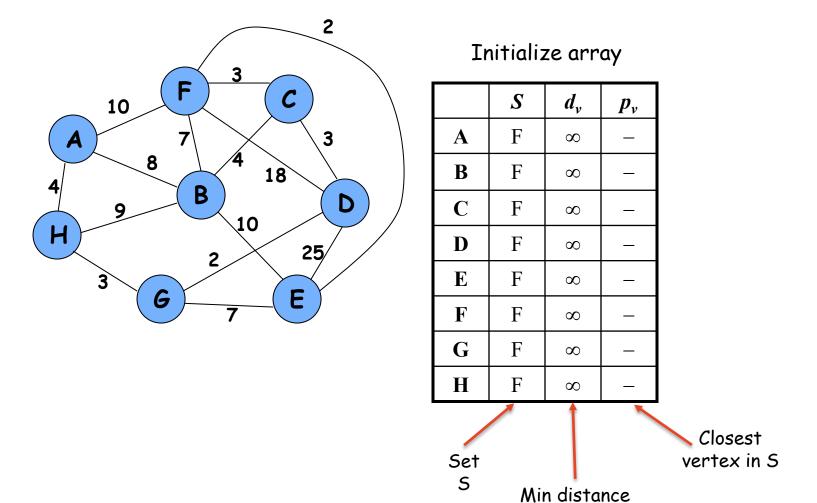
```
def prim(G, c) {
 for v in V do
  d[v] \leftarrow \infty
   parent[v] \leftarrow \emptyset
 u ← arbitrary vertex in V
 d[u] \leftarrow 0
 Q \leftarrow \text{new } PQ \text{ with items } \{ (v, d[v]) \text{ for } v \text{ in } V \}
 S \leftarrow \emptyset
 while Q is not empty do
   u \leftarrow delete min element from Q
   add u to S
  for (u, v) incident to u do
    if v \notin S and c_{u,v} < d[v] then
      parent[v] \leftarrow u
      decrease priority d[v] to c_{u,v}
 return parent
```

Main idea: for every v in $V \setminus S$ we keep

- d[v] = distance to closest neighbor in S
- parent[v] = closest neighbor in S

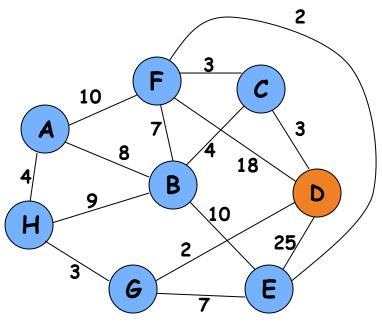


Walk-Through



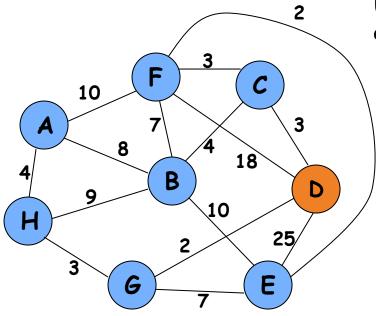
The University of Sydney

to S



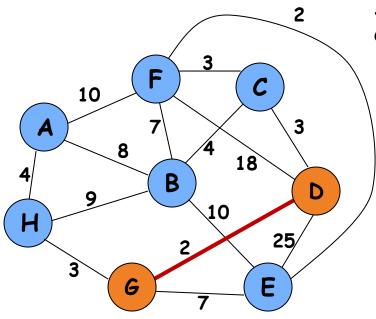
Start with any node, say D

	S	d_v	p_{v}
A			
В			
C			
D	T	0	_
E			
F			
G			
Н			



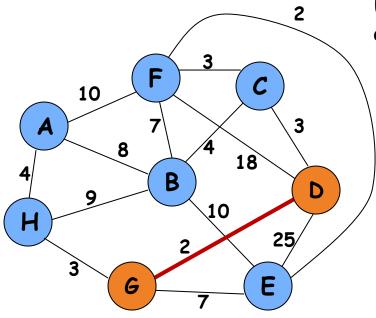
Update distances of adjacent, unselected nodes

	S	d_v	p_{v}
A			
В			
C		3	D
D	Т	0	_
E		25	D
F		18	D
G		2	D
Н			



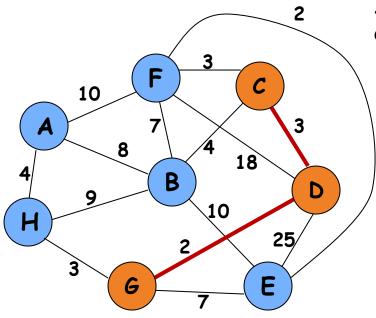
Select node with minimum distance

	S	d_v	p_{v}
A			
В			
C		3	D
D	Т	0	l
E		25	D
F		18	D
G	T	2	D
Н			



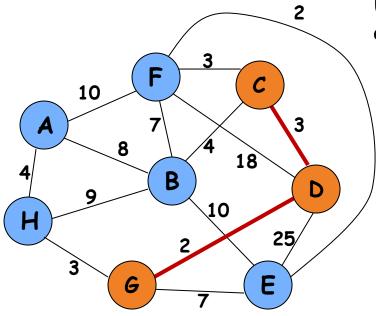
Update distances of adjacent, unselected nodes

	S	d_v	p_{v}
A			
В			
C		3	D
D	T	0	_
E		7	G
F		18	D
G	T	2	D
Н		3	G



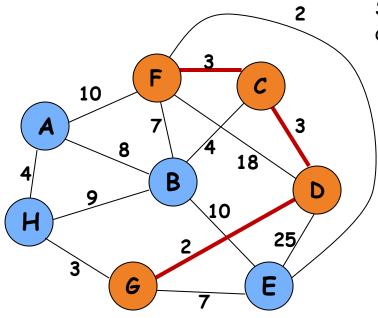
Select node with minimum distance

	S	d_v	p_{v}
A			
В			
С	T	3	D
D	T	0	_
E		7	G
F		18	D
G	T	2	D
Н		3	G



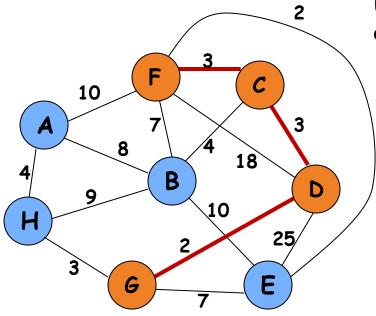
Update distances of adjacent, unselected nodes

	S	d_v	p_{v}
A			
В		4	C
C	T	3	D
D	T	0	
E		7	G
F		3	C
G	T	2	D
Н		3	G



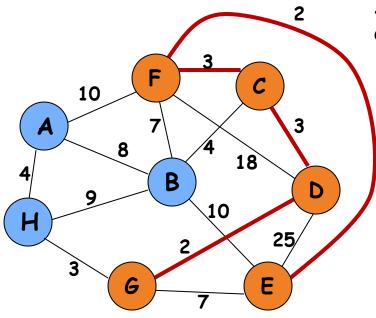
Select node with minimum distance

	S	d_v	p_{v}
A			
В		4	C
C	Т	3	D
D	T	0	_
E		7	G
F	T	3	С
G	T	2	D
Н		3	G



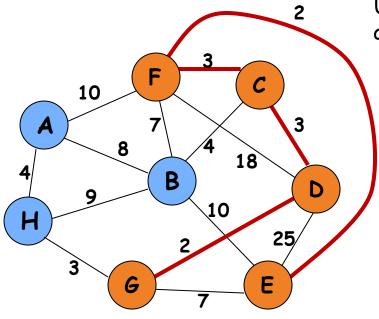
Update distances of adjacent, unselected nodes

	S	d_v	p_{v}
A		10	F
В		4	C
C	Т	3	D
D	T	0	_
E		2	F
F	T	3	С
G	T	2	D
Н		3	G



Select node with minimum distance

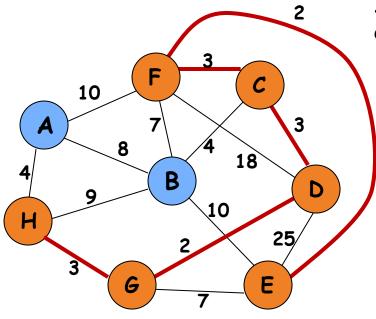
	S	d_v	p_{v}
A		10	F
В		4	C
C	Т	3	D
D	T	0	_
E	T	2	F
F	T	3	С
G	T	2	D
Н		3	G



Update distances of adjacent, unselected nodes

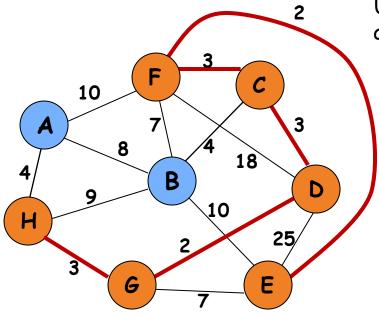
	S	d_v	p_{v}
A		10	F
В		4	С
C	Т	3	D
D	Т	0	_
E	T	2	F
F	T	3	С
G	Т	2	D
Н		3	G

Table entries unchanged



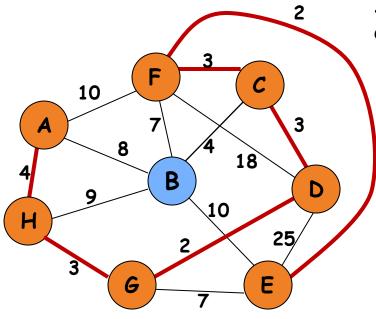
Select node with minimum distance

	S	d_v	p_{v}
A		10	F
В		4	С
C	Т	3	D
D	T	0	_
E	T	2	F
F	T	3	С
G	T	2	D
Н	T	3	G



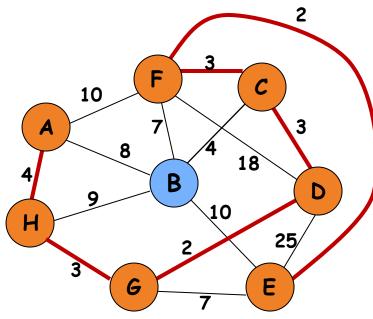
Update distances of adjacent, unselected nodes

	S	d_v	p_{v}
A		4	Н
В		4	C
C	Т	3	D
D	T	0	_
E	T	2	F
F	T	3	С
G	T	2	D
Н	T	3	G



Select node with minimum distance

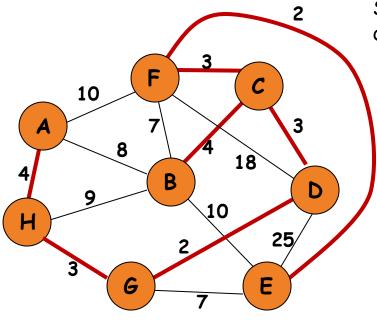
	S	d_v	p_{v}
A	T	4	Н
В		4	С
C	T	3	D
D	T	0	_
E	T	2	F
F	T	3	С
G	T	2	D
Н	T	3	G



Update distances of adjacent, unselected nodes

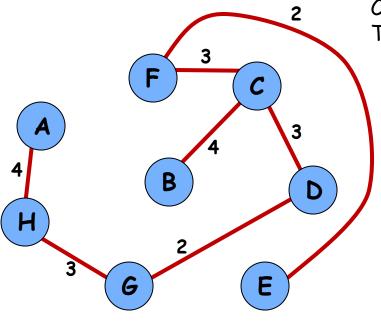
	S	d_v	p_{v}
A	T	4	Н
В		4	C
C	Т	3	D
D	T	0	_
E	T	2	F
F	T	3	С
G	T	2	D
Н	T	3	G

Table entries unchanged



Select node with minimum distance

	S	d_v	p_{v}
A	T	4	Н
В	T	4	C
C	T	3	D
D	T	0	_
E	T	2	F
F	T	3	С
G	T	2	D
Н	Т	3	G



Cost of Minimum Spanning Tree = $\sum d_v = 21$

	S	d_v	p_{v}
A	T	4	Н
В	T	4	C
C	Т	3	D
D	T	0	_
E	T	2	F
F	T	3	С
G	T	2	D
Н	T	3	G

Done!

Prim's Algorithm complexity

```
def prim(G, c) {
 for v in V do
   d[v] \leftarrow \infty
   parent[v] \leftarrow \emptyset
 u ← arbitrary vertex in V
 d[u] \leftarrow 0
 Q \leftarrow \text{new PQ with items } \{ (v, d[v]) \text{ for } v \text{ in } V \}
 S \leftarrow \emptyset
 while Q is not empty do
   u \leftarrow delete min element from Q
   S \leftarrow S \cup \{u\}
   for (u, v) incident to u do
    if v \notin S and c_{u,v} < d[v] then
      parent[v] \leftarrow u
      decrease priority d[v] to c<sub>uv</sub>
 return parent
```

Similar analysis to Dijkstra's algorithm:

- O(m log n) using a heap
- O(m + n log n) using Fibonacci heap

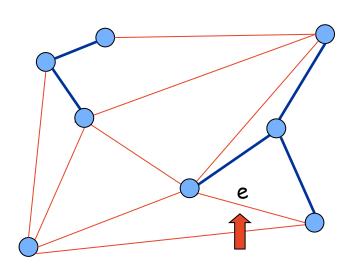
[Jarnik1930s and Prim 1957]

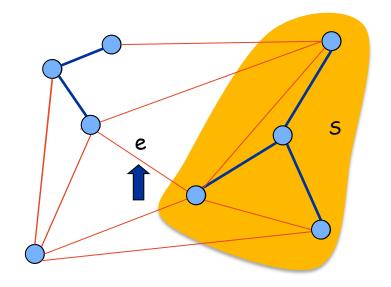
Kruskal's Algorithm

Consider edges in ascending order of weight.

Case 1: If adding e to T creates a cycle, discard e according to cycle property.

Case 2: Otherwise, insert e = (u, v) into T according to cut property where S = set of nodes in u's connected component.



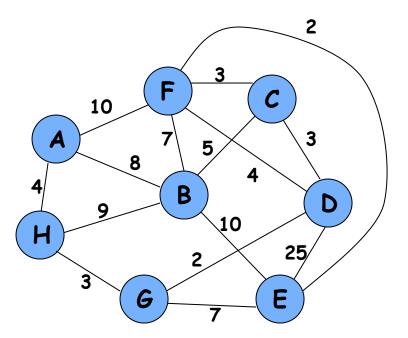


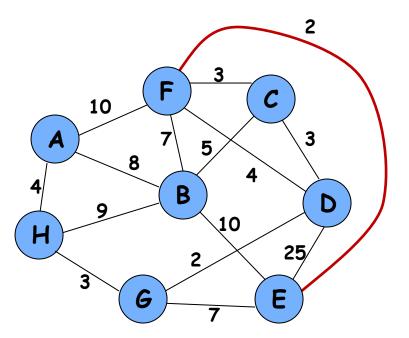
The University of Sydney

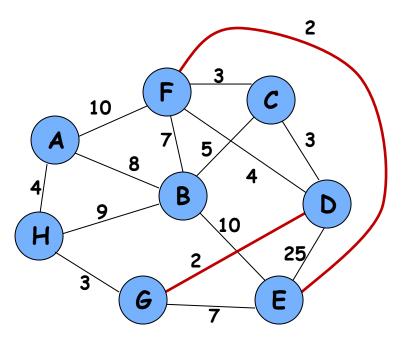
Case 1

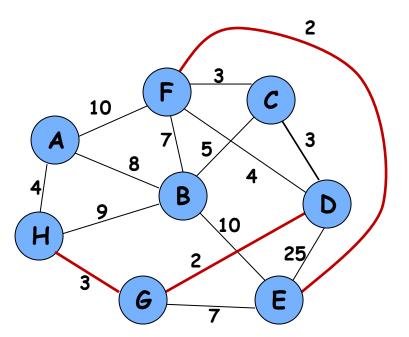
Case 2

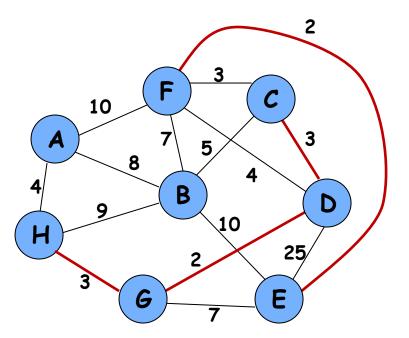
Page 79

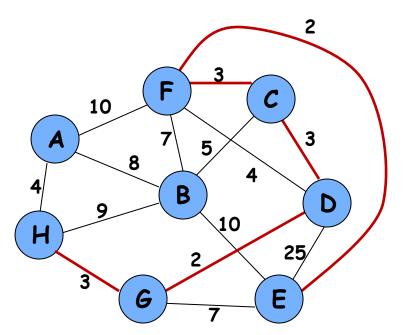


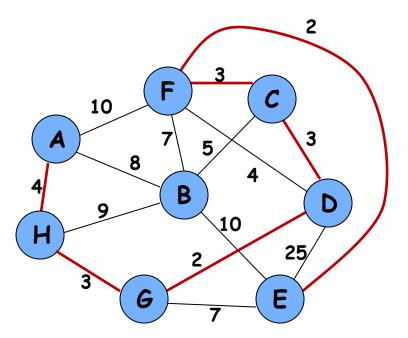


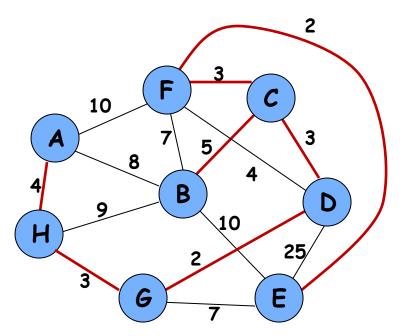












Kruskal's Algorithm: Time complexity

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

One option is to 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

Can we do better?

Yes, keep track of the connected components with a data structure

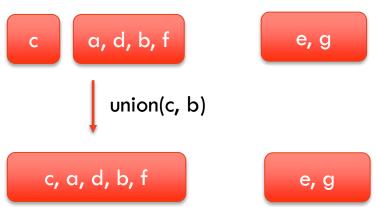
Union Find ADT

Data structure defined on a ground set of elements A

Used to keep track of an evolving partition of A

Supported operations:

- a b c d e f g
- make_sets(A): makes | 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



Kruskal's algorithm implementation

```
def Kruskal(G,c):
    sort E in increasing c-value
    answer ← []
    comp ← make_sets(V)
    for (u,v) in E do
    if comp.find(u) ≠ comp.find(v) then
        answer.append( (u,v) )
        comp.union(u, v)
    return answer
```

Union find operations:

- make_sets(A): one call with |A| = |V|
- find(a): 2m calls
- union(a,b): n-1 calls

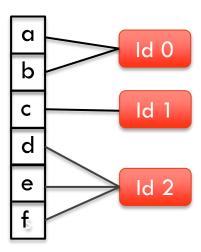
Simple union-find implementation

Sets are represented with lists. And we keep an array mapping elements to the set they belongs to

- make_sets(A) creates and initialized the array
- find(u) is a simple lookup in the array
- union(u,v) adds elements in u's set to v's set

Time complexity:

- make_sets(A) takes O(n) time, where n = |A|
- find(u) takes O(1) time
- union(u,v) take O(n) time



Kruskal's algorithm would run in $O(n^2)$ time after the edge weights are sorted.

Better union-find implementation

Keep track of cardinality of each set. When taking the union of two sets change the smallest.

That way the union of two sets A and B take O(min(|A|, |B|))

This way an element can change sets at most O(log n) time. So a sequence of n union operations takes at most O(n log n) time.

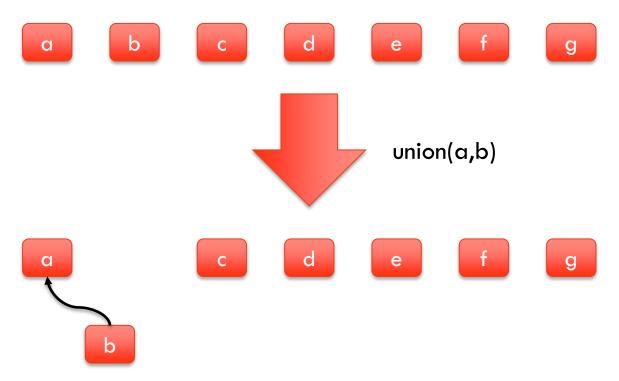
With this implementation, Kruskal's algorithm would run in O(m log n) time after the edge weights are sorted.

[Kruskal 1956]

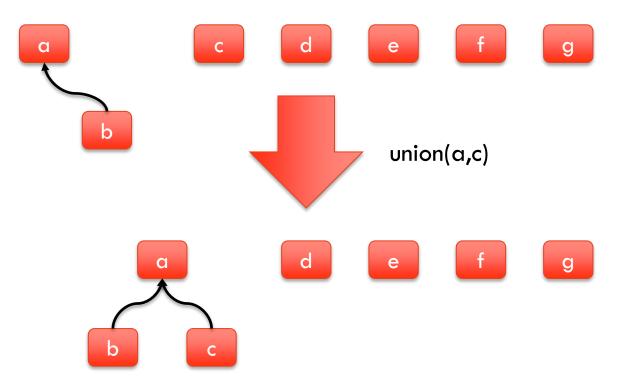
Keep track of sets using trees: For every node u in the graph we keep parent[u]



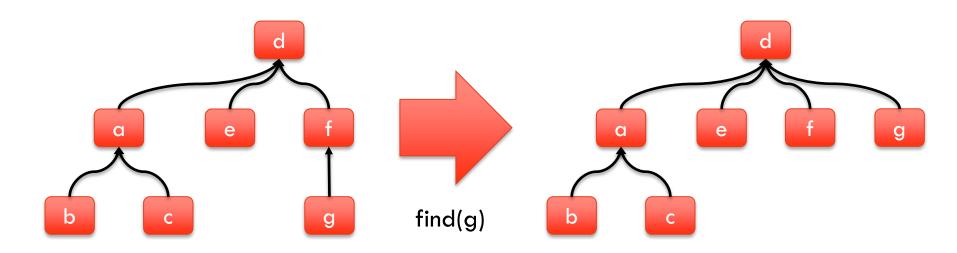
Taking the union of two roots \mathbf{u} and \mathbf{v} we set $parent[\mathbf{u}] = \mathbf{v}$ or $parent[\mathbf{v}] = \mathbf{u}$, depending on who has the largest tree.



Taking the union of two roots \mathbf{u} and \mathbf{v} we set $parent[\mathbf{u}] = \mathbf{v}$ or $parent[\mathbf{v}] = \mathbf{u}$, depending on who has the largest tree.



Perform find(u) by following parent[u] until reaching root r of tree. Set parent[v] to r for every node found along the way



It can be shown that a sequence of n union and m > n find operations takes at most $O(m \alpha(n))$ time, where $\alpha(n)$ is a **very slow** growing function called the *Inverse Ackerman function*.

How slow?

For $\alpha(n)$ to be larger than 4, you need n to be much larger than the number of subatomic particles in the universe.

So not a constant from a theoretical perspective, but for all practical (and impractical) purposes, you can treat it as 4.

With this implementation, Kruskal's algorithm would run in $O(m \alpha(n))$ time after the edge weights are sorted.

[Tarjan 1975]

Lexicographic Tiebreaking

To remove the assumption that all edge costs are distinct: perturb all edge costs by tiny amounts to break any ties.

Impact. Kruskal and Prim only interact with costs via pairwise comparisons. If perturbations are sufficiently small, MST with perturbed costs is MST with original costs.

For example, assuming all costs are integral, if we add i/n^2 to each edge e_i then any MST under the perturbed weights is still an MST under the original weights.

Implementation. Can handle arbitrarily small perturbations implicitly by breaking ties lexicographically, according to index.

Next week

General classes of algorithms:

- Brute force
- Greedy algorithms
 - Knapsack
 - Task scheduling
 - Text compression
- Divide-and-conquer
- Dynamic programming

