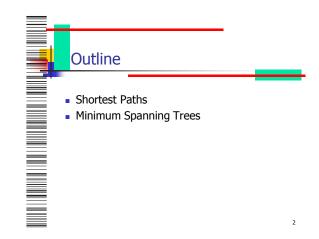
COMP9024: Data Structures and Algorithms

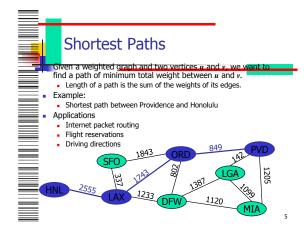
Week Twelve: Graphs (II)

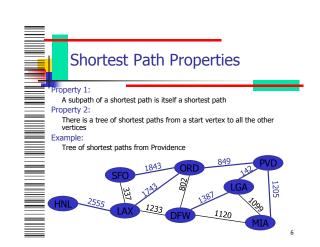
Hui Wu

Session 1, 2015 http://www.cse.unsw.edu.au/~cs9024



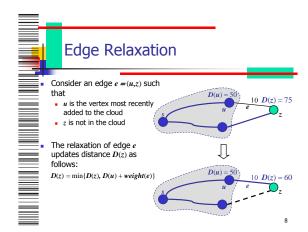
Weighted Graphs In a weighted graph, each edge has an associated numerical value, called the weight of the edge Edge weights may represent, distances, costs, etc. Example: In a flight route graph, the weight of an edge represents the distance in miles between the endpoint airports HNL 2555 LAX 1233 DFW 1120 MIA 4

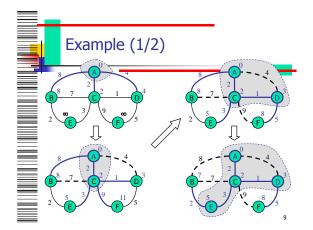


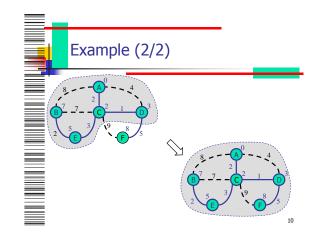


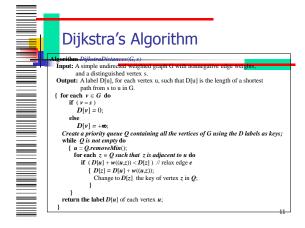
Dijkstra's Algorithm

- The distance of a vertex v from a vertex s is the length of a shortest path between s and v
- Dijkstra's algorithm computes the distances of all the vertices from a given start vertex s
- Assumptions:
 - the graph is connected
 - the edges are undirected
 - the edge weights are
- We grow a "**cloud**" of vertices, beginning with s and eventually covering all the vertices
- We store with each vertex ν a label $D(\nu)$ representing the distance of ν from s in the subgraph consisting of the cloud and its adjacent vertices
- At each step
 - We add to the cloud the vertex u outside the cloud with the smallest distance label, $\boldsymbol{D}(u)$
 - We update the labels of the vertices that are adjacent to *u* and not in the cloud







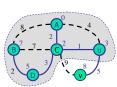


Analysis of Dijkstra's Algorithm Creating the priority queue Q takes $O(n \log n)$ time if using an adaptable priority queue, or O(n) time by using bottom-up heap

- construction.
- At each iteration of the **while** loop, we spend $O(\log n)$ time to remove vertex u from Q and $O(\operatorname{degree}(u)\log n)$ time to perform the relaxation procedure on the edges incident on u.
- The overall running time of the while loop is $O(\Sigma_u(1+\mathrm{degree}(u))\log n) = O((n+m)\log n)$ (Recall that $\Sigma_u \operatorname{degree}(u) = 2m$)
- The running time can also be expressed as $O(m \log n)$ since the graph is connected

Why Dijkstra's Algorithm Works

- Dijkstra's algorithm is based on the greedy method. It adds vertices by increasing distance.
- Suppose it didn't find all shortest distances. Let v be the first wrong vertex the algorithm processed.
- When the previous node, u, on the true shortest path was considered, its distance was correct.
- But the edge (u,v) was relaxed at that time!
- Thus, so long as D(v)≥D(u), v's distance cannot be wrong. That is, there is no wrong vertex.

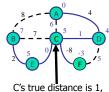


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Why It Doesn't Work for Negative-Weight Edges

 Dijkstra's algorithm is based on the greedy method. It adds vertices by increasing distance.

 If a node with a negative incident edge were to be added late to the cloud, it could mess up distances for vertices already in the cloud.



but it is already in the cloud with D(C)=5!

Bellman-Ford Algorithm (not in book)

- Works even with negativeweight edges
- Must assume directed edges (for otherwise we would have negativeweight cycles)
- Iteration i finds all shortest paths that use i edges.
- Running time: O(nm).

 Can be extended to detect a negative-weight cycle if it exists
 - How?

```
Algorithm BellmanFord(G, s) 

{ for each v \in G do 

  if (v = s) 

D[v] = 0; 

else 

D[v] = +\infty; 

for (i = 1; i \le n - 1; i + +) 

for each e \in G.edges() 

// relax \ edge \ e 

{ u = G.origin(e); 

z = G.opposite(u,e); 

r = D[u] + weight(e); 

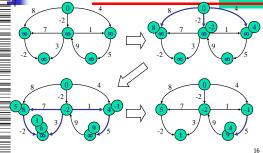
if (r < D[z]) 

D[z] = r; 

}
```

Bellman-Ford Example

Nodes are labeled with their D(v) values

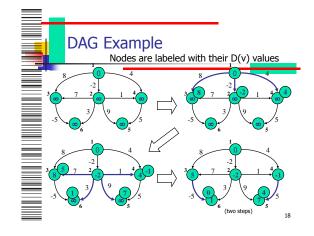


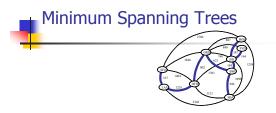
DAG-based Algorithm (not in book)

- Works even with negative-weight edges
- Uses topological orderDoesn't use any fancy
- data structures

 Is much faster than
 Dijkstra's algorithm
- Running time: O(n+m).

Algorithm DagDistances(G, s){ for all $v \in G.vertices()$ if (v = s) D[v] = 0;
 else
 $D[v] = +\infty$;
 $Perform\ a\ topological\ sort\ of\ the\ vertices;$ for $(u = 1;\ u \le n;\ u + +)$ # in topological order
 for each $e \in G.outEdges(u)$ do
 # relax edge e { z = G.opposite(u,e);
 # e = D[u] + weight(e);
 if (r < D[z]) # e = D[z] = r;
 }
}





Minimum Spanning Trees

Spanning subgraph

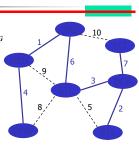
Subgraph of a graph G containing all the vertices of G

Spanning tree

Spanning subgraph that is itself a (free) tree

Minimum spanning tree (MST)

- Spanning tree of a weighted graph with minimum total edge weight
- Applications
 - Communications networks
 - Transportation networks



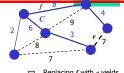
Cycle Property

Cycle Property:

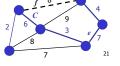
- Let *T* be a minimum spanning tree of a weighted graph *G*
- Let e be an edge of G that is not in T and C let be the cycle formed by e with T
- For every edge f of C, $weight(f) \le weight(e)$

Proof:

- By contradiction
- If weight(f) > weight(e) we can get a spanning tree of smaller weight by replacing e with f



Replacing f with e yields a better spanning tree



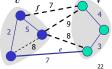
Partition Property

tition Property:

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

Proof: $\bullet \quad \text{Let } T \text{ be an MST of } G$

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



Replacing f with e yields

another MST

Kruskal's Algorithm

A priority queue stores the edges outside the cloud

- Key: weight
- Element: edge
- At the end of the algorithm
 - We are left with one cloud that encompasses the MST
 - A tree *T* which is our MST

Algorithm KruskalMST(G) for each vertex ν in G do define a $Cloud(\nu)$ of $\{\nu\}$; let Q be a priority queue; Insert all edges into Q using their weights as the key; T = Q; $T = \overline{O}$; while T has fewer than n-1 edges \mathbf{do} { edge e = T.removeMin(); Let u, v be the endpoints of e; if $(Cloud(v) \neq Cloud(u))$ { Add edge e to T; Merge Cloud(v) and Cloud(u); return T;

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Data Structure for Kruskal Algorithm

The algorithm maintains a forest of trees

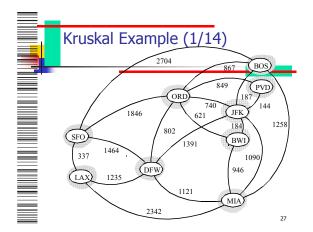
- An edge is accepted it if connects distinct trees
- We need a data structure that maintains a partition, i.e., a collection of disjoint sets, with the operations:
 - -find(u): return the set storing u
 - -union(u,v): replace the sets storing u and v with their union

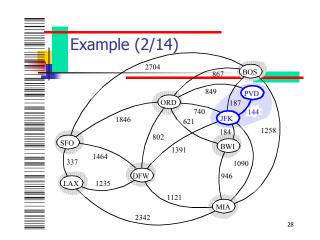


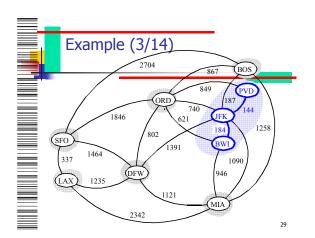
Representation of a Partition

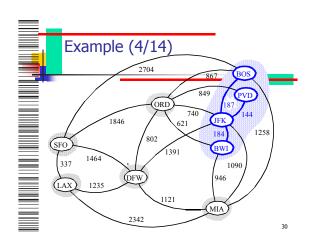
- Each set is stored in a sequence
- Each element has a reference back to the set
 - operation find(u) takes O(1) time, and returns the set of which u is a member.
 - in operation union(u,v), we move the elements of the smaller set to the sequence of the larger set and update their references
 - the time for operation union(u,v) is min(n_u,n_v), where n_u and n_v are the sizes of the sets storing u and v
- Whenever an element is processed, it goes into a set of size at least double, hence each element is processed at most log n times

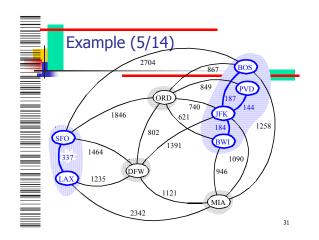
Partition-Based Implementation partition-based version of Kruskal's Algorithm performs cloud merges as unions and tests as finds Algorithm Kruskal(G): Input: A weighted graph G. Output: An MST T for G. { Let ${\it P}$ be a partition of the vertices of ${\it G}$, where each vertex forms a separate set; Let ${\bf Q}$ be a priority queue storing the edges of ${\bf G}$, sorted by their weights; Let T be an initially-empty tree; while Q is not empty do $\{ (u,v) = Q.removeMinElement(); \}$ if (P.find(u) != P.find(v)){ Add (u,v) to T: P.union(u,v): Running time: O((m+n) log n) return T;

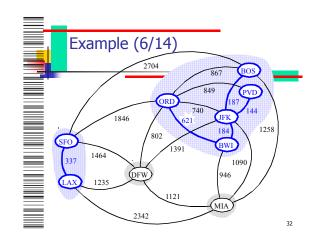


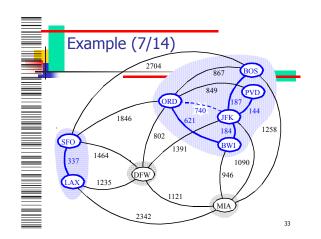


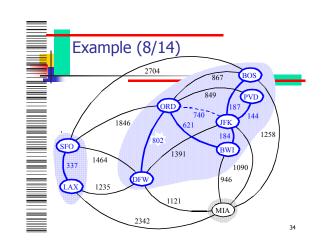


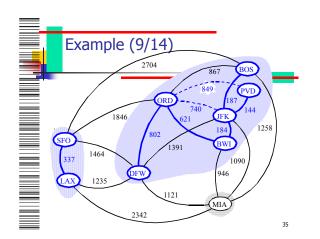


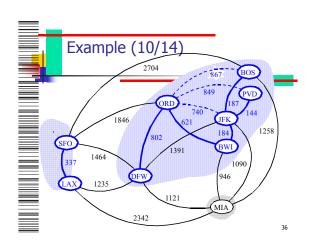


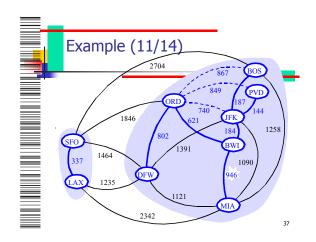


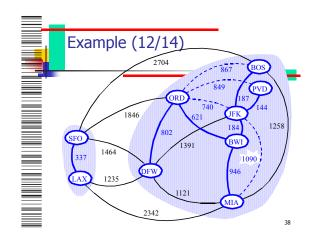


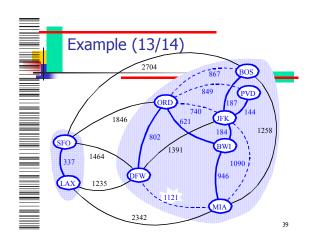


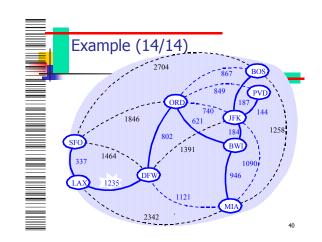








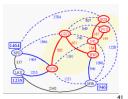


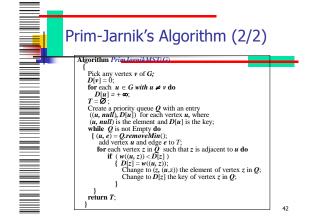


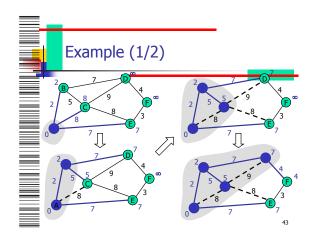
Prim-Jarnik's Algorithm (1/2)

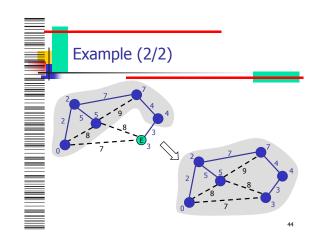
Similar to Dijkstra's algorithm (for a connected graph)

- We pick an arbitrary vertex s and we grow the MST as a cloud of vertices, starting from v
- We store with each vertex u a label D(u) = the smallest weight of an edge connecting u to a vertex in the cloud
- At each step:
 - We add to the cloud the vertex *u* outside the cloud with the smallest distance label
 - We update the labels of the vertices adjacent to \boldsymbol{u}









Analysis

- Graph operations
 - Method incidentEdges is called once for each vertex
- Label operations
 - We set/get the distance, parent and locator labels of vertex z $O(\deg(z))$ times
 - Setting/getting a label takes O(1) time
- Priority queue operations
 - Each vertex is inserted once into and removed once from the priority queue, where each insertion or removal takes $O(\log n)$ time
 - The key of a vertex w in the priority queue is modified at most $\deg(w)$ times, where each key change takes $O(\log n)$ time
- Prim-Jarnik's algorithm runs in $O((n+m)\log n)$ time provided the graph is represented by the adjacency list structure
 - Recall that $\sum_{v} \deg(v) = 2m$
- The running time is $O(m \log n)$ since the graph is connected

Baruvka's Algorithm (not in book)

 Like Kruskal's Algorithm, Baruvka's algorithm grows many "clouds" at once.

Algorithm BaruvkaMST(G) { T = V; // just the vertices of G while T has fewer than n-1 edges do for each connected component C in T do { Let edge e be the smallest-weight edge from C to another component in T; if e is not already in T then Add edge e to T; } } return T;

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

The running time is O(m log n).

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