



# CSE332: Data Abstractions

### Lecture 17: Shortest Paths

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# Single source shortest paths (from some specific point a)

- Done: BFS to find the minimum path length from v to u in O(|E|)
- Actually, can find the minimum path length from **v** to *every node* 
  - ► Still O(|E|)
  - No faster way for a "distinguished" destination in the worst-case
- Now: Weighted graphs

Given a weighted graph and node **v**, find the minimum-cost path from **v** to every node

- As before, asymptotically no harder than for one destination
- Unlike before, BFS will not work
- Aside: We can find the shortest path from every vertex to every other vertex in O(|V|<sup>3</sup>)



Why BFS won't work: Shortest path may not have the fewest edges

We will assume there are no negative weights

- Problem is ill-defined if there are negative-cost cycles
- What is the shortest path here?
- Even without negative cycles, can still get wrong answer if negative weights are involved

# Dijkstra's Algorithm (for shortest paths)

- Named after its creator Edsger Dijkstra (1930-2002)
  - Truly one of the "founders" of computer science; this is just one of his many contributions
  - Quotation: "computer science is no more about computers than astronomy is about telescopes"
- The idea: reminiscent of BFS, but adapted to handle weights
  - A priority queue will prove useful for efficiency (later)
  - Will grow the set of nodes whose shortest distance has been computed
  - Nodes not in the set will have a "best distance so far"



## Dijkstra's Algorithm: Idea



- Conceptually:
  - Grow our 'cloud' of known vertices by 1 each step
  - Pick a vertex outside the cloud, that's closest to our starting point
  - Guaranteed that we have the shortest path to everything within the cloud (more later)

### Dijkstra's Algorithm: Idea



- > Initially, start node has cost 0 and all other nodes have cost  $\infty$
- Mark each vertex as 'unknown' (also referred to as 'unvisited', 'unexplored')
- At each step:
  - Pick closest unknown vertex v (will be start node for first step)
  - Add it to the "cloud" of known vertices
  - Update distances for nodes with edges from v
- That's it! (Have to prove it produces correct answers)

# The Algorithm

- 1. For each node **v**, set **v**.cost=∞ and **v**.known=false
- 2. Set source.cost = 0
- 3. While there are unknown nodes in the graph
  - a) Select the unknown node  $\mathbf{v}$  with lowest cost
  - b) Mark **v** as known
  - c) For each edge (v, u) with weight w,

c1 = v.cost + w // cost of best path through v to u
c2 = u.cost // cost of best path to u previously known
if(c1 < c2) { // if the path through v is better
 u.cost = c1
 u.path = v // for computing actual paths
}</pre>



















- Once a vertex is marked 'known', the cost of the shortest path to that node is known
  - > As is the path itself

While a vertex is still not known, another shorter path to it might still be found

## Interpreting the results

Now that we're done, how do we get the path from, say, A to E?



# **Stopping Short**

How would this have worked differently if we were only interested in the path from A to G?





vertex	known?	cost	path
А		0	
В		??	
С		??	
D		??	
Ш		??	
F		??	
G		??	



vertex	known?	cost	path
А	Y	0	
В		??	
С		≤ <b>2</b>	А
D		≤ <b>1</b>	А
Ш		??	
F		??	
G		??	



vertex	known?	cost	path
А	Y	0	
В		≤6	D
С		≤ <b>2</b>	А
D	Y	1	А
ш		≤ <b>2</b>	D
F		≤7	D
G		≤ <b>6</b>	D



vertex	known?	cost	path
А	Y	0	
В		≤6	D
С	Y	2	А
D	Y	1	А
ш		≤ <b>2</b>	D
F		<b>≤ 4</b>	С
G		≤ <b>6</b>	D



vertex	known?	cost	path
А	Y	0	
В		≤ <b>3</b>	E
С	Y	2	А
D	Y	1	А
ш	Y	2	D
F		≤ <b>4</b>	С
G		≤6	D



vertex	known?	cost	path
А	Y	0	
В	Y	3	E
С	Y	2	А
D	Y	1	A
E	Y	2	D
F		≤ <b>4</b>	С
G		≤ <b>6</b>	D



vertex	known?	cost	path
А	Y	0	
В	Y	3	E
С	Y	2	А
D	Y	1	А
Ш	Y	2	D
F	Y	4	С
G		≤6	D



vertex	known?	cost	path
А	Y	0	
В	Y	3	E
С	Y	2	А
D	Y	1	А
E	Y	2	D
F	Y	4	С
G	Y	6	D

#### True or false:

If we were to count out all the edges 'found' by Dijkstra's, we would have |V|-1 edges



How will the best-cost-so-far from X to Y proceed?



#### How will the best-cost-so-far from X to Y proceed? 90, 81, 72, 63, 54

#### Where are we?

#### Have described Dijkstra's algorithm

- For single-source shortest paths in a weighted graph (directed or undirected) with no negative-weight edges
- An example of a greedy algorithm: at each step, irrevocably does the best thing it can at that step
  - Because of the way the algorithm is structured, the 'apparent best' actually is the best
- What should we do after learning an algorithm?
  - Prove it is correct
    - Not obvious!
    - We will sketch the key ideas
  - Analyze its efficiency

Rough intuition:

All the "known" vertices have the correct shortest path

- True initially: shortest path to start node has cost 0
- If it stays true every time we mark a node "known", then by induction this holds and eventually everything is "known"

Key fact we need: When we mark a vertex "known" we won't discover a shorter path later!

- This holds only because Dijkstra's algorithm picks the node with the next shortest path-so-far
- The proof is by contradiction...

# Correctness: The Cloud (Rough Idea)



Suppose v is the next node to be marked known ("added to the cloud")

- The best-known path to v must have only nodes "in the cloud"
  - Since we've selected it, and we only know about paths through the cloud to a node right outside
- Assume the actual shortest path to **v** is different than the best-known
  - It won't use only cloud nodes, or we would know about it; so it must use noncloud nodes
  - Let w be the *first* non-cloud node on this 'actual shortest path'
  - The part of the path up to w is already known and must be shorter than the
  - best-known path to **v**. So **v** would not have been picked. Contradiction.

```
Efficiency, first approach
 Use pseudocode to determine asymptotic run-time
   Notice each edge is processed only once
dijkstra(Graph G, Node start) {
  for each node: x.cost=infinity, x.known=false
  start.cost = 0
  while(not all nodes are known) {
                                                     O(|V|^2)
    b = find unknown node with smallest cost
    b.known = true
    for each edge (b,a) in G
     if(!a.known)
       if(b.cost + weight((b,a)) < a.cost) 
                                                     O(|E|
         a.cost = b.cost + weight((b,a))
         a.path = b
```

# Improving asymptotic running time

- So far: O(|∨|²)
- We had a similar "problem" with topological sort being O(|V|<sup>2</sup>) due to each iteration looking for the node to process next
  - We solved it with a queue of zero-degree nodes
  - But here we need the lowest-cost node and costs can change as we process edges

#### Solution?

- A priority queue holding all unknown nodes, sorted by cost
- But must support decreaseKey operation
  - Must maintain a reference from each node to its position in the priority queue

## Efficiency, second approach

Use pseudocode to determine asymptotic run-time

```
dijkstra(Graph G, Node start) {
  for each node: x.cost=infinity, x.known=false
                                                       O(|V|)
  start.cost = 0
  build-heap with all nodes
  while (heap is not empty) {
                                                   O(|V|log|V|
    b = deleteMin()
    b.known = true
    for each edge (b,a) in G
     if(!a.known)
      if(b.cost + weight((b,a)) < a.cost){</pre>
                                                   O(|E|\log|V|)
        decreaseKey(a, "new cost - old cost"
         a.path = b
      }
                                           O(|V|\log|V|+|E|\log|V|)
```

#### Dense vs. sparse again

- First approach:  $O(|V|^2)$
- Second approach: O(|V|log|V|+|E|log|V|)
- So which is better?
  - Sparse: O(|V|log|V|+|E|log|V|)
    - ▶ (if |E| > |V|, then *O*(|E|log|V|))
  - Dense:  $O(|V|^2)$
- But, remember these are worst-case and asymptotic
  - Priority queue might have slightly worse constant factors
  - On the other hand, for "normal graphs", we might call decreaseKey rarely (or not percolate far), making |E|log|V| more like |E|