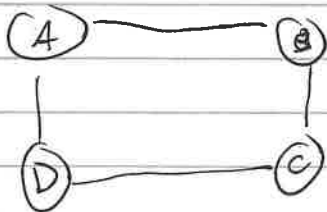


BFS and Dijkstra's Algorithm.

DFS is one way for us to search a graph. By using the stack, we can quickly explore deeply into a graph. It is not particularly good at finding short paths. Here, we mean on an unweighted graph by the number of edges.



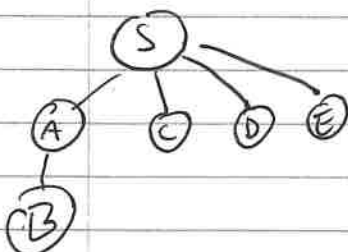
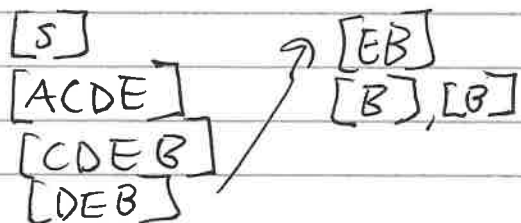
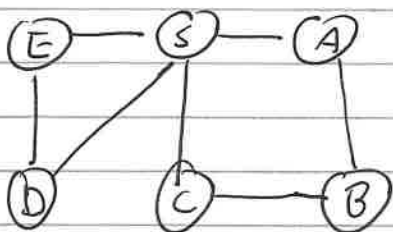
Although D is only distance 1 from A, DFS takes the route A-B-C-D giving a path of length 3.

We need a graph traversal algorithm which prioritizes the closest nodes first. This is called Breadth-First-Search, or BFS.

```

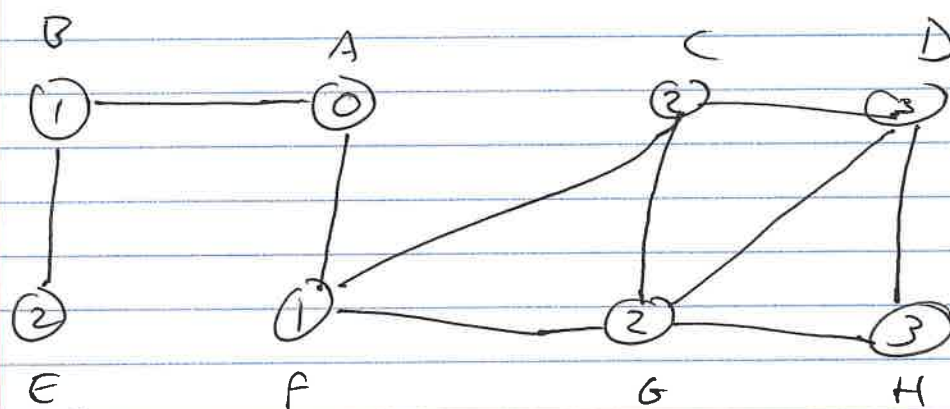
def bfs(G, s)
  for each  $u \in V$ :  $dist(u) = \infty$ 
   $dist(s) = 0$ 
   $Q = [s]$ 
  while  $Q \neq \emptyset$ :
     $u = eject(Q)$ 
    for each  $(u, v) \in E$ :
      if  $dist(v) = \infty$ :
        inject  $v$  into  $Q$ 
         $dist(v) = dist(u) + 1$ 
  
```

Each vertex is pushed and popped from the queue exactly once, which takes $O(1)$ time, so $O(V)$. The adjacency lists are looped over once for each, their sum of lengths being $O(E)$, giving us a total runtime of $O(V + E)$. Takes the same time as DFS but just in a different order.



In the spirit of computing shortest paths, we are only concerned with nodes reachable from our start. The queue structure no more than two levels at a time haven't been fully visited. You cannot see grandchildren before seeing children.

2



E

[A]

[BF]

[FE]

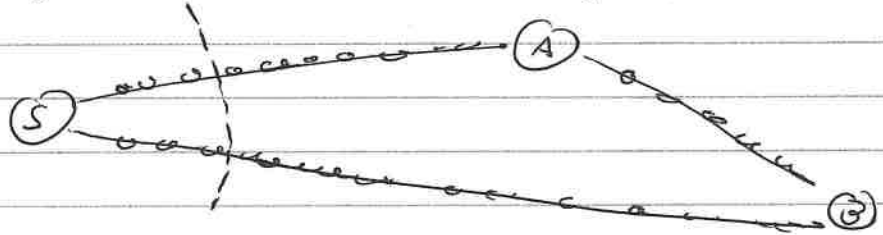
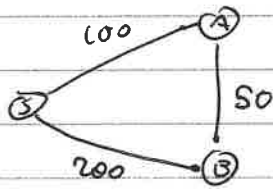
[ECG]

[CG]

[GD]

[DH]

Consider a weighted graph now, with positive weights. we can compute shortest paths again by transforming it into an unweighted graph



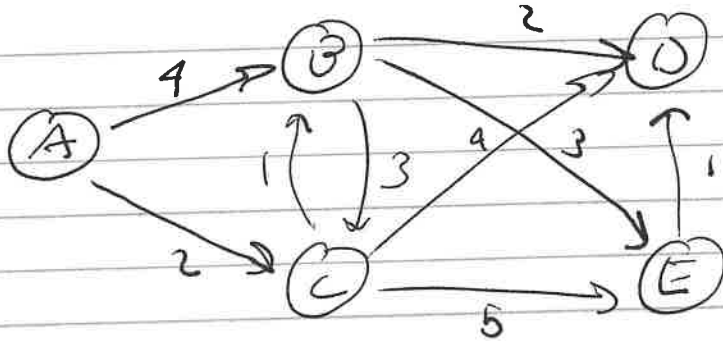
the problem now is it takes 100 steps to see that A is closer to S than B after one edge. why not just directly compare 100 with 200 then choose to BFS on A next? This is the heart of Dijkstra's algorithm.

Instead of a simple queue, we use a priority queue with the following operations:

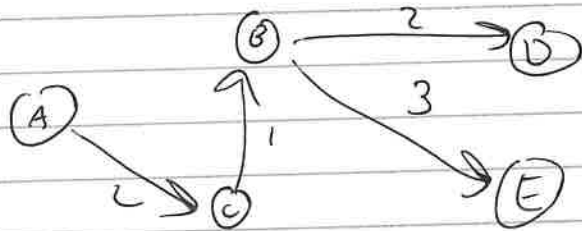
insert	into queue
decrease key	decreases value of some key
delete min	returns element w/ smallest key
make-queue	builds the queue quickly.

```

def dijkstra(G, l, s):
    for each  $u \in V$ 
         $\text{dist}(u) = \infty$ 
         $\text{prev}(u) = \text{null}$ 
     $\text{dist}(s) = 0$ 
     $H = \text{makequeue}(V)$  with  $\text{dist}$  as keys
    while  $H \neq \emptyset$ :
         $u = \text{deletemin}(H)$ 
        for each  $(u, v) \in E$ :
            if  $\text{dist}(v) > \text{dist}(u) + l(u, v)$ 
                 $\text{dist}(v) = \text{dist}(u) + l(u, v)$ 
                 $\text{prev}(v) = u$ 
                decrease key( $H, v$ ).
    
```



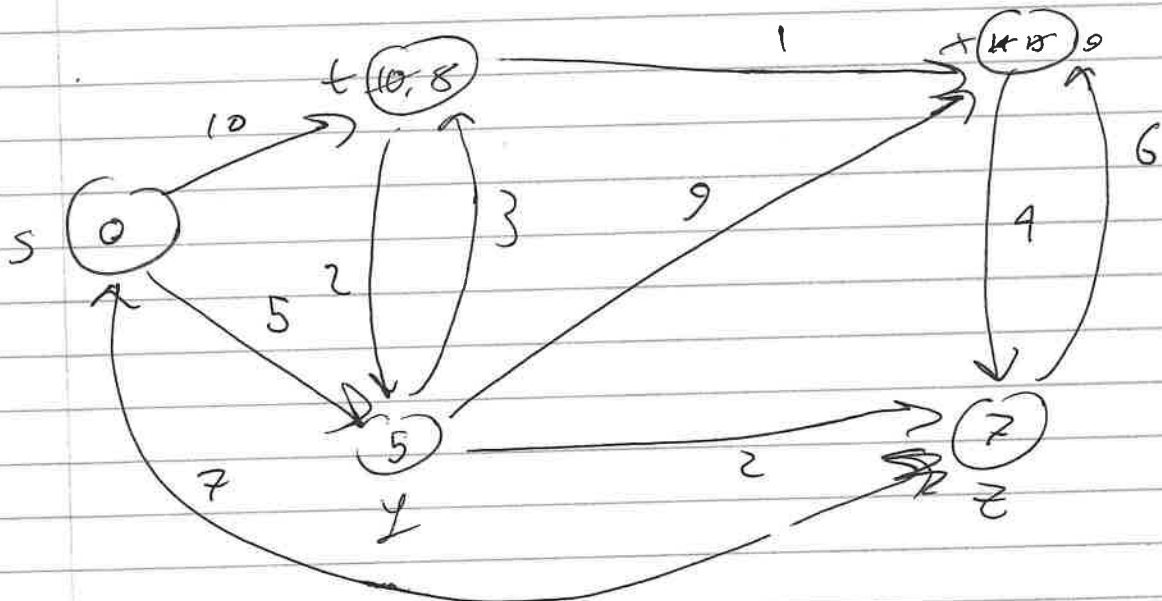
	A	C	B
A:	0		
B:	∞	4	3
C:	∞	2	
D:	∞		6
E:	∞		7



Dijkstra's is just BFS, but it runs the depends on the PQ implementation. There are $|V|$ delete min operations and $|V|+|E|$ insertions/decrease keys.

	delete min	insert	Dijkstra
Array	$O(V)$	$O(1)$	$O(V^2)$
binary heap	$O(\log V)$	$O(\log V)$	$O((V+E) \log V)$
d-ary heap	$(d \log V / \log d)$	$O(\log V / \log d)$	$O((Vd+E) \frac{\log V}{\log d})$

If G is dense, $E \propto V^2$, then array is best. binary heap becomes better and $E \approx V^2 / \log V$



"The shortest path s-t is less than the shortest path s-u-t & u."