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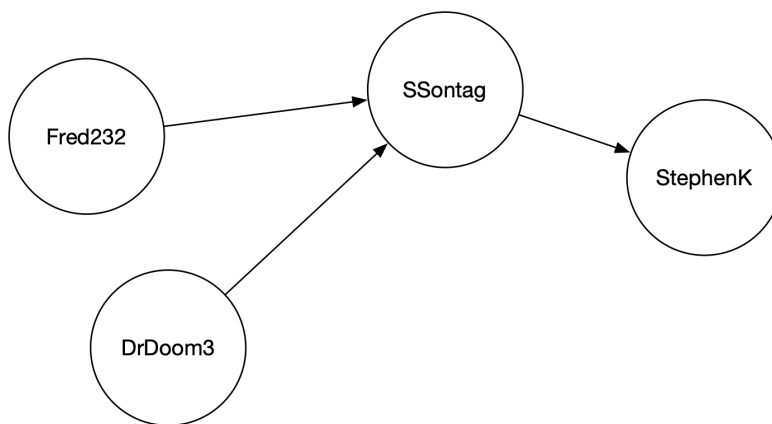
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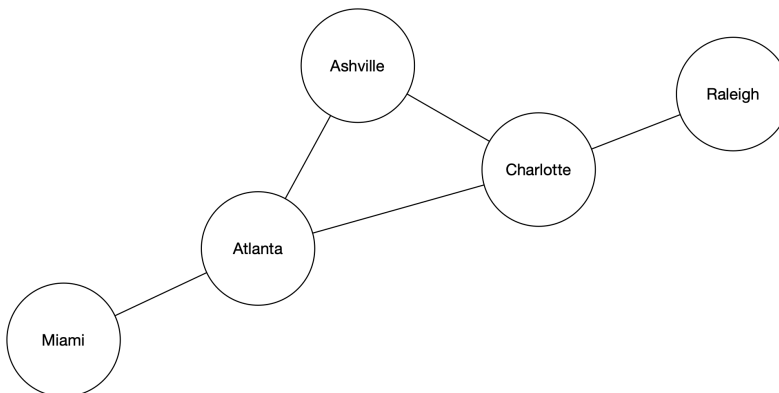
# Introduction to Graphs

Some data is easiest to work with if we imagine it as a set of *nodes* connected by *edges*. For example, on some social networks each user can follow any number of other users. We can think of each user as node and the edge points from the user who follows to the user they follow:



This diagram shows four users and three follows. Following is a directed relationship: Fred232 follows SSontag, but SSontag doesn't follow Fred232. So we would say that this is a *directed graph* with four nodes and three edges.

There are also undirected graphs. for example, you can imagine a graph that represents big data lines between cities. All the big data lines allow communications in both directions:



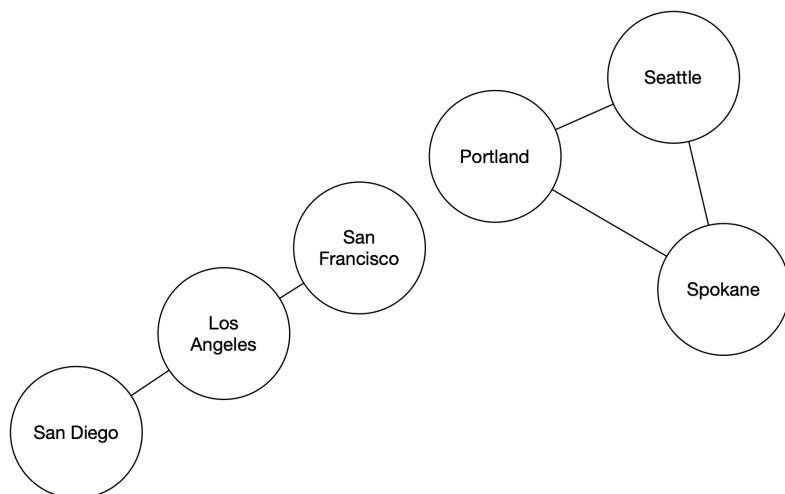
The arrows are gone: if data can flow from Charlotte to Raleigh, then data can flow from Raleigh to Charlotte.

There is a whole branch of mathematics called *Graph Theory* that studies the properties of graphs. Here are two questions that we might ask about this graph:

- What is the shortest number of edges that we would need to follow to get from Miami to Raleigh?
- Does the graph have any paths where you could end up where you started? This is called a *cycle*. This graph has one cycle: Atlanta  $\rightarrow$  Asheville  $\rightarrow$  Charlotte  $\rightarrow$  Atlanta.

There are even database systems that are specifically designed to hold and analyze graph data. Not surprisingly, these are called *Graph Databases*.

Some graphs are *connected*: you can get from one node to any other node by following edges. Is this graph connected?



This graph is *not* connected! You can't follow edges from San Diego to Seattle.

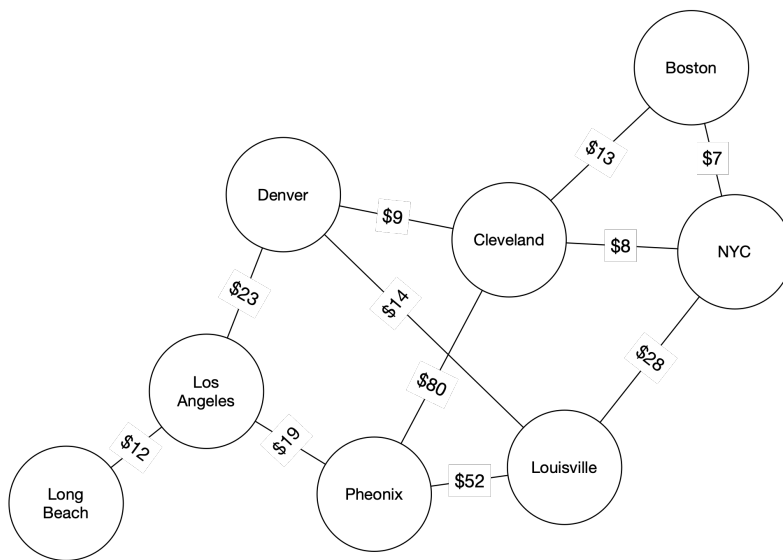
In graph data, the nodes and edges often have attributes. For example, a node representing a city might have a name and a population. An edge representing a data line might have a bandwidth (bits per second) and a latency (how many nanoseconds between when you put a bit into the pipe and when it comes out the other end.).

## 1.1 Finding Good Paths

For a lot of problems, we are trying to find the best path from one node to another. If all the edges are the same, this usually means finding the path that requires walking the

fewest edges.

Sometimes the edges have a cost attribute. For example, you might want to find the cheapest way to ship a container from New York City to Long Beach, Calif. In this case the nodes are train depots. Each train line between the depots has a cost. What is the cheapest path?



When edges have costs like this, we call the *weighted edges*.

The graphs that you see here are really small, so finding efficient paths isn't difficult. – you could just try all of them! However, in many computer programs, we are working with millions of nodes and edges. Efficient graph algorithms are *really* important.

## 1.2 Graphs in Python

In this section you are going to write Python classes that will let you represent an undirected graph with weighted edges, like the shipping problem above.

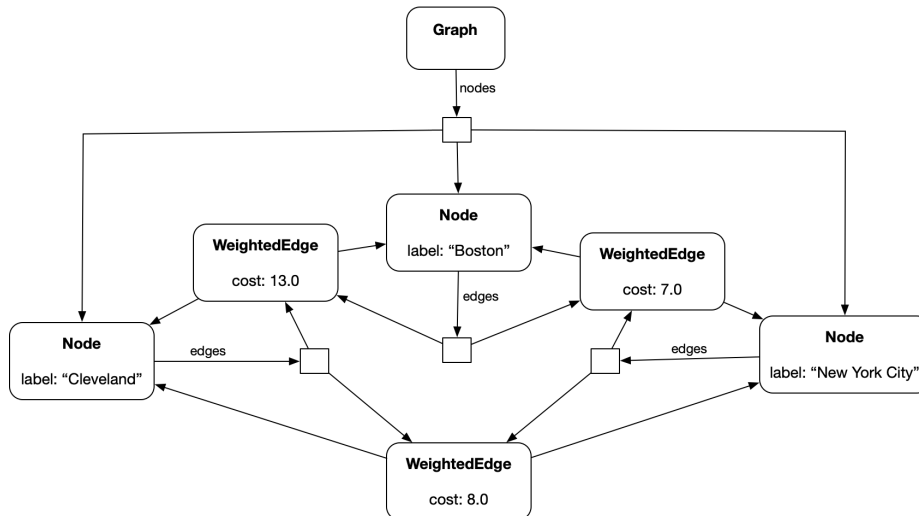
(Naturally things would look a little different if the graph were directed or the edges were unweighted, but this is a good starting place.)

Create a file called `graph.py`. This will hold the code for your `Node` and `WeightedEdge` classes. We will also create a `Graph` class that will just hold onto the list of its nodes.

- A `Node` will have a label string and a list of edges that touch it.
- A `Edge` will have a cost and two nodes: `node_a` and `node_b`.

- A Graph will have a list of nodes.

Here what the object diagram would look like if you had only three cities:



Put this code into graph.py

```
class Node:
    def __init__(self, label):
        self.label = label
        self.edges = []

    def __repr__(self):
        return f"(node:{self.label}, edges:{len(self.edges)})"

class WeightedEdge:
    def __init__(self, cost, node_a, node_b):
        self.cost = cost
        self.node_a = node_a
        node_a.edges.append(self)
        self.node_b = node_b
        node_b.edges.append(self)

    def other_end(self, node_from):
        if self.node_a == node_from:
            return self.node_b
        else:
            return self.node_a

class Graph:
    def __init__(self):
        self.nodes = []

    def add_node(self, new_node):
```

```
self.nodes.append(new_node)

def __repr__(self):
    return f"(Graph:{self.nodes})"
```

Now let's create some instances of `Node` and `WeightedEdge` and wire them together. Create another file in the same directory called `cities.py`. Put in this code:

```
import graph

# Create an empty graph
network = graph.Graph()

# Create city nodes and add to graph
long_beach = graph.Node("Long Beach")
network.add_node(long_beach)
los_angeles = graph.Node("Los Angeles")
network.add_node(los_angeles)
denver = graph.Node("Denver")
network.add_node(denver)
pheonix = graph.Node("Pheonix")
network.add_node(pheonix)
louisville = graph.Node("Louisville")
network.add_node(louisville)
cleveland = graph.Node("Cleveland")
network.add_node(cleveland)
boston = graph.Node("Boston")
network.add_node(boston)
nyc = graph.Node("New York City")
network.add_node(nyc)

# Create edges
graph.WeightedEdge(12, long_beach, los_angeles)
graph.WeightedEdge(23.0, los_angeles, denver)
graph.WeightedEdge(19, los_angeles, pheonix)
graph.WeightedEdge(52, pheonix, louisville)
graph.WeightedEdge(14, denver, louisville)
graph.WeightedEdge(80, pheonix, cleveland)
graph.WeightedEdge(9, denver, cleveland)
graph.WeightedEdge(8, cleveland, nyc)
graph.WeightedEdge(28, louisville, nyc)
graph.WeightedEdge(7, nyc, boston)
graph.WeightedEdge(13, cleveland, boston)

print(network)
```

Run it:

```
python3 cities.py
```

You should see some rather unexciting output:

```
(Graph:[(node:Long Beach, edges:1), (node:Los Angeles, edges:3), (node:Denver, edges:3),  
(node:Pheonix, edges:3), (node:Louisville, edges:3), (node:Cleveland, edges:4),  
(node:Boston, edges:2), (node:New York City, edges:3)])
```

But we will make it more exciting in the next chapter!



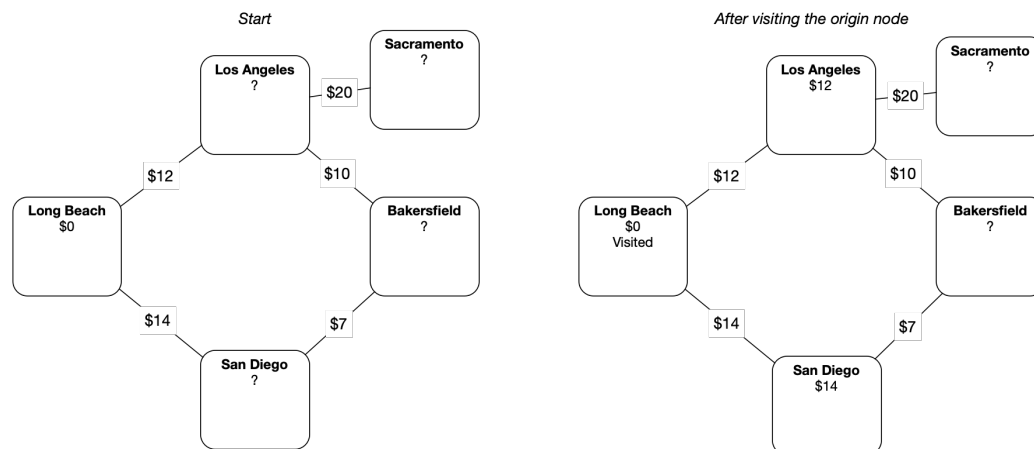
# Dijkstra's Algorithm

Edsger W. Dijkstra was a great Dutch computer scientist. He came up with an algorithm for finding the cheapest path through a graph with weighted edges. Today it is known as *Dijkstra's Algorithm*. It is used in a wide variety of common problems. It is also really pretty simple and elegant.

## 2.1 Algorithm Description

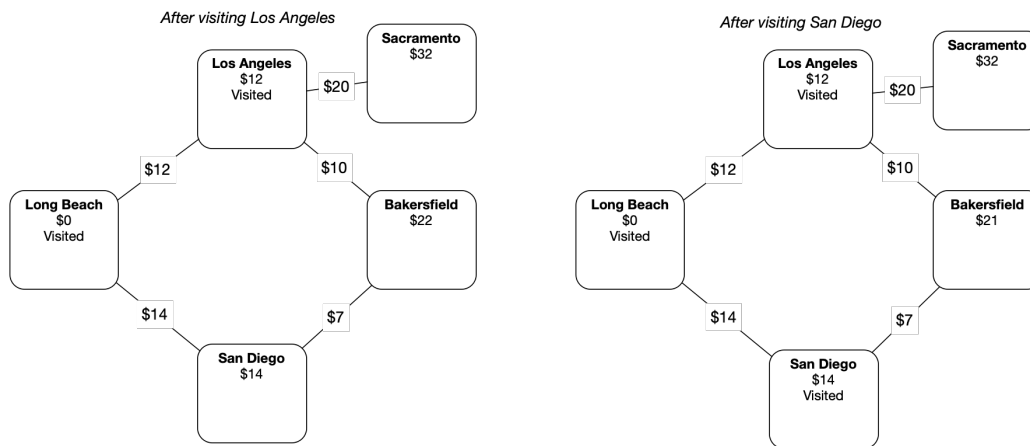
You are going to mark each node with how much it would cost to get there from some origin node. For example, if you are shipping a container from Long Beach, you will mark each city with the cost of getting the container to that city.

You start by marking the price for Long Beach to zero. (The container is already there.) Then, you mark each adjacent city with the cost on the edge. Now you declare Long Beach to be “visited”.



Now, you find the cheapest of the unvisited nodes. In this case, Los Angeles is cheaper than San Diego, so that is the node you will visit next.

You mark all of the unvisited nodes adjacent to Los Angeles, with the price to ship it to Los Angeles plus the cost of shipping the container from Los Angeles to that city. Note that Bakersfield is marked with \$22.



Now the cheapest unvisited node is San Diego. So you mark its neighbors with the cost to ship to San Diego plus the price to ship from San Diego to the neighbor.

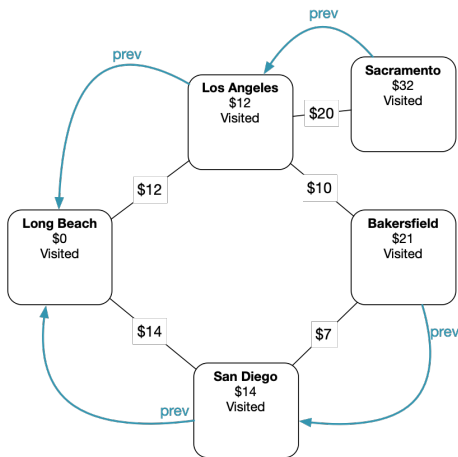
Notice that Bakersfield is already labeled with \$22 from a route through Los Angeles. But the price would be \$21 if you shipped it to Bakersfield via San Diego. Because the new route is cheaper, you change the price to the lower value.

(What does it mean that a node is “visited”? If a node is marked visited, it is marked with a price that won’t get any smaller.)

And you continue visiting the cheapest unvisited node until all the nodes have been visited. Then you know every node has been marked with its lowest price.

In a big graph, each node may be marked several times in this process – each time with a lower price from a cheaper router.

Of course, once you have the price, you will ask “What is the route that gets me that price?” So we will also mark each node with the neighbor from which it would receive the shipment – the previous node. This is easy to do as we execute the algorithm.



Now, to figure out the cheapest route from San Diego to Bakersfield, we start at the destination and follow the `prev` pointer back through San Diego and then to Long Beach.

## 2.2 Implementation

We don't actually want to sully our graph objects with the three additional pieces of information we need:

- The current minimal cost from the origin node. This is usually called the `dist`, from "distance".
- The neighbor who gives us the current minimal cost. This is usually called `prev`, from "previous".
- Whether the city is visited or now.

So we will keep them in collections external to the graph.

For example, to keep track of the `dist`, we will have a `dist` dictionary: Each node will be a key, the current minimal cost will be the value. If the node hasn't received even a first cost, we will put in infinity as the cost.

(After the algorithm is run, if the cost of a node is still infinity, that means that it cannot be reached from the origin node.)

We will also have a `prev` dictionary. The final node will be the key, and its previous neighbor will be the value.

Finally, the graph has a list of all the nodes, so we can just keep a set of the unvisited nodes.

Add a method to the Graph class that implements Dijkstra's algorithm:

```
def cost_from_node(self, origin_node):
    # Cost of cheapest path from origin node discovered so far
    # Initially the origin is zero and all the other are infinity
    dist = {k: math.inf for k in self.nodes}
    dist[origin_node] = 0.0

    # The previous city on that cheapest path
    prev = {}

    # All the nodes start as unvisited
    unvisited = set(self.nodes)

    # While there are still unvisited nodes
    while unvisited:

        # Find unvisited node with lowest cost
        min_cost = math.inf
        for u in unvisited:
            if dist[u] < min_cost:
                current_node = u
                min_cost = dist[u]

        # If none are less than inf, we are done
        # This happens in graphs that are not connected
        if min_cost == math.inf:
            return (dist, prev)

        # Remove the lowest cost node from the unvisited list
        unvisited.remove(current_node)

        # Update all the unvisited neighbors
        for edge in current_node.edges:

            # What node is at the other end of this edge?
            v = edge.other_end(current_node)

            # Visited nodes are already minimized, skip them
            if v not in unvisited:
                continue

            # Is this a shorter route?
            alt = dist[current_node] + edge.cost
            if alt < dist[v]:

                # Update the distance and prev dicts
                dist[v] = alt
                prev[v] = current_node

    return (dist, prev)
```

Append some code to your `cities.py` that test this method:

```
(cost_from_long_beach, prev) = network.cost_from_node(long_beach)
print(f"\nMinimum costs from Long Beach = {cost_from_long_beach}")
print(f"\nLast city before = {prev}")

nyc_cost = cost_from_long_beach[nyc]

if nyc_cost < math.inf:
    print(f"\n*** Total cost from Long Beach to NYC: ${nyc_cost:.2f} ***")
else:
    print("You can't get to NYC from Long Beach")
```

When you run it, you should get a list of how much it costs to ship a container to each city from Long Beach:

```
Minimum costs from Long Beach = {(node:Long Beach, edges:1): 0.0,
(node:Los Angeles, edges:3): 12.0, (node:Denver, edges:3): 35.0,
(node:Pheonix, edges:3): 31.0, (node:Louisville, edges:3): 49.0,
(node:Cleveland, edges:4): 44.0, (node:Boston, edges:2): 57.0,
(node:New York City, edges:3): 52.0}
```

You will also get a collection of node pairs. What are these? For each node, you get the node that you would pass through on the cheapest route from Long Beach:

```
Last city before = {(node:Los Angeles, edges:3):(node:Long Beach, edges:1),
(node:Denver, edges:3):(node:Los Angeles, edges:3),
(node:Pheonix, edges:3):(node:Los Angeles, edges:3),
(node:Louisville, edges:3):(node:Denver, edges:3),
(node:Cleveland, edges:4):(node:Denver, edges:3),
(node:New York City, edges:3):(node:Cleveland, edges:4),
(node:Boston, edges:2): (node:Cleveland, edges:4)}
```

Your users won't want to read this; Give them the shortest path as a list. Add a function to `graph.py` that turns the `prev` table into a path of nodes that lead from the origin to the destination:

```
def shortest_path(prev, destination):

    # Include the destination in the path
    path = [destination]
    current_node = destination

    # Keep stepping backward in the path
    while current_node in prev:

        # What node should come before the current node?
        previous_node = prev[current_node]
```

```
# Insert it at the start of the list
path.insert(0, previous_node)
current_node = previous_node

return path
```

Test that out:

```
if nyc_cost < math.inf:
    print(f"*** Total cost from Long Beach to NYC: ${nyc_cost:.2f} ***")

    path_to_nyc = graph.shortest_path(prev, nyc)
    print(f"*** Cheapest path from Long Beach to NYC: path_to_nyc ***")
else:
    print("You can't get to NYC from Long Beach")
```

This should look like this:

```
*** Cheapest path from Long Beach to NYC: [(node:Long Beach, edges:1),
(node:Los Angeles, edges:3), (node:Denver, edges:3), (node:Cleveland, edges:4),
(node:New York City, edges:3)] ***
```

## 2.3 Making it faster

On really big networks, doing a full Dijkstra's algorithm would take too long. So there are a lot of methods for getting similar results quickly. When you ask for directions from Google Maps, it doesn't do a full Dijkstra's Algorithm for every possible route – it would just take too long.

But there is a way to speed up this implementation. Look at this snippet:

```
# Find unvisited node with lowest cost
min_cost = math.inf
for u in unvisited:
    if dist[u] < min_cost:
        current_node = u
        min_cost = dist[u]
```

We are scanning through the list of all unvisited nodes, one-by-one, looking for the one with the lowest cost. If we kept this list sorted by cost, then the next one to visit would always be the first one in the list. This is done with a *priority queue* – a list that keeps itself sorted by some priority number – in this case the cost. In python, the standard priority queue is `heapq`.

(So why didn't I implement this using `heapq`? For Dijkstra's Algorithm, the nodes' priority – the current cost – changes as we find cheaper routes. `heapq` doesn't handle the changing priority very gracefully.)

In the next chapter, you will make a priority queue class that will work in this case.





# Binary Search

As mentioned in the last chapter, you are going to make a priority queue for use with Dijkstra's Algorithm. Using it will look like this:

```
import kqueue

myqueue = pqueue.PriorityQueue()
myqueue.add(long_beach, 0) # Inserts first city and its cost
myqueue.add(san_diego, 14) # Puts San Diego after Long Beach
myqueue.add(los_angeles, 12) # Inserts LA between Long Beach and San Diego
current_city = myqueue.pop() # Returns first city (Long Beach) and removes it
```

Now if an item gets a new priority, we need to remove it and reinsert it in the new spot.

```
myqueue.add(city_a, 16) # Puts it last in the queue
myqueue.update(city_a, 16, 13) # Moves it to between LA and San Diego
```

### 3.1 A Naive Implementation of the Priority Queue

Create a file called `kqueue.py`. Let's do a simple implementation that stores the priority and the data as tuple. And we will keep it sorted by the priority. If two tuples have the same priority, we'll sort by the data.

Type this in to `kqueue.py`:

```
class PriorityQueue:
    def __init__(self):
        self.list = []

    # Return and remove the first item
    def pop(self):
        if len(self.list) > 0:
            return self.list.pop(0)
        else:
            return None

    def __len__(self):
        return len(self.list)

    def update(self, value, old_priority, new_priority):
```

```
        old_pair = (old_priority, value)
        self.list.remove(old_pair)
        self.add(value, new_priority)

    def add(self, value, priority):
        pair = (priority, value)
        # Add it at the end
        self.list.append(pair)
        # Resort the list
        self.list.sort()
```

This will work fine, but it could be much more efficient:

- Every time we add a single element, we resort the whole list.
- The function `remove` is searching the list sequentially for the item to delete.

In a minute, we will revisit these inefficiencies and make the better.

## 3.2 Using the Priority Queue

We are going to change `graph.py` to use the priority queue. While we are doing, why don't we also shrink the memory footprint of our program a bit.

Notice that as the algorithm is running, each node is in one of three states:

- Unseen: In the earlier implementation, these were the nodes with `math.inf` as their cost.
- Seen, but not finalized: These are “unvisited” but don't have `math.inf` as their cost.
- Finalized: These are the “visited” nodes – we know that their cost won't decrease any more.

We can shrink the memory footprint by not putting the unseen into the `dist` dictionary at all. And instead of a separate set for “unvisited” what if we moved finalized nodes and their distances into a separate dictionary?

Rewrite the `cost_from_node` function in `graph.py`:

```
# Visited nodes are already minimized, skip them
if v in finalized_dist:
    continue

# What is the cost to this neighbor?
alt = current_node_cost + edge.cost
```

```

# Is this the first time I am seeing the node?
if v not in seen_dist:

    # Insert into the seen_dict, prev, and priority queue
    seen_dist[v] = alt
    prev[v] = current_node
    pqueue.add(v, alt)

else: # v has been seen. Is this a cheaper route?
    old_dist = seen_dist[v]
    if alt < old_dist:
        # Update the seen_dict, prev, and priority queue
        seen_dist[v] = alt
        prev[v] = current_node
        pqueue.update(v, old_dist, alt)

return (finalized_dist, prev)

```

This should be have exactly the same except for the unreachable nodes. If you have a graph that is not connected, there will be nodes that can't be reached from the origin. In the old version, these had a cost of `math.inf`. Now they just won't be in the dictionary at all. So, change `cities.py` to deal with this:

```

if nyc in cost_from_long_beach:
    nyc_cost = cost_from_long_beach[nyc]
    print(f"\n*** Total cost from Long Beach to NYC: ${nyc_cost:.2f} ***")

    path_to_nyc = graph.shortest_path(prev, nyc)
    print(f"\n*** Cheapest path from Long Beach to NYC: {path_to_nyc} ***")
else:
    print("You can't get to NYC from Long Beach")

```

If you run `cities.py` now, it should behave exactly like the old version.

But there is a bug. It will rear its head if two cities with the same cost are in the priority queue together. Change `cities.py` so that Denver and Pheonix have the same cost:

```

graph.WeightedEdge(12, long_beach, los_angeles)
graph.WeightedEdge(19, los_angeles, denver)
graph.WeightedEdge(19, los_angeles, pheonix)

```

Now try running it. You should get an error:

```

TypeError: '<' not supported between instances of 'Node' and 'Node'

```

What happened? The `loc_for_pair` method is comparing tuples made up of a `float` and a `Node`. The `float` comes first in the tuple, so that is compared first. However, if the two tuples have the same priority, it then compares nodes.

The error statement say “Nodes don’t have a less-than method; I don’t know how to compare them.”

Each `Node` lives at an address in memory. You can get that address as a number using the `id` function. The ID is unique and constant over the life of the object. It is a rather arbitrary ordering, but it will work for this problem. Add a method to your `Node` class:

```
# Nodes will be ordered by their location in memory
def __lt__(self, other):
    return id(self) < id(other)
```

Fixed.

Now let’s make the priority queue more efficient.

### 3.3 Binary Search

The phone company in every town used to print a thing called a phone book. The names and phone numbers were arranged alphabetically. As you might imagine, these books often had more than a thousand pages.

If you were looking for “John Jeffers”, you wouldn’t start at the first page and read sequentially until you reached his name. You would open the book in the middle, and see a name like “Mac Miller”, and then think “Jeffers comes before Miller”. Then you would split the pages in your left hand in half and see a name like “Hester Hamburg” and think “Jeffers comes after Hamburg”. Then you would split the pages in your right hand, and so on until you found the page with “John Jeffers” on it.

That is binary search.

Binary Search is a search algorithm that finds the position of a target value within a sorted array. The binary search algorithm works by repeatedly dividing the search interval in half. If the target value is equal to the middle element of the array, the position is returned. If the target value is less or greater than the middle element, the search continues in the lower or upper half of the array respectively.

### 3.4 Algorithm

The binary search algorithm can be described as follows:

1. If the array is empty, the search is unsuccessful, so return "Not Found".
2. Otherwise, compare the target value to the middle element of the array.
3. If the target value matches the middle element, return the middle index.
4. If the target value is less than the middle element, repeat the search with the lower half of the array.
5. If the target value is greater than the middle element, repeat the search with the upper half of the array.
6. Repeat steps 2-5 until the target value is found or the array is exhausted.

```
class PriorityQueue:
    def __init__(self):
        self.list = []

    # Return and remove the first item
    def pop(self):
        if len(self.list) > 0:
            return self.list.pop(0)
        else:
            return None

    def __len__(self):
        return len(self.list)

    def add(self, value, priority):
        pair = (priority, value)
        i = self.loc_for_pair(pair)
        self.list.insert(i, pair)

    def update(self, value, old_priority, new_priority):
        old_pair = (old_priority, value)
        i = self.loc_for_pair(old_pair)
        del self.list[i]
        self.add(value, new_priority)

    def loc_for_pair(self, pair):
        # The range where it could be is [lower, upper)
        # Start with the whole list
        lower = 0
        upper = len(self.list)

        while upper > lower:
            next_split = (upper + lower) // 2
```

```
    v = self.list[next_split]
    if pair < v: # pair is to the left
        upper = next_split
    elif pair > v: # pair is to the right
        lower = next_split + 1
    else: # Found pair!
        return next_split
return lower
```

If you try running it now, it should work perfectly.

Now you have a graph class that would find the cheapest path quickly even if it had thousands of nodes with thousands of edges.

# Other Graph Algorithms

Now that you are familiar with Dijkstra's algorithm for finding the shortest path in a graph, then you're well-equipped to understand more graph algorithms. This document will discuss two other important algorithms: Depth-First Search (DFS) and the Bellman-Ford algorithm.

## 4.1 Depth-First Search

Depth-First Search (DFS) is an algorithm for traversing or searching tree or graph data structures. DFS uses a stack (or sometimes recursion which uses the system stack implicitly) to explore the graph in a depthward motion until it hits a node with no unvisited adjacent nodes, then it backtracks.

The procedure is as follows:

1. Push the root node into the stack.
2. Pop a node from the stack, and mark it as visited.
3. Push all unvisited adjacent nodes into the stack.
4. Repeat steps 2 and 3 until the stack is empty.

DFS is particularly useful for solving problems such as connected-component detection in graphs and maze-solving.

## 4.2 Bellman-Ford Algorithm

The Bellman-Ford algorithm is another shortest path algorithm like Dijkstra's. However, unlike Dijkstra's algorithm, Bellman-Ford can handle graphs with negative weight edges.

The algorithm works as follows:

1. Assign a tentative distance value for every node: set it to zero for our initial node and to infinity for all other nodes.
2. For each edge  $(u, v)$  with weight  $w$ , if the current distance to  $v$  is greater than the

distance to  $u$  plus  $w$ , update the distance to  $v$  to be the distance to  $u$  plus  $w$ .

3. Repeat the previous step  $|V| - 1$  times, where  $|V|$  is the number of vertices in the graph.
4. After the above steps, if you can still find a shorter path, there exists a negative cycle.

If the graph does not contain a negative cycle reachable from the source, the shortest paths are well-defined, and Bellman-Ford will correctly calculate them. If a negative cycle is reachable, no solution exists, but Bellman-Ford will detect it.



# Bayesian Networks

A Bayesian network, also known as a Bayes network, belief network, or decision network, is a probabilistic graphical model that represents a set of variables and their conditional dependencies via a directed acyclic graph (DAG).

## 5.1 Components

A Bayesian Network consists of two main components:

1. A directed acyclic graph (DAG) where each node represents a variable, and the absence or presence of a directed edge between nodes denotes the conditional dependence or independence respectively between the variables.
2. A conditional probability table (CPT) associated with each node which contains the conditional probability distribution of that node given its parents in the DAG.

## 5.2 Inferences

Bayesian Networks are typically used for reasoning and making inferences under uncertainty. Given observations of a set of variables, we can compute the posterior probabilities of the other variables using Bayes' rule.

There are three main types of inferences that we can make:

- **Causal reasoning (prediction):** Given the causes, what are the effects?
- **Evidential reasoning (diagnosis):** Given the effects, what are the causes?
- **Intercausal reasoning (explaining away):** Given an effect and some of its causes, what can we say about the other causes?

## 5.3 Learning

Learning a Bayesian Network from data involves two main tasks:

- **Structure learning:** Determining the DAG structure that best fits the data.
- **Parameter learning:** Estimating the parameters (conditional probabilities) of the CPTs given the DAG and data.

# Answers to Exercises





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