## 1 The Union-Find Data Structure

To efficiently implement Kruskal's algorithm, we need to avoid looking for cycles every time. Instead, we shall think of the algorithm as maintaining a set of connected components. In each step, it tries to use an edge to merge two connected components. To store a connected component, we shall use nodes that represent each vertex.

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Initialize: Set P(v) \leftarrow v and L(v) \leftarrow 0 for all vertices v
Function Find(x)

while P(x) \neq x do

| x \leftarrow P(x) |
end
Output x

Function Merge(x,y)

r_x \leftarrow Find(x)
r_y \leftarrow Find(y)
if L(r_x) \geq L(r_y) then

| P(r_y) \leftarrow r_x |
If L(r_x) = L(r_y), set L(r_x) \leftarrow L(r_x) + 1
end
else
| P(r_x) \leftarrow r_y |
end
```

In order to run the algorithm efficiently, we need a data structure that can maintain the connected components of the graph. We do this using the following data structure. The set of connected components will be maintained using a directed rooted tree where every vertex has out-degree 1 except the root. The connected components will be represented by the root of each tree, so two vertices x, y are in the same component if and only if the root of their trees are the same (In the above algorithm this corresponds to Find(x) = Find(y). To merge the components of x, y we simply call Merge(x, y).

For every vertex u, we write L(u) to denote the length of the longest path that ends at u. So, for a root r, L(r) is the depth of the tree rooted at r. To merge two components with roots  $r_x, r_y$ , if  $L(r_x) > L(r_y)$  we make  $r_y$  point to  $r_x$ . Similarly, if  $L(r_y) > L(r_x)$  we make  $r_x$  point to  $r_y$ . If  $L(r_x) = L(r_y)$  we make an arbitrary choice, say  $r_y$  points to  $r_x$ , and then increase  $L(r_x)$  by 1. The following claim implies that any find operation runs in time  $O(\log n)$ .

Claim 1. For every vertex  $v, L(v) \leq \log n$ .

**Proof** We prove by induction that if the root of a component satisfies L(r) = k, then the corresponding component has at least  $2^k$  nodes. Thus if  $L(r) > \log n$  for some node r, then the corresponding component will have more than n nodes, which is a contradiction. We induct on the time.

**Base Case:** At the beginning all components satisfy L(v) = 0 and they all have a single node  $(2^0 = 1)$ .

**IH:** Suppose the claim holds up to some time t.

IS: If at time t+1 we call Find(x,y) it will not change any tree. So, suppose we call Merge(x,y). So, we merge the trees of  $r_x, r_y$ . If  $L(r_x) \neq L(r_y)$  then L(.) of the new root does not increase, during the merge so the conclusion follows immediately. Otherwise, we have  $L(r_x) = L(r_y)$ . Say both of them are equal to k. In this case, by IH, we know that the tree of  $r_x$  has at least  $2^k$  nodes and the tree of  $r_y$  has at least  $2^k$  nodes. So, the new tree has at least  $2^k + 2^k = 2^{k+1}$  nodes. So, we can rightfully increase L(.) of the new root by 1.

## 2 Removing Weight Distinction Assumption

So far, in the description of Kruskal's algorithm, we assumed that edge weights are distinct.

In this section we see that even if the edge weights are not distinct still, the Kruskal's algorithm will give us a MST. Note that if the edge weights are not distinct there could be *multiple* MSTs, and we are only going to find one of them by Kruskal's algorithm.

Claim 2. Suppose the edge weights are not distinct. Kruskal's algorithm still gives us an MST.

**Proof** Note that in the proof we can *no longer* use the Cycle and Cut properties, because these properties only hold/proven when the edge weights are distinct.

Suppose the edge weights are not distinct. Kruskal's algorithm sorts the edges such that

$$c_{e_1} \le c_{e_2} \le \dots \le c_{e_m}$$

and then finds a tree based on this order. Suppose T is the output of the algorithm. For the sake of contradiction, assume that T is not an MST, then there is a tree  $T^*$ , such that  $c(T^*) < c(T)$ .

Perturb the weights of the edges such that all of the above inequalities are strict, i.e.,

$$c'_{e_1} < c'_{e_2} < \dots < c'_{e_m}$$
.

In particular, define  $c'_{e_i} = c_{e_i} + i \cdot \epsilon$  for a very very small  $\epsilon > 0$ .

**Fact 3.** Since the ordering does not change, still T will be the output of the Kruskal's algorithm.

**Fact 4.** Because the edge weights are distinct the output of Kruskal's algorithm is the optimum tree.

Now, if  $\epsilon$  is very small, much smaller than  $(c(T) - c(T^*))/m^2$ , we have  $c'(T^*) < c'(T)$  This is because

$$c'(T^*) \le c(T^*) + \epsilon + 2\epsilon + \dots + m\epsilon \le m^2\epsilon$$
, and,  
 $c'(T) \ge c(T)$ .

Therefore,  $c'(T^*) < c'(T)$ , i.e., Kruskal's algorithm does not output the MST for the edge weights c' which are distinct. But this contradicts Fact 4. Contradiction! Therefore, T is an MST with respect to original weights c as well.