

Advanced Programming Techniques

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Weak and Strong Components

Reminder: Connected components (and so, weak components of directed graphs) can be efficiently computed in parallel.

→ The same is true for blocks (2-connected components).

- However, the situation looks much more complicated for **strong components**: the best known serial algorithms (Tarjan, Kosaraju, etc.) are built on DFS, which (we think) cannot be parallelized.

- Today's lecture: **parallel computation of strong components**.

→ Reduction to *Union-Find* operations.

Disjoint Sets

Reminder

A Disjoint Sets Data Structure maintains a collection of pairwise disjoint sets. It supports the following three basic operations:

- `makeset(x)`: If x is not already present in the collection, then add a new singleton set whose unique elements equals x .
- `find(x)`: outputs the unique identifier of the set containing x .
 - In general, `find(x)` outputs an element of the set, also called its “representative”.
 - In serial implementations, we may force this representative to have special properties (e.g., largest element in the set) with no computational overhead.
- `union(x,y)`: merge the respective sets of x,y into one.

Relation with connected components

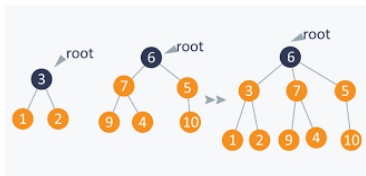
Consider the following algorithm on a graph G :

```
for all vertices  $v$  in parallel  
  makeset( $v$ )  
  
for all edges  $uv$  in parallel  
  union( $u, v$ )  
  
for all vertices  $v$  in parallel  
   $cc[v] = \text{find}(v)$ 
```

Consequence: a parallel implementation for Disjoint Sets leads to a new parallel implementation for connected components computation.

Serial implementation: Representing sets as trees

- The elements of each set are the nodes of a tree, whose root is the representative of this set.



- makeset(x):** create a new tree whose unique node is x
- find(x):** find the root of the tree containing x as a node (climb up)
- union(x,y):** link together the trees that contain x, y as nodes (one root becomes the child of the other).

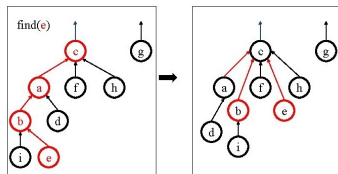
In a naive implementation, complexity depends on the height of the trees.

Improvements: Path Compression

Operation find

In order to access to the root (representative), we climb in the tree. *On our way, all visited nodes are reconnected as children of the root.*

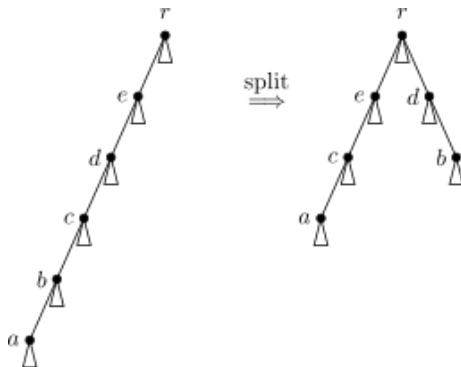
→ Speed-up of subsequent find operations.



Improvements: Path Splitting

Operation find

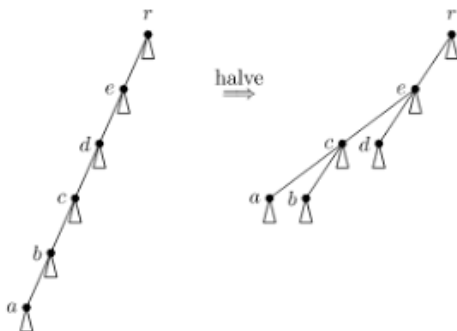
All nodes on the path to the root are reconnected to their grandparent.



Improvements: Path Halving

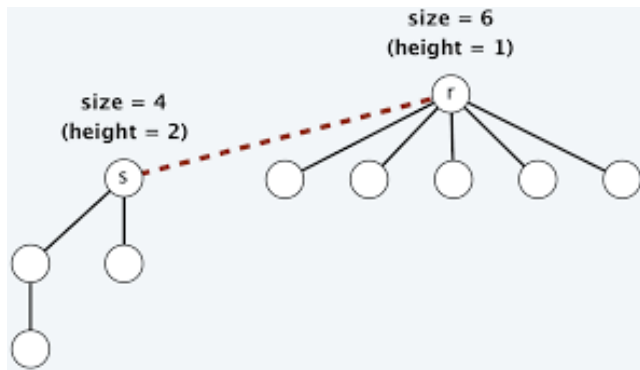
Operation find

On our way to the root, we reconnect each traversed node to its grandparent. Former parent nodes are skipped.



Improvements: Union by size

- Each node stores the size of its rooted subtree. If we merge two sets, then the root of the new set is the root of the biggest tree.

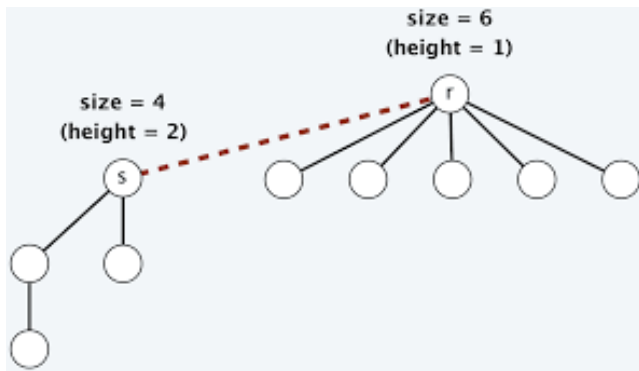


Remark: Ensures logarithmic depth.

Improvements: Union by rank

- Each node keeps a rank: **upper bound** on its depth. If we merge two sets, then the root of the new set is the root of larger rank.

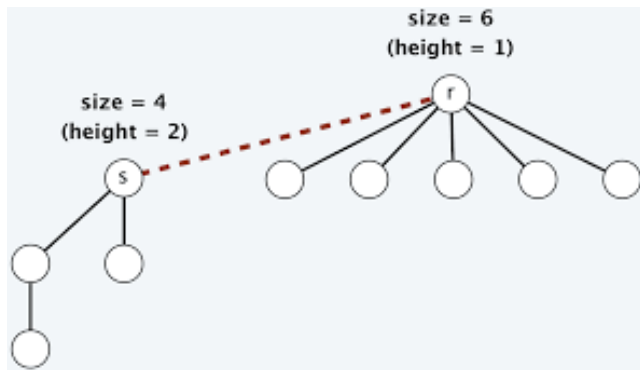
→ The real depth might be smaller than the rank because of path compression/splitting/halving. . .



Also ensures logarithmic depth.

Improvements: Union by **random index**

- Node first select random distinct indices. If we merge two sets, then the root of the new set is the root of larger (or smaller) index.



Also ensures logarithmic depth (in expectation).

Computation of a random index

Assumption: the universe of n elements is fixed, and known to us from the beginning.

- Every element just selects u.a.r. a number between 1 and n^3 .
- For any fixed element, the probability for another element to select its same index is at most $n \times 1/n^3 = 1/n^2$.
- Therefore, all indices are distinct with probability at least $1 - 1/n$.

Remark: can be done in parallel (assuming the random seed has no correlation with our scheduler).

Toward a parallel implementation: The CAS primitive

CAS(x, y, z):

Input: an address x in the shared memory; two values y,z.

Output: true if x was containing value y before the operation (after the operation, x stores the new value z); false otherwise.

This is an **atomic** compare-and-swap operation (it requires synchronization).

- OpenMP implementation:

```
#pragma omp atomic compare  
if(x == y) x = z;
```

Concurrent Disjoint Sets

The need for more operations

- In a serial implementation, we only have three operations: `makeset`, `find`, `union`.
- A `union` can be seen as two `find` operations (that can be done in parallel), followed by a `link` operation between two roots. **In a concurrent setting, a `link` can sometimes fail!** Therefore, we need to define `link` separately from `union`, and to repeatedly call `link` until it becomes successful.
- We mostly rely on `find` in order to decide whether two elements are in the same set. However, this approach may fail in a concurrent setting. Therefore, we need to define a `same_set` operation separately.

Link

- Linking by size or rank is problematic because of concurrent operations. However, a simple implementation can still be achieved using Linking by random index.

```
1: procedure link( $u, v$ )  
2:   if  $u < v$  then CAS( $u.p, u, v$ )  
3:   else CAS( $v.p, v, u$ )
```

Remark: the operation fails if $u \neq u.p$, i.e., u is no more a root because of concurrent operations.

Find (1/2)

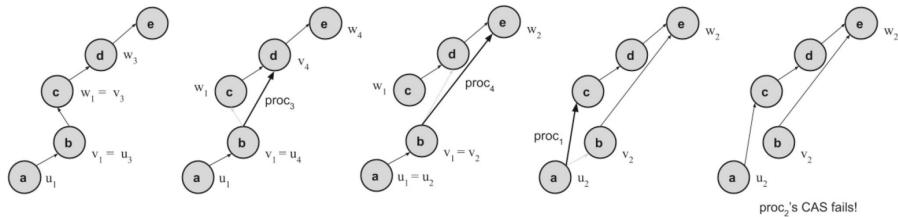
- Path compaction (reconnection of nodes on the path to the root) may fail in a concurrent setting.
- However, “*intuitively*”, when it does fail, it means that compaction has been done by another processor.

⇒ we continue climbing up into the tree *even if the compaction has failed!*

```
1: procedure find(x)  
2:    $u \leftarrow x; v \leftarrow u.p; w \leftarrow v.p$   
3:   while  $v \neq w$  do  
4:      $\text{CAS}(u.p, v, w); u \leftarrow v; v \leftarrow u.p; w \leftarrow v.p$   
5:   return  $v$ 
```


Find (2/2)

Example



Same_set

- Concurrent calls to `find` may output two different nodes, even if they belong to the same set.
- We repeatedly need to check whether the root has changed since our last call to `find`.

```
1: procedure same-set( $x, y$ )
2:    $u \leftarrow \text{find}(x); v \leftarrow \text{find}(y)^*$ 
3:   while  $u \neq v$  do
4:      $w \leftarrow u.p$ 
5:     if  $u = w$  then return false
6:      $u \leftarrow \text{find}(u); v \leftarrow \text{find}(v)^*$ 
7:   return true
```

Union

- Repeated calls to `link` until it becomes successful.

```
1: procedure unite( $x, y$ )  
2:    $u \leftarrow \text{find}(x); v \leftarrow \text{find}(y)$   
3:   while  $u \neq v$  do  
4:      $\text{link}(u, v)$   
5:      $u \leftarrow \text{find}(u); v \leftarrow \text{find}(v)$ 
```

Strong components




Purdom-Monroe Algorithm

```
1:  $\forall v \in V : \mathcal{S}(v) := \{v\}$ 
2: DEAD := VISITED :=  $\emptyset$ 
3:  $R := \emptyset$ 
4: SETBASED( $v_0$ )
5: procedure SETBASED( $v$ )
6:   VISITED := VISITED  $\cup \{v\}$ 
7:    $R.PUSH(v)$ 
8:   for each  $w \in POST(v)$  do
9:     if  $w \in DEAD$  then continue ..... [already completed SCC]
10:    else if  $w \notin VISITED$  then ..... [unvisited node  $w$ ]
11:      SETBASED( $w$ )
12:    else while  $\mathcal{S}(v) \neq \mathcal{S}(w)$  do ..... [cycle found]
13:       $r := R.POP()$ 
14:      UNITE( $\mathcal{S}, r, R.TOP()$ )
15:   if  $v = R.TOP()$  then ..... [completely explored SCC]
16:     report SCC  $\mathcal{S}(v)$ 
17:     DEAD := DEAD  $\cup \mathcal{S}(v)$ 
18:      $R.POP()$ 
```

- *Partial* strong components are stored. This is \neq from Tarjan's algorithm where each strong component is fully computed at once.

Strong components

Purdom-Monroe Algorithm

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2: DEAD := VISITED :=  $\emptyset$ 
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4: SETBASED( $v_0$ )
5: procedure SETBASED( $v$ )
6:   VISITED := VISITED  $\cup \{v\}$ 
7:    $R.PUSH(v)$ 
8:   for each  $w \in \text{POST}(v)$  do
9:     if  $w \in \text{DEAD}$  then continue
10:    else if  $w \notin \text{VISITED}$  then
11:      SETBASED( $w$ )
12:    else while  $\mathcal{S}(v) \neq \mathcal{S}(w)$  do  same_set(u,v)
13:       $r := R.POP()$ 
14:      UNITE( $\mathcal{S}, r, R.TOP()$ )  union
15:   if  $v = R.TOP()$  then
16:     report SCC  $\mathcal{S}(v)$ 
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```

- *Partial* strong components are stored. This is \neq from Tarjan's algorithm where each strong component is fully computed at once.

UFSCC Algorithm

A modified Purdom-Monroe Algorithm is executed in parallel by all processors, with one shared Disjoint Sets data structures for partial strong components.

```
1:  $\forall v \in V : S(v) := \{v\}$  ..... [global  $S$ ]
2:  $DEAD := DONE := \emptyset$  ..... [global  $DEAD$  and  $DONE$ ]
3:  $\forall p \in [1 \dots p] : R_p := \emptyset$  ..... [local  $R_p$ ]
4:  $UFSCC_1(v_0) || \dots || UFSCC_p(v_0)$ 
5: procedure  $UFSCC_p(v)$ 
6:    $R_p.PUSH(v)$ 
7:   while  $v' \in S(v) \setminus DONE$  do
8:     for each  $w \in RANDOM(POST(v'))$  do
9:       if  $w \in DEAD$  then continue ..... [DEAD]
10:      else if  $\nexists w' \in R_p : w \in S(w')$  then ..... [NEW]
11:         $UFSCC_p(w)$ 
12:      else while  $S(v) \neq S(w)$  do ..... [LIVE]
13:         $r := R_p.POP()$ 
14:         $UNITE(S, r, R_p.TOP())$ 
15:       $DONE := DONE \cup \{v'\}$ 
16:   if  $S(v) \not\subseteq DEAD$  then  $DEAD := DEAD \cup S(v)$ ; report SCC  $S(v)$ 
17:   if  $v = R_p.TOP()$  then  $R_p.POP()$ 
```

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```

repeated iterations
over the partial strong
component of v
(because of parallelism)

UFSCC Algorithm

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9:       if  $w \in DEAD$  then continue
10:      else if  $w' \in R_p \wedge w \in S(w')$  then
11:         $UFSCC_p(w)$ 
12:        else while  $S(v) \neq S(w)$  do
13:           $r := R_p.POP()$ 
14:           $UNITE(S, r, R_p.TOP())$ 
15:         $DONE := DONE \cup \{v'\}$ 
16:   if  $S(v) \not\subseteq DEAD$  then  $DEAD := DEAD \cup S(v)$ ; report SCC  $S(v)$ 
17:   if  $v = R_p.TOP()$  then  $R_p.POP()$ 
```

w' was pushed in the stack before v. Therefore, v and w are in the same scc.

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9:       if  $w \in DEAD$  then continue
10:      else if  $\nexists w' \in R_p : w \in S(w')$  then
11:         $UFSCC_p(w)$ 
12:      else while  $S(v) \neq S(w)$  do
13:         $r := R_p.POP()$ 
14:         $UNITE(S, r, R_p.TOP())$  critical section
15:         $DONE := DONE \cup \{v\}$ 
16:      if  $S(v) \not\subseteq DEAD$  then  $DEAD := DEAD \cup S(v)$ ; report SCC  $S(v)$ 
17:      if  $v = R_p.TOP()$  then  $R_p.POP()$ 
```

Some Terminology

For a vertex v in a directed graph G , let us define:

- $\text{pred}(G, v)$: all vertices that can reach v in G
- $\text{desc}(G, v)$: all vertices that can be reached from v in G

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Proposition 2: every strong component must either be fully in $\text{pred}(G, v)$, be fully in $\text{desc}(G, v)$, or not intersect $\text{pred}(G, v) \cup \text{desc}(G, v)$.

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Proposition 2: every strong component must either be fully in $\text{pred}(G, v)$, be fully in $\text{desc}(G, v)$, or not intersect $\text{pred}(G, v) \cup \text{desc}(G, v)$.

Remark: $\text{pred}(G, v)$ and $\text{desc}(G, v)$ can be computed using BFS (no need for DFS).

Forward-Backward Algorithm

```
1: procedure FWBW( $G$ )
2:   if  $G = \emptyset$  then
3:     return  $\emptyset$ 
4:   select pivot  $v$ 
5:    $D \leftarrow DESC(G, v)$ 
6:    $P \leftarrow PRED(G, v)$ 
7:    $R \leftarrow (G \setminus (P \cup D))$ 
8:    $S \leftarrow (D \cap P)$ 
9:    $FWBW(D \setminus S)$ 
10:   $FWBW(P \setminus S)$ 
11:   $FWBW(R)$ 
```

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10:   $FWBW(P \setminus S)$ 
11:   $FWBW(R)$ 
```

Two graph searches (BFS)
in parallel

Forward-Backward Algorithm

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8:    $S \leftarrow (D \cap P)$ 
9:   FWBW( $D \setminus S$ )
10:  FWBW( $P \setminus S$ )
11:  FWBW( $R$ )
```

Can be computed in parallel

Forward-Backward Algorithm

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11:   $FWBW(R)$ 
```

In parallel

Coloring Algorithm

- The graph is vertex coloured. Each colour c is ‘owned’ by a vertex (the one c_v of which c represents the identifier).
- **Key invariant:** all vertices with same colour c are contained in $\text{desc}(G, c_v)$. Furthermore, no vertex of $\text{desc}(G, c_v)$ with another colour than c can be in the same strong component as c_v .

```
1: while  $G \neq \emptyset$  do
2:   initialize  $\text{colors}(v_{id}) = v_{id}$ 
3:   while at least one vertex has changed colors do
4:     for all  $v \in G$  do
5:       for all  $u \in N(v)$  do
6:         if  $\text{colors}(v) > \text{colors}(u)$  then
7:            $\text{colors}(u) \leftarrow \text{colors}(v)$ 
8:   for all unique  $c \in \text{colors}$  do
9:      $\text{SCC}(c_v) \leftarrow \text{PRED}(G(c_v), c)$ 
10:     $G \leftarrow (G \setminus \text{SCC}(c_v))$ 
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8:   for all unique  $c \in \text{colors}$  do In parallel
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In parallel

Questions

