# Linear Time Distributed Swap Edge Algorithms<sup>\*</sup>

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Abstract. In this paper, we consider the all best swap edges problem in a distributed environment. We are given a 2-edge connected positively weighted network X, where all communication is routed through a rooted spanning tree T of X. If one tree edge  $e = \{x, y\}$  fails, the communication network will be disconnected. However, since X is 2-edge connected, communication can be restored by replacing e by non-tree edge e', called a swap edge of e, whose ends lie in different components of T - e. Of all possible swap edges of e, we would like to choose the best, as defined by the application. The all best swap edges problem is to identify the best swap edge for every tree edge, so that in case of any edge failure, the best swap edge can be activated quickly. There are solutions to this problem for a number of cases in the literature. A major concern for all these solutions is to minimize the number of messages. However, especially in fault-transient environments, time is a crucial factor. In this paper we present a novel technique that addresses this problem from a time perspective; in fact, we present a distributed solution that works in linear time with respect to the height h of T for a number of different criteria, while retaining the optimal number of messages. To the best of our knowledge, all previous solutions solve the problem in  $O(h^2)$  time in the cases we consider.

## 1 Introduction and Preliminaries

For a communication network, low cost and high reliability can be conflicting goals. For example, a *spanning tree* of a network could have minimum cost, but will not survive even a single link failure. We consider the problem of restoring connectivity when one link of a spanning tree fails.

One recent technique, particularly efficient in case of transient faults, consists in pre-computing a *replacement spanning tree* for each possible link or node failure, by computing the best replacement edge (or edges) which reconnects the tree. A number of studies have been done for this problem, both for the sequential [1–6] and distributed [7–10] models of computation, for different types of spanning trees and failures.

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In this paper, we consider the all best swap edges problem in the distributed setting. We are given a positively weighted 2-edge connected network X of processes, where w(x, y) denotes the weight of any edge  $\{x, y\}$  of X, together with a spanning tree T of X, rooted at a process r. Suppose that all communication between processes is routed through T. If one tree edge  $e = \{x, p(x)\}$  fails (where p(x) denotes the parent of x in T) we say that x is the point of failure. Since X is 2-edge connected, communication can be restored by replacing e by an edge e' of X whose ends lie in different components of T - e. We call such an edge e' a swap edge of x (or a swap edge of e), and we define SwapEdges(x)(or SwapEdges(e)) to be the set of all swap edges of x (refer to the example depicted in Figure 1.(b) and (c)). Of all possible swap edges of x, we would like to choose the best, as defined by the application. The all best swap edges problem is to identify the best swap edge for every tree edge, so that in case of any edge failure, the best replacement edge can be activated quickly.

Notation. Given T a spanning tree of X, we refer to an edge of T as a tree edge, and any other edge of X as a cross edge (see also Figure 1.(a)).

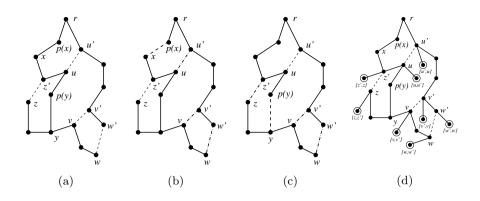
If  $x \neq r$  is a process, we denote the set of children of x by Chldrn(x), and the subtree of T rooted at x by  $T_x$ ; the *level* of a process x is defined to be the *hop-distance* from x to r. We write  $x \leq y$  or  $y \geq x$  to indicate that x is an ancestor of y, *i.e.*,  $y \in T_x$ , and x < y or y > x if x is a proper ancestor of y.

If S is any subgraph of X, we let  $path_{s}(x, y)$  denote the shortest (least weight) path through S from x to y, and let  $W_{S}(x, y)$  denote the weighted length of  $path_{s}(x, y)$ . (We write simply path(x, y) and W(x, y) if S is understood.)

We will denote by  $T^*$  the *augmented tree*, whose nodes consist of all processes of T, together with a node for each directed cross edge of T, which we call an *augmentation node* of  $T^*$ . (See Figure 1.(d).) In particular, if  $\{y, y'\}$  is a cross edge in T, we will denote by [y, y'] and [y', y] its corresponding nodes in  $T^*$ ; the parent of [y, y'] is y. For any process x, define  $T^*_x$  to be the subtree of the augmented tree rooted at x; in particular,  $T^*_x$  consists of  $T_x$  together with all the augmentation nodes [y, y'] such that  $y \in T_x$ .

Related Work and Our Contribution. In [2, 9], several different criteria for defining the "best" swap edge for a tree edge e have been considered. In each case, the best swap edge for e is that swap edge e' for which some penalty function F is minimized. We consider three penalty functions in this paper. In each case, let T' = T - e + e' be the spanning tree of X obtained by deleting e and adding e', where  $e = \{x, p(x)\}$  is a tree edge,  $y \in T_x$ , and  $e' = \{y, y'\}$  a swap edge for e.

- 1.  $F_{\text{wght}}(x, y, y') = w(e')$ , the weight of the swap edge. Note that if T is a minimum spanning tree of X and e' is that swap edge for e such that w(e') is minimum, then T' = T e + e' is a minimum spanning tree of X e.
- 2.  $F_{\text{dist}}(x, y, y') = W_{T'}(r, x)$ , the distance from the root to the point of failure in T'.
- 3.  $F_{\max}(x, y, y') = \max \{ W_{T'}(r, u) : u \in T_x \}$ , the maximum distance, in T', from the root to any process in  $T_x$ .



**Fig. 1.** (a) An example of a network X and its spanning tree T: Tree edges are bold, cross edges are dotted. (b) Failure at x.  $\{u, u'\}$ ,  $\{v, v'\}$ , and  $\{w, w'\}$  are the swap edges of x. (c) Failure at y.  $\{v, v'\}$ ,  $\{w, w'\}$ , and  $\{z, z'\}$  are the swap edges of y. (d) The augmented tree of T; the augmentation nodes are double circled.

If F is any of the above penalty functions, we define  $F(x, y, y') = \infty$  for any  $\{y, y'\}$  which is not a swap edge of x. The output of the problem is then  $M_F(x) = \min \{F(x, y, y') : (y, y') \in T_x^*\}$ .

In [7], Flocchini *et al.* give an algorithm for solving the  $F_{\text{dist}}$  version of all best swap edge problem. In [9], Flocchini *et al.* give a general algorithm for the all best swap edges problem, and then give specific versions of the technique for the  $F_{\text{max}}$  version. In [8], the  $F_{\text{wght}}$  version is solved both for the failure of a link and for the failure of a node and all its incident links.

All the above mentioned distributed solutions have the same general form, and have message complexity  $O(n^*)$ , with  $n^*$  the number of edges of the transitive closure of  $T_r \setminus \{r\}$ . The time complexity of each is  $O(h^2)$ , where h is the height of T. In particular, each of those solutions consists of two waves for each level  $\ell$  of T, with  $1 \le \ell \le h$ : A broadcast wave followed by a converge ast wave. In particular, the general schema of previous solutions consists of two nested loops, where the outer loop is indexed by  $\ell$ , and for each  $\ell$ , the inner loop computes  $M_F(x)$  for all x at level  $\ell$  using two waves; a top-down wave that computes F(x, y, y') for all  $(y, y') \in T_x^*$ , and a bottom-up wave that computes  $M_F(x)$ . Each wave takes O(h)time in the worst case, hence the overall strategy leads to a final cost in time of  $O(h^2)$ . This is mainly due to the fact that the waves needs to be executed one after the other. In this paper we present a novel technique that finds a solution in linear time, for each of the penalty functions listed above. In particular, our strategy distributes the information and the computation among processes so that the waves can be pipelined. This reduces the final time of the execution to O(h), using the same number of messages as the previous solutions.

As a final remark, we note that in [2], Gfeller *et al.* study the problem of finding the optimal swap edges of a minimum spanning tree having minimum diameter: They provide a distributed algorithm that already works in linear time. The general technique presented here can be also adapted to this case.

The paper is organized as follows. The overall structure of our paradigm is given in Section 2. The various phases are described in Sections 3, 4, and 5. Due to space constraints, some of the proofs are given in the appendix.

### 2 The Linear Time Solution

In this section, we present the strategy that allows to devise O(h)-time distributed algorithms to solve the five versions of the all best swap edges problem introduced in the previous section. We call these algorithms LINEAR<sub>dist</sub>, LINEAR<sub>wght</sub>, LINEAR<sub>max</sub>, respectively. Each can be considered to be a version of a general algorithm, which we call LINEAR, whose structure is given as Algorithm 1. LINEAR is structured in phases; the actual number of phases depends on the specific version of the problem. However, in all cases, the number of phases is at least three: a *preprocessing* phase, a *ranking* phase, and an *optimization* phase. In the last optimization phase, a piece of information, denoted by *up\_package*( $y, \ell$ ), is computed in a convergecast wave. The content of this package is different for each of the versions of the problem (details in Section 5).

- 1: Preprocessing Phase
- 2: Ranking Phase
- 3: If LINEAR<sub>max</sub> Then Additional Critical Level Phase(s)
- 4: Optimization Phase

Each of the phases of LINEAR uses at most  $O(\delta_x)$  space for each x, where  $\delta_x$  is the degree of x. The space complexity of LINEAR is thus  $O(\delta_x)$  for each x.

Our linear time algorithms make use of the concept of *critical level*. Informally, a critical level function is a function that can be computed top-down, which enables another function – whose computation would otherwise require independent top-down followed by bottom-up waves for all processes – to be computed in a single bottom-up wave for each process, thus allowing the waves to be pipelined. In particular, for each of the versions of the best swap edge problem we consider, one or more critical levels are computed, depending on the specific penalty function. Due to space constraints, all the proofs will be omitted.

The Role of Critical Levels. A critical level function is a function  $\Lambda$  on the augmentation nodes of  $T^*$  such that  $0 \leq \Lambda(y, y') \leq y.level$ , and which aids in the computation of F(x, y, y') for any  $x \leq y$ . More specifically, the computation of F(x, y, y') contains a branch which depends on the comparison between x.level and  $\Lambda(y, y')$ . For example, the function rank, defined in Section 4, has the property that  $F(x, y, y') = \infty$  if and only if  $rank(y, y') \geq x.level$ , where F is any one of the penalty functions defined above.

## 3 Preprocessing Phase

In the preprocessing phase, which takes O(h) time, each process x computes and retains a set of variables, some of which are the same as in [7, 9]. All the variables listed below are needed for LINEAR<sub>max</sub>, but only *level*, *index*, and *depth* are needed for LINEAR<sub>max</sub> and LINEAR<sub>dist</sub>.

- 1. x.level, the level of x, which is the hop-distance from r to x.
- 2.  $x.index = (x.pre\_index, x.post\_index)$ , the index of x, where  $x.pre\_index$  is the index of x in the pre-order visit of T, and  $x.post\_index$  is the index of x in the reverse postorder of T (see Figure 2(a)).
- 3. x.depth = W(r, x), the depth of x.
- 4.  $x.height = \max \{ W(x, u) : u \in T_x \}, \text{ the height of } x.$
- 5.  $x.best\_child$ , the best child of x, defined to be the process  $y \in Chldrn(x)$  such that w(x, y) + y.height > w(x, z) + z.height for any other child z of x. Note that, since we use a strict inequality in this definition, a process can have at most one best child. If  $Chldrn(x) = \emptyset$ , or if there is more than one choice of y for which w(x, y) + y.height is maximum,  $best\_child(x)$  is undefined.
- 6. x.eta, for  $x \neq r$ , the largest weight of any path in  $T_{p(x)} T_x$  from p(x); that is,  $x.eta = \max \{w(p(x), y) + y.height : y \neq x \text{ and } y \in Chldrn(p(x))\}$ . If x is the only child of its parent, then x.eta defaults to 0.
- 7. x.secondary\_height, the length of the longest path which does not contain  $x.best\_child$  from x to any leaf of  $T_x$ . In the case that  $x.best\_child$  is undefined, let  $x.secondary\_height = x.height$ .

Note that all of the above variables can be computed with a constant number of broadcast and converge ast waves, in O(h) total time.

## 4 Ranking Phase

The ranking phase is the same for all versions of the best swap edge problem. In this phase, we compute the *rank* of every cross edge  $\{y, y'\}$ , defined to be the level of the nearest common ancestor of y and y' in T. This value is stored by both y and y'. Ranks are used to distinguish swap edges of x from other cross edges in  $T_x^*$ .

Remark 1.

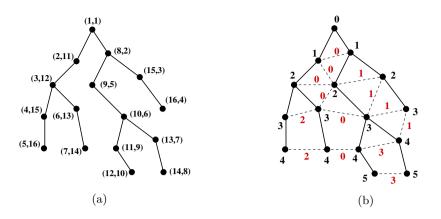
(a) A process x is an ancestor of y if and only if  $x.index \le y.index$ .

(b) If  $[y, y'] \in T_x^*$ , then  $\{y, y'\} \in SwapEdges(x)$  if and only if  $x.index \leq y'.index$ .

From the previous remark, it follows that:

Remark 2. Let  $x \neq r$  be a process and  $e' = \{z, z'\}$  a cross edge, where  $z \in T_x$ . Then, e' is a swap edge for x if and only if rank(z, z') < x.level.

The ranking phase is given as Algorithm 2. In particular, there is a main loop that cycles over the levels of the tree in increasing order. The phase consists of



**Fig. 2.** (a) Processes are labeled with their indices. A process x is an ancestor of y if and only if  $x.index \leq y.index$ . (b) Levels of processes and ranks of cross edges.

Algorithm 2. Ranking Phase: Rank of every Cross Edge of T is Computed			
1: <b>I</b>	For $0 \le \ell \le h$ in increasing order <b>Do</b> % Wave $\ell$ %		
2:	For all y such that $y.level \ge \ell$ in top-down order <b>Do</b>		
3:	If $y.level = \ell$ Then $ancestor\_index(y, \ell) \leftarrow y.index$		
4:	<b>Else</b> ancestor_index $(y, \ell) \leftarrow ancestor_index(p(y), \ell)$		
5:	For all cross edges $\{y, y'\}$ Do		
6:	If $y'$ .index $\geq$ ancestor_index $(y, \ell)$ Then $rank(y, y') \leftarrow \ell$		

a top-down wave for each  $0 \leq \ell \leq h$ , denoted by Wave  $\ell$ . For each  $\ell$ , the inner loop computes, for each process y whose level is greater than or equal to  $\ell$ , the value ancestor\_index $(y, \ell)$ , which is x.index where x is the ancestor of y at level  $\ell$ . Then, for each  $[y, y'] \in T_x^*$ , the value  $\ell$  is assigned to rank(y, y') if  $y' \notin T_x$ , *i.e.*, y'.index  $\geq$  ancestor\_index $(y, \ell)$  (refer to Remark 1).

The inner loop is executed as a top-down wave; hence the waves can be pipelined, so that the total time of the ranking phase is O(h).

**Lemma 1.** If  $rank(y, y') = \ell$ , then, for all  $\ell' \leq \ell$ , the computed value of rank(y, y') will be set to  $\ell'$  during Wave  $\ell'$  of Algorithm 2 and thus the final computed value of rank(y, y') will be  $\ell$ .

### 5 Optimization Phase

The optimization phase is implemented as a bottom-up wave for each level  $\ell$ . (Refer to Algorithm 1.) In this phase, all best swap edges are computed. In particular, the phase consists of an outer loop, indexed by decreasing values of  $1 \leq \ell \leq h$ , where each iteration consists of an inner loop which computes  $M_F(x)$ for all x at level  $\ell$ . For each x such that  $x.level = \ell$ , the inner loop consists of a convergecast wave, which computes a set of variables we call  $up\_package(y, \ell)$  for each  $y \in T_x$ ; each process y is able to compute  $up\_package(y, \ell)$  by using the information computed and stored at y during the earlier phases, as well as the contents of  $up\_package(z, \ell)$  received from all  $z \in Chldrn(y)$ . The final value of  $M_F(x)$  is then computed using  $up\_package(x, \ell)$ . To save space, each up-package is deleted as soon as it is no longer needed. The convergecast waves can be pipelined, and thus the entire optimization phase can be executed in O(h) time.

The specific content of  $up\_package(y, \ell)$  depends on the specific version of LINEAR that is solved.

#### 5.1 LINEAR<sub>wght</sub> and LINEAR<sub>dist</sub>

For each  $\ell \geq 1$  and each  $y \in T$  at level  $\geq \ell$ , let x be the unique ancestor of y at level  $\ell$ , and let  $e = \{x, p(x)\}$ . We define  $Swap_N(y, \ell)$  to be the set of all neighbors y' of y such that  $\{y, y'\}$  is a swap edge for e. In order to compute this set, the test established by Remark 2 is used.

For both LINEAR<sub>wght</sub> and LINEAR<sub>dist</sub>,  $up\_package(y, \ell)$  consists of just the value  $sbtree\_min(y, \ell)$ , defined as follows. If x is the unique ancestor of y at level  $\ell$ , then

- 1. In LINEAR<sub>wght</sub>: sbtree\_min(y,  $\ell$ ) = min{w(z, z')}, such that  $(z, z') \in T_y^* \cap SwapEdges(x)$ .
- 2. In LINEAR<sub>dist</sub>: sbtree\_min(y,  $\ell$ ) = min{W(x, z) + w(z, z') + z'.depth}, such that  $(z, z') \in T_y^* \cap SwapEdges(x)$ .

At the end of the iteration for  $\ell$ , the value of  $M_F(x)$  is set to  $sbtree\_min(x, \ell)$  for all x at level  $\ell$ .

The pseudo-code of the optimization phase, for the functions  $F_{\text{wght}}$  and  $F_{\text{dist}}$  is given as Algorithm 3 and 4, respectively. In both cases, the waves of the optimization phase are pipelined, permitting the total time complexity of the phase to be O(h).

Concerning the number of messages of both LINEAR<sub>wght</sub> and LINEAR<sub>dist</sub>, note that the information sent along the tree either in the ranking phase or in the optimization phase, is composed of messages of constant size. In the ranking phase, the information consists of node indices, and in the optimization phase of "subtree minimum" values. Thus, the communication complexity, corresponding to the transitive closure of the tree edges, is  $O(n^*)$  (limited by  $O(n^2)$ ) in both cases.

### 5.2 LINEAR<sub>max</sub>

If  $S \subseteq X$  is connected and  $x \in S$ , define  $radius(S, x) = \max \{W_S(x, s) : s \in S\}$ , the radius of S based at s. Note that, we can write  $F_{\max}(x, y, y') = \max\{W_{T'}(r, u) : u \in T_x\} = radius(T_x, y) + w(y, y') + y'.depth$  if  $\{y, y'\} \in SwapEdges(x)$ . Thus, in the case of LINEAR<sub>max</sub> we face with the problem of computing  $radius(T_x, y)$ : This computation is handled by an additional phase before the actual optimization phase. In this phase, we compute a variable called  $critical\_level(y)$ , for all  $y \in T_x$ .

#### **Algorithm 3.** Algorithm LINEAR<sub>webt</sub>

1: Prep	rocessing	phase
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- 2: Ranking phase
- 3: For all  $1 < \ell < h$  Do %*Optimization Phase*%
- 4: For all y such that y.level >  $\ell$  in bottom-up order **Do**
- 5: $Swap_N(y, \ell) \leftarrow \{y' : \{y', y\} \text{ is a cross edge and } rank(y, y') < \ell\}$
- 6:
- $sbtree\_min(y,\ell) \leftarrow \min \begin{cases} w(y,y') : y' \in Swap\_N(y,\ell) \\ \min \{sbtree\_min(z,\ell) : z \in Chldrn(y)\} \end{cases}$
- For all x such that  $x.level = \ell$  Do 7:
- 8:  $M_F(x) = sbtree\_min(x, \ell)$

#### Algorithm 4. Algorithm LINEAR<sub>dist</sub>

lS€

- 2: Ranking phase
- 3: For all  $1 < \ell < h$  Do %*Optimization Phase*%
- For all y such that y.level >  $\ell$  in bottom-up order **Do** 4:
- $\mathit{Swap\_N}(y,\ell) \gets \{y': \{y',y\} \text{ is a cross edge and } \mathit{rank}(y,y') < \ell\}$ 5:
- 6:
- $sbtree\_min(y, \ell) \leftarrow \min \begin{cases} w(y, y') + depth(y') : y' \in Swap\_N(y, \ell) \\ \min \{w(y, z) + sbtree\_min(z, \ell) : z \in Chldrn(y)\} \end{cases}$
- For all x such that  $x.level = \ell$  Do 7:
- 8:  $M_F(x) = sbtree\_min(x, \ell)$

Additional Critical Level Phase. For  $y \in T_x$ , define  $\mu(y, x)$  to be the weight of the longest path in  $T_x$  from y to any node of  $T_x - T_y$ . We let  $\mu(x, x) = 0$  by default. It follows from these definitions that

$$radius(T_x, y) = \max \begin{cases} y.height\\ \mu(y, x) \end{cases}$$
(1)

Since we want LINEAR to use only constant space per process, y can hold only  $O(\delta_y)$  values; hence, it could not be possible for y to store all the values  $\{\mu(y, x) : x \leq y\}$ . We tackle this problem by executing in LINEAR<sub>max</sub> an extra phase before the optimization phase (called *critical level phase* in Algorithm 1). In particular, as the convergecast wave moves up the tree, we compute the *critical level* of y, that determines not the actual value of  $radius(T_x, y)$ , but rather which of the two choices given in Equation (1) is larger, together with enough additional information to calculate the actual value of  $M_F(x)$  when the wave reaches x.

We now explain critical levels in greater detail. Let

$$critical\_level(y) = \min \{x.level : y \in T_x \text{ and } radius(T_x, y) = y.height\}.$$

Note that  $critical\_level(y) = \min \{x.level : y \in T_x \text{ and } \mu(y, x) \le y.height\}.$ 

**Lemma 2.** For any processes  $x' \le x \le y$ ,  $\mu(y, x') \ge \mu(y, x)$ .

**Corollary 1.** If  $y \in T_x$ , then  $radius(T_x, y) = y$ .height if and only if x.level  $\geq$  $critical\_level(y).$ 

Critical levels are calculated by Algorithm 5. Recall, from Section 3, that y.eta, for  $y \neq r$ , is the largest weight of any path in  $T_{p(y)} - T_y$  from p(y); this value is computed during the preprocessing phase. Note that, once again, the waves of the inner loop of Algorithm 5 can be pipelined, so that the total time required for this phase is again O(h).

Algorithm 5. Critical Level Phase			
1:	For $0 \leq \ell \leq h$ in decreasing order <b>Do</b> % Wave $\ell$ %		
2:	For all x such that $x.level = \ell$ concurrently <b>Do</b>		
3:	$\mu(x,x) \leftarrow 0$		
4:	For all $y \in T_x - x$ in top down order <b>Do</b>		
5:	$\mu(y, x) \leftarrow \max \begin{cases} \mu(p(y), x) + w(y, p(y)) \\ y.eta \end{cases}$		
6:	If $\mu(y, x) \leq y.height$ Then $critical\_level(y) \leftarrow \ell$		

*Optimization Phase.* Before introducing the optimization phase, we need to introduce the notion of *Spine*, which is strictly related to the notion of critical level. (Refer also to Figure 3.)

**Definition 1 (Spine).** Given any process x, we define the Spine of x:

 $Spine(x) = \{y \in T_x : radius(T_x, y) = y.height\}.$ 

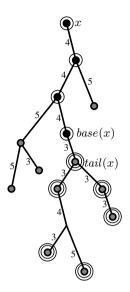
We extend this definition to a specific level  $\ell$  as follows:  $Spine(\ell) = \bigcup \{Spine(x) : x.level = \ell\}.$ 

We will denote by Others(x) the nodes in  $T_x$  that are not in Spine(x) (*i.e.*,  $Others(x) = T_x - Spine(x)$ ), and by  $Others(\ell) = \bigcup \{Others(x) : x.level = \ell\}$ . Furthermore, we define the *base process* of x, denoted by base(x), as the process in Spine(x) of greatest level; again, given a specific level  $\ell$ , we let  $Base(\ell) = \{base(x) : x.level = \ell\}$ . We define the *tail process* of x as  $tail(x) = best\_child$  (base(x)) (note that tail(x) might be not defined), and  $Tail(\ell) = \{tail(x) : x.level\} = \ell$ . Finally, we let  $Fan(x) = T_{tail(x)}$  and  $Fan(\ell) = \bigcup \{Fan(x) : x.level = \ell\}$ ; if tail(x) is undefined, we let  $Fan(x) = \emptyset$ . We now give few properties of Spine(x).

**Lemma 3.** For any process x

- (a)  $x \in Spine(x)$ .
- (b) If  $y \in Spine(x)$  and  $y \neq x$ , then  $p(y) \in Spine(x)$  and  $y = best\_child(p(y))$ .
- (c) Spine(x) is a chain.

For any  $s \in S$ , where S is connected, let  $longest\_path(S, s)$  denote the simple path of weight radius(S, s) in S starting at s. In the next lemma, we give a characterization of  $longest\_path(T_x, y)$ .



**Fig. 3.** Def. 1: Black (single circled) nodes are in Spine(x), while the light gray nodes are in Others(x); base(x), tail(x) are also shown; nodes in Fan(x) are double circled

**Lemma 4.** Let  $y \in T_x$ , and let u be the process of minimum level on longest\_path  $(T_x, y)$ . Then, the following properties hold: (a)  $u \in Spine(x)$ .

(b) If  $y \in Fan(x)$ , then longest\_path $(T_x, y) = path(y, u) +$ 

 $secondary_down_path(u)$ , where "+" denotes concatenation of paths.

(c) If  $y \notin Fan(x)$ , then  $longest\_path(T_x, y) = path(y, u) + longest\_path(T_u, u)$ ,

Let  $F_{\ell}$  be the forest given by the union of all  $T_x$ , where  $x.level = \ell$ . Thus,  $radius(F_{\ell}, y) = radius(T_x, y)$  if  $x.level = \ell$  and  $y \in T_x$ . The critical level of a process y enables y to determine whether it lies in  $Others(\ell)$  for any given  $\ell$ , as shown by the following lemma.

**Lemma 5.**  $y \in Spine(\ell)$  if and only if  $critical\_level(y) \le \ell \le y.level$ .

**Corollary 2.** Given any y such that  $\ell \leq y$ .level, the following properties hold: (a)  $y \in Others(\ell)$  if and only if critical\_level $(y) > \ell$ .

(b)  $y \in Spine(\ell)$  if and only if critical\_level $(y) \leq \ell$ .

(c)  $y \in Base(\ell)$  if and only if  $y \in Spine(\ell)$ , and either best\_child $(y) \in Others(\ell)$ , or best\_child(y) is undefined.

(d)  $y \in Tail(\ell)$  if and only if  $p(y) \in Base(\ell)$  and  $y = best\_child(p(y))$ .

Corollary 2 is used during the optimization phase of LINEAR<sub>max</sub> to determine the content of  $up\_package(y, \ell)$  (See Algorithms 1 and 6). In particular, the optimization phase proceeds bottom-up in the tree, with two nested loops. Let

$$local\_cost(y, \ell) = \min \{ w(y, y') + depth(y') : y' \in Swap\_N(y, \ell) \}$$

where  $Swap_N(y, \ell)$  is as defined in Section 5.1.

#### Algorithm 6. Algorithm LINEAR<sub>max</sub>

1:	Preprocessing phase
2:	Ranking phase
3:	Critical Level Phase (Algorithm 5)
4:	For all $1 \leq \ell \leq h$ Do $\%$ Optimization Phase $\%$
5:	For all y such that $y.level \ge \ell$ in bottom-up order <b>Do</b>
6:	$Swap_N(y, \ell) \leftarrow \{y' : \{y', y\} \text{ is a cross edge and } rank(y, y') < \ell\}$
7:	$local\_cost(y, \ell) = \min\left\{w(y, y') + depth(y') : y' \in Swap\_N(y, \ell)\right\}$
8:	If $y \in Others(\ell)$ Then
9:	If $y \in Others(\ell)$ Then $\min\_up\_cost(y, \ell) \leftarrow \min \begin{cases} local\_cost(y, \ell) \\ \min \{min\_up\_cost(z, \ell) + w(y, z) : z \in Chldrn(y)\} \end{cases}$ Else $\forall y \in Snine(\ell) \forall y$
5.	$\min \{\min_{y \in cost}(y, z) : z \in Chldrn(y)\}$
10	Else $\% y \in Spine(\ell)\%$ $min\_normal\_cost(y, \ell) \leftarrow \min \begin{cases} local\_cost(y, \ell) \\ min \{min\_up\_cost(z, \ell) + w(y, z) : z \in Normal\_Chldrn(y)\} \end{cases}$
11:	min normal cost(u, $\ell$ ) ( min $\int local\_cost(y, \ell)$
11.	$\min \{\min_{z \in \mathcal{O}} (z, \ell) \in \min \} \min \{\min_{z \in \mathcal{O}} (z, \ell) + w(y, z) : z \in Normal_Chldrn(y) \}$
12	If $best\_child(y)$ is defined Then
13	$z \leftarrow best\_child(y)$
14	If $z \in Spine(\ell)$ Then
15	$min\_fan\_cost(y,\ell) \leftarrow min\_fan\_cost(z,\ell) + w(z,y)$
16	Else
17	$min\_fan\_cost(y, \ell) \leftarrow min\_up\_cost(z, \ell) + w(z, y)$
	$sbtree\_min(y, \ell) \leftarrow \min \begin{cases} \min\_normal\_cost(y, \ell) + y.height\\ \min\_fan\_cost(y, \ell) + secondary\_height(y)\\ sbtree\_min(z, \ell) \end{cases}$
18:	$sbtree\_min(y, \ell) \leftarrow min \left\{ min\_fan\_cost(y, \ell) + secondary\_height(y) \right\}$
	$sbtree\_min(z,\ell)$
19	Else $\% y = base(x)$ , and $tail(x)$ undefined $\%$
20	$min\_fan\_cost(y, \ell) \leftarrow \infty$
21	$sbtree\_min(y, \ell) \leftarrow min\_normal\_cost(y, \ell) + y.height$
22	For all x such that $x.level = \ell$ Do
23	$M_F(x) = sbtree\_min(x, \ell)$

If  $y \in Others(\ell)$ , then  $radius(F_{\ell}, y)$  is not computed going down in the tree; hence, the only information that needs to be propagated (that is, the content of  $up\_package(y, \ell)$ ) is  $min\_up\_cost(y, \ell) = min \{local\_cost(z, \ell) + W(y, z) : z \in T_y\}$ , *i.e.*, the minimum value of W(path(y, z)) + w(z, z') + depth(z') over all  $z \in T_y$ such that  $\{z, z'\} \in SwapEdges(x)$ .

If  $y \in Spine(\ell)$ , first the value of  $min\_normal\_cost(y, \ell)$  is computed: It is equal to  $min\{local\_cost(z, \ell) + W(y, z)\}$ , such that  $z \in T_y$  and  $z \notin T_{best\_child(y)}$ . Then, the algorithm branches according to whether  $best\_child(y)$  is defined or not. If  $z = best\_child(y)$  is defined, then  $T_y \cap Fan(\ell) \neq \emptyset$ ; in this case we compute  $min\_fan\_cost(y, \ell) = min\{local\_cost(z, \ell) + W(y, z) : z \in T_y \cap Fan(\ell)\}$ , *i.e.*, the min value of W(path(y, z)) + w(z, z') + depth(z') over all  $\{z, z'\} \in SwapEdges(x)$ such that  $z \in T_y \cap Fan(\ell)$  (note that the algorithm computes  $min\_fan\_cost(y, \ell)$ differently, according to whether  $z \in Spine(\ell)$  or not). Finally, the actual cost of the swap edge is computed:  $sbtree\_min(y, \ell)$ , which is the minimum value of  $radius(T_y, z) + w(z, z') + depth(z')$  over all  $\{z, z'\} \in SwapEdges(x)$  such that  $z \in T_y$  (this is the value that is propagated in  $up\_package(y, \ell)$ ). By Lemma 4,  $sbtree\_min(y, \ell)$  is the minimum between the value of  $sbtree\_min(z, \ell)$ obtained from z,  $min\_normal\_cost(y, \ell) + y.height$ , and  $min\_fan\_cost(y, \ell) + secondary\_height(y)$ . If  $z = best\_child(y)$  is not defined, then  $T_y \cap Fan(\ell) = \emptyset$ . In this case,  $min\_fan\_cost(y, \ell)$  is set to  $\infty$ , and  $sbtree\_min(y, \ell)$  is set to  $min\_normal\_cost(y, \ell) + y.height$ .

When the  $\ell^{th}$  wave terminates, it is possible to compute the best swap edge for x:  $M_F(x) = sbtree\_min(x, \ell)$  for all x such that  $x.level = \ell$ . Again, as in the previous cases, the waves are executed in pipeline, and thus the overall time complexity is O(h). Also, in this case, it is not difficult to see that the number of messages used by Algorithm 6 is the same as for the quadratic time versions, *i.e.*,  $O(n^*)$ .

**Theorem 1.** The overall time complexity for LINEAR is O(h).

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