# The Tree Inclusion Problem: In Linear Space and Faster\*

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#### Abstract

Given two rooted, ordered, and labeled trees P and T the tree inclusion problem is to determine if P can be obtained from T by deleting nodes in T. This problem has recently been recognized as an important query primitive in XML databases. Kilpeläinen and Mannila [SIAM J. Comput. 1995] presented the first polynomial time algorithm using quadratic time and space. Since then several improved results have been obtained for special cases when P and T have a small number of leaves or small depth. However, in the worst case these algorithms still use quadratic time and space. Let  $n_S$ ,  $l_S$ , and  $d_S$  denote the number of nodes, the number of leaves, and the depth of a tree  $S \in \{P, T\}$ . In this paper we show that the tree inclusion problem can be solved in space  $O(n_T)$  and time:

$$O\left(\min\left\{\begin{array}{l} l_P n_T \\ l_P l_T \log \log n_T + n_T \\ \frac{n_P n_T}{\log n_T} + n_T \log n_T \end{array}\right\}\right)$$

This improves or matches the best known time complexities while using only linear space instead of quadratic. This is particularly important in practical applications, such as XML databases, where the space is likely to be a bottleneck.

# 1 Introduction

Let T be a rooted tree. We say that T is labeled if each node is assigned a character from an alphabet  $\Sigma$  and we say that T is ordered if a left-to-right order among siblings in T is given. All trees in this paper are rooted, ordered, and labeled. A tree P is included in T, denoted  $P \sqsubseteq T$ , if P can be obtained from T by deleting nodes of T. Deleting a node v in T means making the children of v children of the parent of v and then removing v. The children are inserted in the place of v in the left-to-right order among the siblings of v. The tree inclusion problem is to determine if P can be included in T and if so report all subtrees of T that include P.

Recently, the problem has been recognized as an important query primitive for XML data and has received considerable attention, see e.g., [33, 34, 37, 39–41]. The key idea is that an XML document can be viewed as a tree and queries on the document correspond to a tree inclusion problem. As an example consider Figure 1. Suppose that we want to maintain a catalog of books for a bookstore. A fragment of the tree, denoted D, corresponding to the catalog is shown in (b). In addition to supporting full-text queries, such as find all documents containing the word "John", we can also utilize the tree structure of the catalog to ask more specific queries, such as "find all books written by John with a chapter that has something to do with XML". We can model this query by constructing the tree, denoted Q, shown in (a) and solve the tree inclusion problem: is  $Q \sqsubseteq D$ ? The answer is yes and a possible way to include Q in D is indicated

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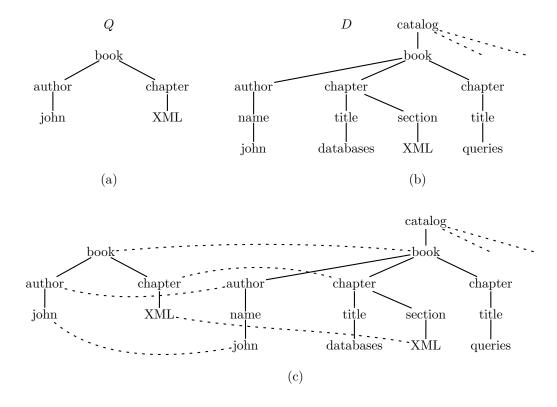


Figure 1: (a) The tree Q corresponding to the query. (b) A fragment of the tree D. Can the tree Q be included in the tree D? It can and an embedding is given in (c).

by the dashed lines in (c). If we delete all the nodes in D not touched by dashed lines the trees Q and D become isomorphic. Such a mapping of the nodes from Q to D given by the dashed lines is called an *embedding* (formally defined in Section 3). We note that the ordering of the XML document, and hence the left-to-right order of siblings, is important in many cases. For instance, in the above example, the relative order of contents of the chapters is most likely important. Also, in biological databases, order is of critical importance. Consequently, standard XML query languages, such as XPath [13] and XQuery [10], require the output of queries to be ordered.

The tree inclusion problem was initially introduced by Knuth [27, exercise 2.3.2-22] who gave a sufficient condition for testing inclusion. Motivated by applications in structured databases [24,29] Kilpeläinen and Mannila [25] presented the first polynomial time algorithm using  $O(n_P n_T)$  time and space, where  $n_P$  and  $n_T$  is the number of nodes in P and T, respectively. During the last decade several improvements of the original algorithm of [25] have been suggested [1,11,23,32]. The previously best known bound is due to Chen [11] who presented an algorithm using  $O(l_P n_T)$  time and  $O(l_P \cdot \min\{d_T, l_T\})$  space. Here,  $l_S$  and  $d_S$  denote the number of leaves and the depth of a tree S, respectively. This algorithm is based on an algorithm of Kilpeläinen [23]. Note that the time and space is still  $\Theta(n_P n_T)$  for worst-case input trees.

In this paper we present three algorithms which combined improve all of the previously known time and space bounds. To avoid trivial cases we always assume that  $1 \le n_P \le n_T$ . We show the following theorem:

**Theorem 1** For trees P and T the tree inclusion problem can be solved in  $O(n_T)$  space with the following

running time:

$$O\left(\min\left\{\begin{array}{l} l_P n_T \\ l_P l_T \log \log n_T + n_T \\ \frac{n_P n_T}{\log n_T} + n_T \log n_T \end{array}\right\}\right)$$

Hence, when P has few leaves we obtain a fast algorithm and even faster if both P and T have few leaves. When both trees have many leaves and  $n_P = \Omega(\log^2 n_T)$ , we instead improve the previous quadratic time bound by a logarithmic factor. Most importantly, the space used is linear. In the context of XML databases this will likely make it possible to query larger trees and speed up the query time since more of the computation can be kept in main memory.

The extended abstract of this paper [9] contained an error. The algorithms in the paper [9] did not use linear space. The problem was due to a recursive traversal of P which stored too many sets of nodes leading to a worst-case space complexity of  $\Omega(d_P l_T)$ . In this paper we fix this problem by recursively visiting the nodes such that the child with the largest number of descendant leaves is visited first, and by showing that the size of the resulting stored node sets exponentially decrease. With these ideas we show that all of our algorithms use  $O(n_T)$  space. Additionally, our improved analysis of the sizes of the stored node sets also leads to an improvement in the running time of the algorithm in the second case above. In the previous paper the running time was  $O(n_p l_T \log \log n_T + n_T)$ .

## 1.1 Techniques

Most of the previous algorithms, including the best one [11], are essentially based on a simple dynamic programming approach from the original algorithm of [25]. The main idea behind this algorithm is the following: Let v be a node in P with children  $v_1, \ldots, v_i$  and let w be a node in T. Consider the subtrees rooted at v and w, denoted by P(v) and T(w). To decide if P(v) can be included in T(w) we try to find a sequence  $w_1, \ldots, w_i$  of left-to-right ordered descendants of w such that  $P(v_k) \sqsubseteq T(w_k)$  for all  $k, 1 \le k \le i$ . The sequence is computed greedily from left-to-right in T(w) effectively finding the leftmost inclusion of P(v) in T(w). Applying this approach in a bottom-up fashion we can determine, if  $P(v) \sqsubseteq T(w)$ , for all pairs of nodes v in P and w in T.

In this paper we take a different approach. The main idea is to construct a data structure on T supporting a small number of procedures, called the *set procedures*, on subsets of nodes of T. We show that any such data structure implies an algorithm for the tree inclusion problem. We consider various implementations of this data structure which all use linear space. The first simple implementation gives an algorithm with  $O(l_P n_T)$  running time. As it turns out, the running time depends on a well-studied problem known as the *tree color problem*. We show a direct connection between a data structure for the tree color problem and the tree inclusion problem. Plugging in a data structure of Dietz [16] we obtain an algorithm with  $O(l_P l_T \log \log n_T + n_T)$  running time.

Based on the simple algorithms above we show how to improve the worst-case running time of the set procedures by a logarithmic factor. The general idea used to achieve this is to divide T into small trees called clusters of logarithmic size which overlap with other clusters in at most 2 nodes. Each cluster is represented by a constant number of nodes in a macro tree. The nodes in the macro tree are then connected according to the overlap of the cluster they represent. We show how to efficiently preprocess the clusters and the macro tree such that the set procedures use constant time for each cluster. Hence, the worst-case quadratic running time is improved by a logarithmic factor.

Our algorithms recursively traverse P top-down. For each node  $v \in V(P)$  we compute a set of nodes representing all of the subtrees in T that include P(v). To avoid storing too many of these node sets the traversal of P visits the child with the largest number of descendant leaves first. For the first two algorithms this immediately implies a space complexity of  $O(l_T \log l_P)$ , however, by carefully analyzing the sizes of stored node sets we are able to show that they decrease exponentially leading to the linear space bound. In the last algorithm the node sets are compactly encoded in  $O(n_T/\log n_T)$  space and therefore our recursive traversal alone implies a space bound of  $O(n_T/\log n_T \cdot \log l_P) = O(n_T)$ .

Throughout the paper we assume a unit-cost RAM model of computation with word size  $\Theta(\log n_T)$  and a standard instruction set including bitwise boolean operations, shifts, addition, and multiplication. All space complexities refer to the number of words used by the algorithm.

## 1.2 Related Work

For some applications considering unordered trees is more natural. However, in [25,30] this problem was proved to be NP-complete. The tree inclusion problem is closely related to the tree pattern matching problem [14,17,22,28]. The goal is here to find an injective mapping f from the nodes of P to the nodes of T such that for every node v in P the ith child of v is mapped to the ith child of f(v). The tree pattern matching problem can be solved in  $(n_P + n_T) \log^{O(1)}(n_P + n_T)$  time. Another similar problem is the subtree isomorphism problem [12,35], which is to determine if T has a subgraph isomorphic to P. The subtree isomorphism problem can be solved efficiently for ordered and unordered trees. The best algorithms for this problem use  $O(\frac{n_P^{1,5}n_T}{\log n_P} + n_T)$  time for unordered trees and  $O(\frac{n_Pn_T}{\log n_P} + n_T)$  time for ordered trees [12,35]. Both use  $O(n_Pn_T)$  space. The tree inclusion problem can be considered a special case of the tree edit distance problem [15,26,36,42]. Here one wants to find the minimum sequence of insert, delete, and relabel operations needed to transform P into T. Currently the best algorithm for this problem uses  $O(n_T n_P^2 (1 + \log \frac{n_T}{n_P}))$  time [15]. For more details and references see the survey [8].

## 1.3 Outline

In Section 2 we give notation and definitions used throughout the paper. In Section 3 a common framework for our tree inclusion algorithms is given. Section 4 presents two simple algorithms and then, based on these results, we show how to get a faster algorithm in Section 5.

## 2 Notation and Definitions

In this section we define the notation and definitions we will use throughout the paper. For a graph G we denote the set of nodes and edges by V(G) and E(G), respectively. Let T be a rooted tree. The root of T is denoted by root(T). The size of T, denoted by  $n_T$ , is |V(T)|. The depth of a node  $v \in V(T)$ , depth(v), is the number of edges on the path from v to root(T) and the depth of T, denoted  $d_T$ , is the maximum depth of any node in T. The parent of v is denoted parent(v) and the set of children of v is denoted child(v). We define parent $(root(T)) = \bot$ , where  $\bot \not\in V(T)$  is a special null node. A node with no children is a leaf and otherwise an internal node. The set of leaves of T is denoted L(T) and we define  $l_T = |L(T)|$ . We say that T is labeled if each node v is a assigned a character, denoted label(v), from an alphabet  $\Sigma$  and we say that T is ordered if a left-to-right order among siblings in T is given. Note that we do not require that the size of the alphabet is bounded by a constant. All trees in this paper are rooted, ordered, and labeled.

Ancestors and Descendants Let T(v) denote the subtree of T rooted at a node  $v \in V(T)$ . If  $w \in V(T(v))$  then v is an ancestor of w, denoted  $v \preceq w$ , and if  $w \in V(T(v)) \setminus \{v\}$  then v is a proper ancestor of w, denoted  $v \prec w$ . If v is a (proper) ancestor of w then w is a (proper) descendant of v. A node z is a common ancestor of v and w if it is an ancestor of both v and w. The nearest common ancestor of v and v, ncav, is the common ancestor of v and v of greatest depth. The first ancestor of v labeled v denoted flv, v is the node v such that  $v \preceq w$ , label(v) = v, and no node on the path between v and v is labeled v. If no such node exists then flv, v is the flv is a proper ancestor of v and v is labeled v. If no such node exists then flv, v is the flv in the flv in the flv in the flv in the flv is a proper ancestor of v and v is labeled v. If no such node exists then flv in the flv

**Traversals and Orderings** Let T be a tree with root v and let  $v_1, \ldots, v_k$  be the children of v from left-to-right. The *preorder traversal* of T is obtained by visiting v and then recursively visiting  $T(v_i)$ ,  $1 \le i \le k$ , in order. Similarly, the *postorder traversal* is obtained by first visiting  $T(v_i)$ ,  $1 \le i \le k$ , in order and then v. The *preorder number* and *postorder number* of a node  $w \in T(v)$ , denoted by  $\operatorname{pre}(w)$  and  $\operatorname{post}(w)$ , are the

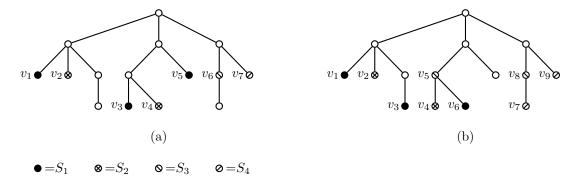


Figure 2: In (a) we have  $\{(v_1, v_2, v_3, v_6, v_7), (v_1, v_2, v_5, v_6, v_7), (v_1, v_4, v_5, v_6, v_7), (v_3, v_4, v_5, v_6, v_7)\}$  =  $\Phi(S_1, S_2, S_1, S_3, S_4)$  and thus  $\text{mop}(S_1, S_2, S_1, S_3, S_4) = \{(v_3, v_7)\}$ . In (b) we have  $\Phi(S_1, S_2, S_1, S_3, S_4) = \{(v_1, v_2, v_3, v_5, v_7), (v_1, v_2, v_6, v_8, v_9), (v_1, v_2, v_3, v_8, v_9), (v_1, v_2, v_3, v_5, v_9), (v_1, v_4, v_6, v_8, v_9), (v_3, v_4, v_6, v_8, v_9)\}$  and thus  $\text{mop}(S_1, S_2, S_1, S_3, S_4) = \{(v_1, v_7), (v_3, v_9)\}$ .

number of nodes preceding w in the preorder and postorder traversal of T, respectively. The nodes to the left of w in T is the set of nodes  $u \in V(T)$  such that  $\operatorname{pre}(u) < \operatorname{pre}(w)$  and  $\operatorname{post}(u) < \operatorname{post}(w)$ . If u is to the left of w, denoted by  $u \triangleleft w$ , then w is to the right of u. If  $u \triangleleft w$  or  $u \preceq w$  or  $w \prec u$  we write  $u \unlhd w$ . The null node  $\bot$  is not in the ordering, i.e.,  $\bot \not \lhd v$  for all nodes v.

Minimum Ordered Pairs A set of nodes  $X \subseteq V(T)$  is deep if no node in X is a proper ancestor of another node in X. For k deep sets of nodes  $X_1, \ldots, X_k$  let  $\Phi(X_1, \ldots, X_k) \subseteq (X_1 \times \cdots \times X_k)$ , be the set of tuples such that  $(x_1, \ldots, x_k) \in \Phi(X_1, \ldots, X_k)$  iff  $x_1 \lhd \cdots \lhd x_k$ . If  $(x_1, \ldots, x_k) \in \Phi(X_1, \ldots, X_k)$  and there is no  $(x'_1, \ldots, x'_k) \in \Phi(X_1, \ldots, X_k)$ , where either  $x_1 \lhd x'_1 \lhd x'_k \unlhd x_k$  or  $x_1 \unlhd x'_1 \lhd x'_k \lhd x_k$  then the pair  $(x_1, x_k)$  is a minimum ordered pair. Intuitively,  $(x_1, x_k)$  is a closest pair of nodes from  $X_1$  and  $X_k$  in the left-to-right order for which we can find  $x_2, \ldots, x_{k-1}$  such that  $x_1 \lhd \cdots \lhd x_k$ . The set of minimum ordered pairs for  $X_1, \ldots, X_k$  is denoted by  $mop(X_1, \ldots, X_k)$ . Figure 2 illustrates these concepts on a small example. For any set of pairs Y, let  $Y|_1$  and  $Y|_2$  denote the projection of Y to the first and second coordinate, that is, if  $(y_1, y_2) \in Y$  then  $y_1 \in Y|_1$  and  $y_2 \in Y|_2$ . We say that Y is deep if  $Y|_1$  and  $Y|_2$  are deep. The following lemma shows that given deep sets  $X_1, \ldots, X_k$  we can compute  $mop(X_1, \ldots, X_k)$  iteratively by first computing  $mop(X_1, X_2)$  and then  $mop(mop(X_1, X_2)|_2, X_3)$  and so on.

**Lemma 1** For any deep sets of nodes  $X_1, \ldots, X_k$ , k > 2, we have,  $(x_1, x_k) \in \text{mop}(X_1, \ldots, X_k)$  iff there exists a node  $x_{k-1}$  such that  $(x_1, x_{k-1}) \in \text{mop}(X_1, \ldots, X_{k-1})$  and  $(x_{k-1}, x_k) \in \text{mop}(\text{mop}(X_1, \ldots, X_{k-1})|_2, X_k)$ .

*Proof.* We start by showing that if  $(x_1, x_k) \in \text{mop}(X_1, \dots, X_k)$  then there exists a node  $x_{k-1}$  such that  $(x_1, x_{k-1}) \in \text{mop}(X_1, \dots, X_{k-1})$  and  $(x_{k-1}, x_k) \in \text{mop}(\text{mop}(X_1, \dots, X_{k-1})|_2, X_k)$ .

First note that  $(z_1,\ldots,z_k)\in\Phi(X_1,\ldots,X_k)$  implies  $(z_1,\ldots,z_{k-1})\in\Phi(X_1,\ldots,X_{k-1})$ . Since  $(x_1,x_k)\in \operatorname{mop}(X_1,\ldots,X_k)$  there must be a minimum node  $x_{k-1}$  such that the tuple  $(x_1,\ldots,x_{k-1})$  is in  $\Phi(X_1,\ldots,X_{k-1})$ . We have  $(x_1,x_{k-1})\in\operatorname{mop}(X_1,\ldots,X_{k-1})$ . We need to show that  $(x_{k-1},x_k)\in\operatorname{mop}(\operatorname{mop}(X_1,\ldots,X_{k-1})|_2,X_k)$ . Since  $(x_1,x_k)\in\operatorname{mop}(X_1,\ldots,X_k)$  there exists no node  $z\in X_k$  such that  $x_{k-1}\lhd z\lhd x_k$ . If such a z existed we would have  $(x_1,\ldots,x_{k-1},z)\in\Phi(X_1,\ldots,X_k)$ , contradicting that  $(x_1,x_k)\in\operatorname{mop}(X_1,\ldots,X_k)$ . Assume there exists a node  $z\in\operatorname{mop}(X_1,\ldots,X_{k-1})|_2$  such that  $x_{k-1}\lhd z\lhd x_k$ . Since  $(x_1,x_{k-1})\in\operatorname{mop}(X_1,\ldots,X_{k-1})$  this implies that there is a node  $z'\rhd x_1$  such that  $(z',z)\in\operatorname{mop}(X_1,\ldots,X_{k-1})$ . But this implies that the tuple  $(z',\ldots,z,x_k)$  is in  $\Phi(X_1,\ldots,X_k)$  contradicting that  $(x_1,x_k)\in\operatorname{mop}(X_1,\ldots,X_k)$ .

We will now show that if there exists a node  $x_{k-1}$  such that  $(x_1, x_{k-1}) \in \text{mop}(X_1, \dots, X_{k-1})$  and  $(x_{k-1}, x_k) \in \text{mop}(\text{mop}(X_1, \dots, X_{k-1})|_2, X_k)$  then the pair  $(x_1, x_k) \in \text{mop}(X_1, \dots, X_k)$ . Clearly, there exists a tuple  $(x_1, \dots, x_{k-1}, x_k) \in \Phi(X_1, \dots, X_k)$ . Assume that there exists a tuple  $(z_1, \dots, z_k) \in \Phi(X_1, \dots, X_k)$  such that  $x_1 \triangleleft z_1 \triangleleft z_k \unlhd x_k$ . Among the tuples satisfying these constraints let  $(y_1, \dots, y_{k-1}, y_k)$  be the one with maximum  $y_1$ , minimum  $y_{k-1}$ , and maximum  $y_k$ . It follows that  $(y_1, y_{k-1}) \in \text{mop}(X_1, \dots, X_{k-1})$ .

Since  $(x_1, x_{k-1}) \in \text{mop}(X_1, ..., X_{k-1})$  we must have  $x_{k-1} \triangleleft y_{k-1}$ . But this contradicts  $(x_{k-1}, x_k) \in \text{mop}(\text{mop}(X_1, ..., X_{k-1})|_2, X_k)$ , since node  $y_{k-1} \in \text{mop}(X_1, ..., X_{k-1})|_2$ .

Assume that there exists a tuple  $(z_1, \ldots, z_k) \in \Phi(X_1, \ldots, X_k)$  such that  $x_1 \leq z_1 < z_k < x_k$ . Since  $(x_1, x_{k-1}) \in \text{mop}(X_1, \ldots, X_{k-1})$  we have  $x_{k-1} \leq z_{k-1}$  and thus  $x_{k-1} < z_k < x_k$  contradicting  $(x_{k-1}, x_k) \in \text{mop}(\text{mop}(X_1, \ldots, X_{k-1})|_2, X_k)$ .

The following lemma is the reverse of the previous lemma and shows that given deep sets  $X_1, \ldots, X_k$  we also can compute  $mop(X_1, \ldots, X_k)$  iteratively from right to left. The proof is similar to the proof of Lemma 1.

**Lemma 2** For any deep sets of nodes  $X_1, \ldots, X_k$ , k > 2, we have,  $(x_1, x_k) \in \text{mop}(X_1, \ldots, X_k)$  iff there exists a node  $x_2$  such that  $(x_2, x_k) \in \text{mop}(X_2, \ldots, X_k)$  and  $(x_1, x_2) \in \text{mop}(X_1, \text{mop}(X_2, \ldots, X_k)|_1)$ .

**Heavy Leaf Path Decomposition** We construct a heavy leaf path decomposition of P as follows. We classify each node as heavy or light. The root is light. For each internal node v we pick a child  $v_j$  of v with maximum  $l_{P(v_j)}$  and classify it as heavy. The remaining children of v are light. An edge to a light node is a light edge, and an edge to a heavy node is a heavy edge. The heavy child of a node v is denoted heavy (v). Let ldepth(v) denote the number of light edges on the path from v to root(P).

Note that a heavy leaf path decomposition is the same as the classical heavy path decomposition [21] except that the heavy child is defined as the child with largest number of the descendant leaves and not the child with the largest number of descendants. This distinction is essential for achieving the linear space bound of our algorithms. Note that heavy path decompositions have previously been used in algorithms for the related tree edit distance problem [26].

**Lemma 3** For any tree P and node  $v \in V(P)$ ,

$$l_{P(v)} \le \frac{l_P}{2^{\operatorname{ldepth}(v)}}$$
.

*Proof.* By induction on ldepth(v). For ldepth(v) = 0 it is trivially true. Let  $ldepth(v) = \ell$ . Assume wlog. that v is light. Let w be the unique light ancestor of v with  $ldepth(w) = \ell - 1$ . By the induction hypothesis  $l_{P(w)} \leq l_P/2^{\ell-1}$ . Now v has a sibling heavy(parent(v)) and thus at most half of the leaves in P(parent(v)) can be in the subtree rooted at v. Therefore,  $l_{P(v)} \leq l_{P(w)}/2 \leq l_P/2^{\ell}$ .

Corollary 1 For any tree P and node  $v \in V(P)$ ,  $ldepth(v) < log l_P$ .

**Notation** When we want to specify which tree we mean in the above relations we add a subscript. For instance,  $v \prec_T w$  indicates that v is an ancestor of w in T.

# 3 Computing Deep Embeddings

In this section we present a general framework for answering tree inclusion queries. As in [25] we solve the equivalent tree embedding problem. Let P and T be rooted labeled trees. An embedding of P in T is an injective function  $f: V(P) \to V(T)$  such that for all nodes  $v, u \in V(P)$ ,

- (i) label(v) = label(f(v)). (label preservation condition)
- (ii)  $v \prec u$  iff  $f(v) \prec f(u)$ . (ancestor condition)
- (iii)  $v \triangleleft u$  iff  $f(v) \triangleleft f(u)$ . (order condition)

An example of an embedding is given in Figure 1(c).

**Lemma 4 (Kilpeläinen and Mannila [25])** For any trees P and T,  $P \sqsubseteq T$  iff there exists an embedding of P in T.

We say that the embedding f is deep if there is no embedding g such that  $f(\text{root}(P)) \prec g(\text{root}(P))$ . The deep occurrences of P in T, denoted emb(P,T) is the set of nodes,

$$emb(P,T) = \{ f(root(P)) \mid f \text{ is a deep embedding of } P \text{ in } T \}.$$

By definition the set of ancestors of nodes in  $\operatorname{emb}(P,T)$  is exactly the set of nodes  $\{u \mid P \sqsubseteq T(u)\}$ . Hence, to solve the tree inclusion problem it is sufficient to compute  $\operatorname{emb}(P,T)$  and then, using additional  $O(n_T)$  time, report all ancestors of this set. We note that Kilpeläinen and Mannila [25] used the similar concept of *left embeddings* in their algorithms. A left embedding of P in T is an embedding such that the root of P is mapped to the node in P with the smallest postorder number, i.e., the deepest node among the nodes furthest to the left. Our definition of  $\operatorname{emb}(P,T)$  only requires that the root is mapped to a deepest node.

In the following we show how to compute deep embeddings. The key idea is to build a data structure for T allowing a fast implementation of the following procedures. For all  $X \subseteq V(T)$ ,  $Y \subseteq V(T) \times V(T)$ , and  $\alpha \in \Sigma$  define:

PARENT(X): Return the set  $\{parent(x) \mid x \in X\}$ .

NCA(Y): Return the set  $\{ nca(y_1, y_2) \mid (y_1, y_2) \in Y \}$ .

DEEP(X): Return the set  $\{x \in X \mid \text{there is no } z \in X \text{ such that } x \prec z\}.$ 

MOPRIGHT(Y, X): Return the set of pairs R such that for any pair  $(y_1, y_2) \in Y$ ,  $(y_1, x) \in R$  iff  $(y_2, x) \in \text{mop}(Y|_2, X)$ .

MOPLEFT(X,Y): Return the set of pairs R such that for any pair  $(y_1,y_2) \in Y$ ,  $(x,y_2) \in R$  iff  $(x,y_1) \in \text{mop}(X,Y|_1)$ .

 $FL(X, \alpha)$ : Return the set  $DEEP(\{fl(x, \alpha) \mid x \in X\})$ .

Collectively we call these procedures the set procedures. The procedures Parent and NCA are self-explanatory. Deep(X) returns the set of all nodes in X that have no descendants in X. Hence, the returned set is always deep. MopRight and MopLeft are used to iteratively compute minimum ordered pairs.  $FL(X, \alpha)$  returns the deep set of first ancestors with label  $\alpha$  of all nodes in X. If we want to specify that a procedure applies to a certain tree T we add the subscript T. With the set procedures we can compute deep embeddings. The following procedure EMB(v),  $v \in V(P)$ , recursively computes the set of deep occurrences of P(v) in T. Figure 3 illustrates how EMB works on a small example.

## Procedure Emb(v)

```
1 Let v_1, \ldots, v_k be the sequence of children of v ordered from left to right. There are three cases:
 2 case 1. k = 0 // v is a leaf
    Compute R := FL(L(T), label(v)).
 4 case 2. k = 1
       Recursively compute R_1 := EMB(v_1).
 5
       Compute R := FL(DEEP(PARENT(R_1)), label(v)).
 7 case 3. k > 1
 8
       Let v_i be the heavy child of v.
       Recursively compute R_i := \text{Emb}(v_i) and set U_i := \{(r, r) \mid r \in R_i\}.
 9
       for i := j + 1 to k do
10
          Recursively compute R_i := \text{EmB}(v_i) and set U_i := \text{MOPRIGHT}(U_{i-1}, R_i).
11
       end
12
       Set U_i := U_k.
13
       for i := j - 1 downto 1 do
14
        Recursively compute R_i := \text{Emb}(v_i) and set U_i := \text{MopLeft}(R_i, U_{i+1}).
15
       end
16
       Compute R := FL(DEEP(NCA(U_1)), label(v)).
17
18
19 if R = \emptyset then
      stop and report that there is no deep embedding of P(v) in T.
21 else
       Return R.
22
\mathbf{23}
```

To prove the correctness of the EMB procedure we need the following two propositions. The first proposition characterizes for node  $v \in V(P)$  the set  $\mathrm{emb}(v,T)$  using mop, nca, and fl. The second proposition shows that the set  $U_1$  computed in case 3 of the EMB procedure is the set  $\mathrm{mop}(\mathrm{EMB}(v_1),\ldots,\mathrm{EMB}(v_k))$ .

**Proposition 1** Let v be a node in P and let  $v_1, \ldots, v_k$  be the sequence of children of v ordered from left to right, where  $k \geq 2$ . For any node  $w \in \text{emb}(P(v), T)$ , there exists a pair of nodes  $(w_1, w_k) \in \text{mop}(\text{emb}(P(v_1), T), \ldots, \text{emb}(P(v_k), T))$  such that  $w = \text{fl}(\text{nca}(w_1, w_k), \text{label}(v))$ .

*Proof.* Since w is the root of an occurrence of P(v) in T there must exist a set of disjoint occurrences of  $P(v_1), \ldots, P(v_k)$  in T(w) with roots  $w_1 \triangleleft \ldots \triangleleft w_k$ , such that w is an ancestor of  $w_1, \ldots, w_k$ . Since the  $w_i$ 's are ordered w must be an ancestor of  $\operatorname{nca}(w_1, w_k)$ . Since w is the root of a *deep* occurrence of P(v) in T it follows that  $w = \operatorname{fl}(\operatorname{nca}(w_1, w_k), \operatorname{label}(v))$ .

It remains to show that we can assume  $(w_1, w_k) \in \text{mop}(\text{emb}(P(v_1), T), \dots, \text{emb}(P(v_k), T))$ . It follows from the previous discussion that  $(w_1, \dots, w_k) \in \Phi(\text{emb}(P(v_1), T), \dots, \text{emb}(P(v_k), T))$ . Assume for the sake of contradiction that  $(w_1, w_k)$  is not a minimum ordered pair. Then there exists a set of disjoint occurrences of  $P(v_1), \dots, P(v_k)$  in T(w) with roots  $u_1 \triangleleft \dots \triangleleft u_k$ , such that either  $w_1 \triangleleft u_1 \triangleleft u_k \triangleleft w_k$  or  $w_1 \triangleleft u_1 \triangleleft u_k \triangleleft w_k$ , and  $(u_1, u_k) \in \text{mop}(\text{emb}(P(v_1), T), \dots, \text{emb}(P(v_k), T))$ . Therefore  $u = \text{fl}(\text{nca}(u_1, u_k), \text{label}(v))$  is an embedding of P(v) in T. Now either  $w \triangleleft u$  contradicting the assumption that w is a deep embedding or w = u in which case  $(u_1, u_k)$  satisfies the properties of the lemma (see also Figure 4(a)).

**Proposition 2** For  $j + 1 \le l \le k$ ,

$$U_l = \text{mop}(\text{EmB}(v_i), \dots, \text{EmB}(v_l)), \tag{1}$$

For 
$$1 \le l \le j - 1$$
,
$$U_l = \operatorname{mop}(\operatorname{EmB}(v_l), \dots, \operatorname{EmB}(v_k)). \tag{2}$$

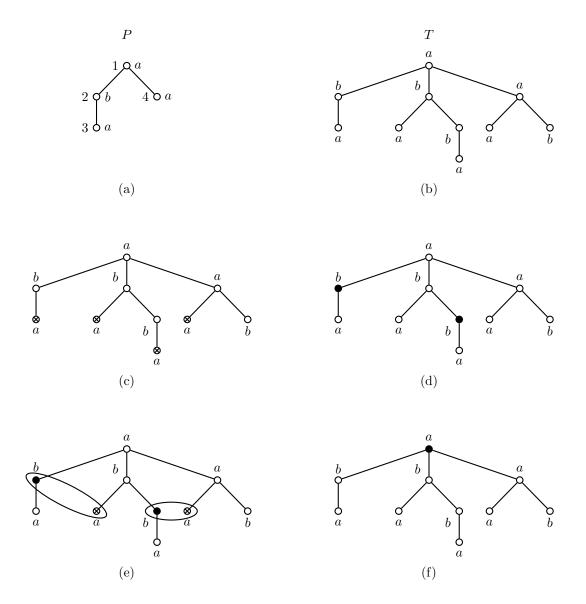


Figure 3: Computing the deep occurrences of P into T depicted in (a) and (b) respectively. The nodes in P are numbered 1–4 for easy reference. (c) Case 1 of EMB: The crossed nodes are the nodes in the set EMB(3). Since 3 and 4 are leaves and label(3) = label(4) we have EMB(3) = EMB(4). (d) Case 2 of EMB: The black nodes are the nodes in the set EMB(2). Note that the middle child of the root of T is not in the set since it is not a deep occurrence. (e) and (f) illustrates the computation of EMB(1) and case 3 of EMB: (e) The two minimal ordered pairs of the sets from (d) and (c). In the procedure  $R_1$  is the set from (d) and  $R_2$  is the set from (c). The set  $U_1 = \{(v,v) \mid v \in R_1\}$  and the set  $U_2 = \text{MopRight}(U_1, R_2)$  which corresponds to the pairs shown in (e). The black nodes in the pairs are the nodes from  $R_1$  and the crossed nodes are the nodes from  $R_2$ . Since k = 2 we set  $U_1 = U_2$ . (f) The nearest common ancestors of both pairs shown in (e) is the root node of T which is the only (deep) occurrence of P.

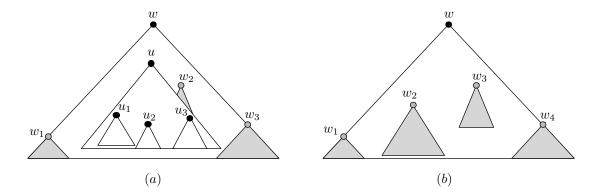


Figure 4: (a) For all i,  $w_i$  and  $u_i$  are roots of occurrences of  $P(v_i)$  in T, and w and u is the nearest common ancestor of  $(w_1, w_3)$  and  $(u_1, u_3)$ , respectively. Since  $w_1 \triangleleft u_1$  and  $u_3 \triangleleft w_3$  we cannot have  $u \triangleleft w$ . (b) For all i,  $w_i$  is an embedding of  $P(v_i)$  in T,  $(w_1, w_4)$  is a minimum ordered pair, and w is the nearest common ancestor of all the  $w_i$ 's. The number of leaves in T(w) is at least  $\sum_{i=1}^4 l_{T(w_i)} \ge \sum_{i=1}^4 l_{P(v_i)}$ .

*Proof.* We first show Equation (1) by induction on l. For l = j+1 it follows from the definition of MopRight that  $U_l$  is the set of minimum ordered pairs of  $\text{EMB}(v_j)$  and  $\text{EMB}(v_{j+1})$ , i.e.,  $U_l = \text{mop}(\text{EMB}(v_j), \text{EMB}(v_l))$ . Hence, assume that l > j+1. By the induction hypothesis we have

$$U_l = \text{MopRight}(U_{l-1}, \text{EmB}(v_l)) = \text{MopRight}(\text{mop}(\text{EmB}(v_i), \dots, \text{EmB}(v_{l-1})), R_l)$$
.

By definition of Mopricht,  $U_l$  is the set of pairs such that for any pair  $(r_j, r_{l-1}) \in \text{mop}(\text{EmB}(v_j), \dots, \text{EmB}(v_{l-1})), (r_j, r_l) \in U_l$  iff  $(r_{l-1}, r_l) \in \text{mop}(\text{mop}(\text{EmB}(v_j), \dots, \text{EmB}(v_{l-1}))|_2, R_l)$ . By Lemma 1 it follows that  $(r_j, r_l) \in U_l$  iff  $(r_j, r_l) \in \text{mop}(\text{EmB}(v_j), \dots, \text{EmB}(v_l))$ .

We can now similarly show Equation (2) by induction on j' = j - l. By Equation (1) we have  $U_j = \text{mop}(\text{EMB}(v_j), \dots, \text{EMB}(v_k))$  when we begin computing  $U_{j-1}$ . For j' = 1 (l = j - 1) it follows from the definition of Mopleft that  $U_{j-1} = \text{mop}(\text{EMB}(v_{j-1}), \text{EMB}(v_j))$ . Hence, assume that j' > 1. Using Lemma 2 the Equation follows similarly to the proof of Equation (1).

By Proposition 2,  $U_1 = \text{mop}(\text{EmB}(v_1), \dots, \text{EmB}(v_k))$ . We can now show the correctness of procedure Emb.

**Lemma 5** For trees P and T and node  $v \in V(P)$ , Emb(v) computes the set of deep occurrences of P(v) in T.

*Proof.* By induction on the size of the subtree P(v). If v is a leaf, emb(v,T) is the deep set of nodes in T with label label(v). It immediately follows that emb(v,T) = FL(L(T), label(v)) and thus case 1 follows.

Suppose that v is an internal node with  $k \geq 1$  children  $v_1, \ldots, v_k$ . We show that emb(P(v), T) = Emb(v). Consider cases 2 and 3 of the algorithm.

For k = 1 we have that  $w \in EmB(v)$  implies that label(w) = label(v) and there is a node  $w_1 \in EmB(v_1)$  such that  $f(parent(w_1), label(v)) = w$ , that is, no node on the path between  $w_1$  and w is labeled label(v). By induction  $EmB(v_1) = emb(P(v_1), T)$  and therefore w is the root of an embedding of P(v) in T. Since EmB(v) is the deep set of all such nodes it follows that  $w \in emb(P(v), T)$ . Conversely, if  $w \in emb(P(v), T)$  then label(w) = label(v), there is a node  $w_1 \in emb(P(v_1), T)$  such that  $w \prec w_1$ , and no node on the path between w and  $w_1$  is labeled label(v), that is,  $f(w_1, label(v)) = w$ . Hence,  $w \in EmB(v)$ .

Next consider the case k > 1. By Proposition 2 and the induction hypothesis

$$U_1 = \operatorname{mop}(\operatorname{emb}(P(v_1), T), \dots, \operatorname{emb}(P(v_k), T))$$
.

We first show that  $w \in \text{emb}(P(v), T)$  implies that  $w \in \text{EMB}(v)$ . By Proposition 1 there exists a pair of nodes  $(w_1, w_k) \in \text{mop}(\text{emb}(P(v_1), T), \dots, \text{emb}(P(v_k), T))$  such that  $w = \text{fl}(\text{nca}(w_1, w_k), \text{label}(v))$ . We have

 $(w_1, w_k) \in U_1$  and it follows directly from the implementation that  $w \in \text{EMB}(v)$ . To see that we do not loose w by taking DEEP of NCA $(U_1)$  assume that  $w' = \text{nca}(w_1, w_k)$  is removed from the set in this step. This means there is a node u in NCA $(U_1)$  which is a descendant of w' and which is still in the set. Since w is the root of a deep occurrence we must have w = fl(w', label(v)) = fl(u, label(v)).

Let  $w \in \text{EMB}(v)$ . Then w is the first ancestor with label label (v) of a nearest common ancestor of a pair in  $U_1$ . That is, label (w) = label(v) and there exists nodes  $(w_1, w_k) \in \text{mop}(\text{emb}(P(v_1), T), \ldots, \text{emb}(P(v_k), T))$  such that  $w = \text{fl}(\text{nca}(w_1, w_k), \text{label}(v))$ . Clearly, w is the root of an embedding of P(v) in T. Assume for contradiction that w is not a deep embedding, that is,  $w \prec u$  for some node  $u \in \text{emb}(P(v), T)$ . We have just shown that this implies  $u \in \text{EmB}(v)$ . Since EmB(v) is a deep set this contradicts  $w \in \text{EmB}(v)$ .

The set L(T) is deep and in all three cases of Emb(V) the returned set is also deep. By induction it follows that the input to Parent, Fl, Nca, and Mopright is always deep. We will use this fact to our advantage in the following algorithms.

## 4 A Simple Tree Inclusion Algorithm

In this section we a present a simple implementation of the set procedures which leads to an efficient tree inclusion algorithm. Subsequently, we modify one of the procedures to obtain a family of tree inclusion algorithms where the complexities depend on the solution to a well-studied problem known as the *tree color problem*.

## 4.1 Preprocessing

To compute deep embeddings we require a data structure for T which allows us, for any  $v, w \in V(T)$ , to compute  $\text{nca}_T(v, w)$  and determine if  $v \prec w$  or  $v \triangleleft w$ . In linear time we can compute pre(v) and post(v) for all nodes  $v \in V(T)$ , and with these it is straightforward to test the two conditions. Furthermore,

**Lemma 6 (Harel and Tarjan [21])** For any tree T there is a data structure using  $O(n_T)$  space and preprocessing time which supports nearest common ancestor queries in O(1) time.

Hence, our data structure uses linear preprocessing time and space (see also [2,7] for more recent nearest common ancestor data structures).

### 4.2 Implementation of the Set Procedures

To answer tree inclusion queries we give an efficient implementation of the set procedures. The idea is to represent sets of nodes and sets of pairs of nodes in a left-to-right order using linked lists. For this purpose we introduce some helpful notation. Let  $X = [x_1, \ldots, x_k]$  be a linked list of nodes. The length of X, denoted |X|, is the number of elements in X and the list with no elements is written []. The ith node of X, denoted X[i], is  $x_i$ . Given any node y the list obtained by appending y to X, is the list  $X \circ y = [x_1, \ldots, x_k, y]$ . If for all i,  $1 \le i \le |X| - 1$ ,  $X[i] \triangleleft X[i+1]$  then X is ordered and if  $X[i] \unlhd X[i+1]$  then X is semiordered. Recall that  $X[i] \unlhd X[i+1]$  means that we can have  $X[i] \triangleleft X[i+1]$  or either of the nodes can be an ancestor of the other  $(X[i] \triangleleft X[i+1])$  or  $X[i] \succeq X[i+1]$  or  $X[i] \succeq X[i+1]$ ). A list  $Y = [(x_1, z_k), \ldots, (x_k, z_k)]$  is a node pair list. By analogy, we define length, append, etc. for Y. For a pair  $Y[i] = (x_i, z_i)$  define  $Y[i]_1 = x_i$  and  $Y[i]_2 = z_i$ . If the lists  $[Y[1]_1, \ldots, Y[k]_1]$  and  $[Y[1]_2, \ldots, Y[k]_2]$  are both ordered or semiordered then Y is ordered or semiordered, respectively.

The set procedures are implemented using node lists. All lists used in the procedures are either ordered or semiordered. As noted in Section 3 we may assume that the inputs to all of the procedures, except DEEP, represent deep sets, that is, the corresponding node list or node pair list is ordered. We assume that the input list given to DEEP is semiordered and the output, of course, is ordered. Hence, the output of all the other set procedures must be semiordered. In the following let X be a node list, Y a node pair list, and  $\alpha$  a character in  $\Sigma$ . The detailed implementation of the set procedures is given below. We show the correctness in Section 4.3 and discuss the complexity in Section 4.4.

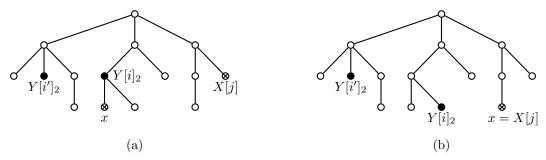


Figure 5: Case 1 and 2 from the implementation of MOPRIGHT. (a) We have  $x \triangleleft X[j]$  and therefore  $Y[i]_2 \not \triangleleft x$ . So (y,x) is inserted in R. (b) We have  $Y[i']_2 \triangleleft Y[i]_2 \triangleleft x = X[j]$ .

## Procedure Parent(X)

1 Return the list  $[parent(X[1]), \ldots, parent(X[|X|])]$ 

## Procedure NCA(Y)

1 Return the list [nca(Y[1]), ..., nca(Y[|Y|])].

#### Procedure DEEP(X)

```
1 Initially, set x := X[1] and R := [].
 2 for i := 2 to |X| do
       Compare x and X[i]. There are three cases:
 3
 4
       case 1. x \triangleleft X[i]
          Set R := R \circ x and x := X[i].
 5
       case 2. x \prec X[i]
 6
           Set x := X[i].
 7
       case 3. X[i] \leq x
 8
 9
           Do nothing.
10
11 end
12 Return R \circ x.
```

The implementation of procedure DEEP takes advantage of the fact that the input list is semiordered. In case 1 the node X[i] is to the right of our "potential output node" x. Since any node that is a descendant of x must also be to the left of X[i] it cannot appear later in the list X than X[i]. We can thus safely add x to R at this point. In case 2 the node x is an ancestor of X[i] and thus x cannot be in DEEP(X). In case 3 the node X[i] is an ancestor of x and can therefore not be in DEEP(X).

In procedure Moprich we have a "potential pair" (y,x) where  $y=Y[i']_1$  for some i' and  $Y[i']_2 \triangleleft x$ . In case 1 we have  $x \triangleleft X[j]$  and also  $Y[i']_2 \triangleleft Y[i]_2$  since the input lists are ordered and i' < i (see Figure 5(a)). Therefore, (y,x) is inserted into R. In case 2 we have x=X[j], i.e.,  $Y[i]_2 \triangleleft x$ , and as before  $Y[i']_2 \triangleleft Y[i]_2$  (see Figure 5(b)). Therefore (y,x) cannot be in Moprich (y,x), and we set  $(Y[i]_1,x)$  to be the new potential pair.

We can implement MopLeft(X, Y) similarly to MopRight replacing smallest by largest,  $\triangleleft$  by  $\triangleright$ , and traversing the lists backwards:

### **Procedure** MopRight(Y,X)

```
1 Initially, set R := [].
 2 Find the smallest j such that Y[1]_2 \triangleleft X[j] and set y := Y[1]_1, x := X[j]. If no such j exists stop and
   return R.
 3 for i := 2 to |Y| do
       until Y[i]_2 \triangleleft X[j] or j > |X| do
 4
           set j := j + 1.
       if j > |X| then
 6
           stop and return R := R \circ (y, x).
 7
       else
 8
           Compare X[i] and x. There are two cases:
 9
           case 1. x \triangleleft X[j]
10
            set R := R \circ (y, x), y := Y[i]_1, and x := X[j].
11
           case 2. If x = X[j]
12
            set y := Y[i]_1.
13
14
15
16 end
17 Return R := R \circ (y, x).
```

#### Procedure $FL(X,\alpha)$

```
1 Initially, set L := X, Z := L.
 2 while Z \neq [] do
       for i := 1 to |Z| do
 3
           case 1. label(Z[i]) = \alpha
 4
              Delete Z[i] from Z (but keep it in L).
 5
           case 2. label(Z[i]) \neq \alpha and parent(Z[i]) \neq \bot
 6
              Replace Z[i] with parent(Z[i]) in both Z and L.
 7
           case 3. label(Z[i]) \neq \alpha and parent(Z[i]) = \bot
 8
              Delete Z[i] from both Z and L.
 9
10
11
       Set (Z, L) := \text{Deep}^*(Z, L).
12
13 end
14 Return L.
```

The procedure FL computes the set  $DEEP(\{fl(x,\alpha)|x\in X\})$  bottom up. The list Z contains ancestors of the elements of X for which we have not yet found an ancestor with label  $\alpha$ . In each step it considers each node z in the list Z. If it has the right label then  $x\in FL(X,\alpha)$  and we remove it from Z but keep it in L. Otherwise we replace it with its parent (unless it is the root). Thus L contains both the elements in Z and the part of  $FL(X,\alpha)$  found until now.

To keep the running time down we wish to maintain the invariant that L is deep at the beginning of each iteration of the outer loop. To do this procedure FL calls an auxiliary procedure DEEP\*(Z, L) which takes two ordered lists Z and L, where  $Z \subseteq L$ , and returns two ordered lists representing the set DEEP $(L) \cap Z$  and DEEP(L), i.e., DEEP\* $(Z, L) = ([z \in Z | \nexists x \in L : z \prec x], \text{DEEP}(L))$ . If we use the procedure DEEP to calculate DEEP\* it takes time O(|Z| + |L|) = O(|L|). Instead we will show how to calculate it in time O(|Z|) using a linked list representation for Z and L. We will need this in the proof of Lemma 12, which shows that the total running time of all calls to FL from EMB takes time  $O(n_T)$ . Below we describe in more detail how to implement FL together with the auxiliary procedures.

## **Procedure** MopLeft(X,Y)

```
1 Initially, set R := [].
 2 Find the largest j such that Y[|Y|]_1 \triangleright X[j] and set y := Y[|Y|]_2 and x := X[j]. If no such j exists
   stop and return R.
 3 for i := |Y| - 1 \ to \ 1 \ do
       until Y[i]_1 \triangleright X[j] or j < 1 do
 4
           set j := j - 1.
 5
       if j < 1 then
 6
           stop and return R := (x, y) \circ R.
 7
       else
 8
           compare X[j] and x. There are two cases:
 9
           case 1. x \triangleright X[j]
10
            set R := (x, y) \circ R, y := Y[i]_2, and x := X[j].
11
12
           case 2. x = X[j]
13
            set y := Y[i]_2.
14
15
       end
16
17 end
18 Return R := (x, y) \circ R.
```

We use a doubly linked list to represent L and extra pointers in this list to represent Z. Each element in the list has pointers  $\operatorname{Succ}_L$  and  $\operatorname{Pred}_L$  pointing to its predecessor and successor in L. Similarly, each element in Z has pointers  $\operatorname{Succ}_Z$  and  $\operatorname{Pred}_Z$  pointing to its predecessor and successor in Z (right after the initialization these are equal to  $\operatorname{Succ}_L$  and  $\operatorname{Pred}_L$ ). In the for loop we use the  $\operatorname{Succ}_Z$  pointers to find the next element in Z. To delete Z[i] from Z in case 1 we set  $\operatorname{Succ}_Z(\operatorname{Pred}_Z(Z[i])) = \operatorname{Succ}_Z(Z[i])$  and  $\operatorname{Pred}_Z(\operatorname{Succ}_Z(Z[i])) = \operatorname{Pred}_Z(Z[i])$ . The L pointers stay the same. In case 2 we simply replace Z[i] with its parent in the linked list. The  $\operatorname{Succ}$  and  $\operatorname{Pred}$  pointers stay the same. To delete Z[i] from both Z and L in case 3 we set  $\operatorname{Succ}_j(\operatorname{Pred}_j(Z[i])) = \operatorname{Succ}_j(Z[i])$  and  $\operatorname{Pred}_j(\operatorname{Succ}_j(Z[i])) = \operatorname{Pred}_j(Z[i])$  for  $j \in \{Z, L\}$ . Finally, to compute  $\operatorname{DEEP}^*(Z, L)$  walk through Z following the  $\operatorname{Succ}_Z$  pointers. At each node Z compare  $\operatorname{Pred}_L(Z)$  and  $\operatorname{Succ}_L(Z)$  with Z. If one of them is a descendant of Z remove Z from the doubly linked list Z and L as in case 3. Note that instead of calling  $\operatorname{DEEP}^*(Z, L)$  this comparison can also be done directly in step 2, which is the only place where we insert nodes that might be an ancestor of another node in L. We will show in the next section that it is enough to compare Z to its neighbors in the list L.

### 4.3 Correctness of the Set Procedures

Clearly, PARENT and NCA are correct. The following lemmas show that DEEP, FL, and MOPRIGHT are also correctly implemented. For notational convenience we write  $x \in X$ , for a list X, if x = X[i] for some i,  $1 \le i \le |X|$ .

**Lemma 7** Procedure DEEP(X) is correct.

*Proof.* Let x be the variable in the procedure. We will first prove the following invariant on x:

**Invariant** At the beginning of each iteration of the for loop in line 2 we have  $x \not\prec X[j]$  for any  $1 \le j \le i-1$ .

We prove the invariant by induction on i. The invariant obviously holds for the base case i=2. For the induction step let  $i \geq 3$ . Let iteration k denote the iteration of the for loop when i=k. By the induction hypothesis we have  $x \not\prec X[j]$  for any  $1 \leq j \leq i-2$  at the beginning of iteration i-1. Let x' denote the value of the variable x at the beginning of iteration i-1. Consider the value of variable x at the beginning of the iteration i. There are two cases:

- 1. If x = x' then by the induction hypothesis  $x = x' \not\prec X[j]$  for any  $1 \le j \le i 2$ . Since x was not changed in iteration i 1 we have  $X[i 1] \preceq x$  (case 3 of the procedure) and thus  $x \not\prec X[j]$  for any  $1 \le j \le i 1$ .
- 2. If  $x \neq x'$  then x was set in either case 1  $(x' \triangleleft x)$  or case 2  $(x' \triangleleft x)$  in iteration i-1. Therefore, x = X[i-1] and by the induction hypothesis  $x' \not\prec X[j]$  for any  $1 \leq j \leq i-2$ . There are two subcases:
  - (a) If  $x' \prec x$  it follows immediately from the induction hypothesis that  $x \not\prec X[j]$  for any  $1 \le j \le i-1$ , since all descendants of x also are descendants of x'.
  - (b) If  $x' \triangleleft x$  we note that x is the first node to the left of x' occurring after x' in X (otherwise x would have been reset in case 1 of the procedure in an earlier iteration, contradicting that x' is the value of variable x at the beginning of iteration i-1). Since X is semiordered no node X[j] with smaller index in X than x' can be to the right of x'. Thus no node X[j],  $1 \le j < i-2$ , can be to the right of x'. Since all descendants of x must be to the right of x' we have  $x \not\prec X[j]$  for any  $1 \le j \le i-1$ .

We are now ready to prove that  $y \in \text{DEEP}(X)$  iff there exists no  $z \in X$  such that  $y \prec z$ . We first argue that if  $y \in \text{DEEP}(X)$  then  $\nexists z \in X$  such that  $y \prec z$ . Let y be an element in DEEP(X). Only elements that have been assigned to x during the procedure are in the output. Consider the iteration where x = y is appended to R. This only happens in case 1 of the procedure and thus  $y = x \triangleleft X[i]$ . Since X is semiordered this implies that  $x \triangleleft X[j]$  for  $i \le j \le |X|$ , and therefore  $y = x \not\prec X[j]$  for  $i \le j \le |X|$ . By the above invariant it follows that  $y = x \not\prec X[j]$  for  $1 \le j \le i-1$ . Thus if  $y \in \text{DEEP}(X)$  then  $\nexists z \in X$  such that  $y \prec z$ .

Let  $y \in X$  be an element such that  $X \cap V(T(y)) = \{y\}$ . Let j be the smallest index such that X[j] = y. When comparing y and x during the iteration where i = j we are in case 1 or 2, since j is the smallest index such that X[j] = y (implying  $x \neq y$ ) and  $X \cap V(T(y)) = \{y\}$  (implying  $y \not\prec x$ ). In either case x is set to y. Since there are no descendants of y in X, the variable x remains equal to y until added to R. If y occurs several times in X we will have x = y each time we meet a copy of y (except the first) and it follows from the implementation that y will occur exactly once in R.

To show that the implementation of Moprish is correct we will use the following proposition.

**Proposition 3** Before the first iteration of the for loop in line 3 of MOPRIGHT we have  $y = Y[1]_1$ , x = X[j] and either  $X[j-1] \triangleleft Y[1]_2 \triangleleft X[j]$  (if j > 1) or  $Y[1]_2 \triangleleft X[j]$  if (j = 1).

At the end of each iteration of the for loop then, unless  $Y[i]_2 \not \triangleleft X[|X|]$ , we have  $y = Y[i]_1$ , x = X[j] and either  $X[j-1] \triangleleft Y[i]_2 \triangleleft X[j]$  (if j > 1) or  $Y[i]_2 \triangleleft X[j]$  if (j = 1).

*Proof.* The first statement  $(y = Y[1]_1, x = X[j])$  and either  $X[j-1] \triangleleft Y[1]_2 \triangleleft X[j]$  (if j > 1) or  $Y[1]_2 \triangleleft X[j]$  if (j = 1)) follows immediately from the implementation of the procedure line 2 and the fact that the input lists are ordered.

We prove the second statement by induction on i. Base case i=2. By the first statement we have  $y=Y[1]_1, \ x=X[j]$  and either  $X[j-1] \triangleleft Y[1]_2 \triangleleft X[j]$  (if j>1) or  $Y[1]_2 \triangleleft X[j]$  if (j=1) before this iteration. Let j' be the value of j before this iteration. It follows immediately from the implementation that  $y=Y[2]_1$  since y is set to this in both case 1 and 2. If  $Y[2]_2 \triangleleft X[j']$  then j=j'. Since Y is ordered it follows that  $X[j-1] \triangleleft Y[1]_2 \triangleleft Y[2]_2 \triangleleft X[j]$  (if j>1) or  $Y[2]_2 \triangleleft X[j]$  if (j=1). If  $Y[2]_2 \not A[j']$  then j is increased until  $Y[2]_2 \triangleleft X[j]$  implying  $X[j-1] \triangleleft Y[2]_2 \triangleleft X[j]$  unless j>|X|, since X is ordered.

Induction step i > 2. It follows immediately from the implementation that  $y = Y[i]_1$  since y is set to this in both case 1 and 2. By the induction hypothesis we have  $y = Y[i]_1$ , x = X[j] and  $Y[i]_2 \triangleleft X[j]$  right before this iteration. Let j' be the value of j before this iteration. If  $Y[i]_2 \triangleleft X[j']$  then j = j'. Since Y is ordered it follows that  $X[j-1] \triangleleft Y[i-1]_2 \triangleleft Y[i]_2 \triangleleft X[j]$  (if j > 1) or  $Y[i]_2 \triangleleft X[j]$  if (j = 1). If  $Y[i]_2 \not \triangleleft X[j']$  then j is increased until  $Y[i]_2 \triangleleft X[j]$  implying  $X[j-1] \triangleleft Y[i]_2 \triangleleft X[j]$  unless j > |X|.

**Lemma 8** Procedure MopRight(Y, X) is correct.

*Proof.* We want to show that for any  $1 \le i' \le |Y|$ ,  $1 \le j' \le |X|$ :

$$(Y[i']_1, X[j']) \in R \quad \Leftrightarrow \quad (Y[i']_2, X[j']) \in \operatorname{mop}(Y|_2, X) .$$

Since  $Y|_2$  and X are ordered lists we have  $(Y[i']_2, X[j']) \in \text{mop}(Y|_2, X)$  if and only if

- (1)  $\operatorname{arg\,min}_{j} Y[i']_{2} \triangleleft X[j] = j'$
- (2)  $\arg \max_i Y[i]_2 \triangleleft X[j'] = i'$

We will first show that (1) and (2) implies  $(Y[i']_2, X[j']) \in R$ . We start by showing that when i is about to be incremented to i' + 1 then  $y = Y[i']_1$  and x = X[j']. There are two cases to consider.

- i' = 1. After line 2 is executed, y is set to  $Y[1]_1$ , j is set to j' and x is set to X[j'].
- i' > 1. Consider the step in the iteration when i = i'. At the beginning of this iteration, y = Y[i' 1] and j is the minimal index such that  $Y[i' 1] \triangleleft X[j]$ . By (1) this implies that  $j \leq j'$ , and that after line 4 and 5 are executed, j is set to j'. At the end of the iteration y = Y[i'] (y is assigned to Y[i'] in both cases) and x = X[j']. If j = j' then x set to X[j'] in case 1, otherwise we had j = j' (case 2) and then x was set to X[j'] already).

We have established that when i is about to be incremented to i' + 1 then  $y = Y[i']_1$  and x = X[j']. To show that  $(Y[i']_2, X[j']) \in R$  we consider the following two cases.

- i' < |Y|. Consider the  $(i'+1)^{th}$  iteration. By condition (2)  $X[j'] \le Y[i'+1]_2$  and therefore j is increased in line 5. So now j > j'. If j > |X| then  $(y,x) = (Y[i']_2, X[j'])$  is added to R in line 7. Otherwise, since X is ordered,  $x = X[j'] \triangleleft X[j]$ . We are therefore in case 1 and  $(y,x) = (Y[i']_2, X[j'])$  is added to R.
- i' = |Y|. Then  $(y, x) = (Y[i']_2, X[j'])$  is added to R in line 15.

We will now show that  $(Y[i']_1, X[j']) \in R$  implies (1) and (2). Since  $(Y[i']_1, X[j']) \in R$  we had  $(y, x) = (Y[i']_1, X[j'])$  at some point during the execution. The pair (y, x) can be added to R only in the for loop before changing the values of y and x or at the execution of the last line of the procedure. Therefore  $(y, x) = (Y[i']_1, X[j'])$  at the beginning of some execution of the for loop, or after the last iteration (i = |Y|). It follows by Proposition 3 that  $X[j-1] \triangleleft Y[i]_2 \triangleleft X[j]$  if j > 1 or  $Y[i]_2 \triangleleft X[j]$  if j = 1. It remains to show that  $X[j'] \triangleleft Y[i'+1]_2$  for i' < |Y|. It follows from the implementation that (y, x) only is added to R inside the for loop if j is increased. Thus j was increased in the next iteration (i = i'+1) implying  $X[j'] \triangleleft Y[i'+1]_2$ .

**Lemma 9** Procedure MopLeft(X, Y) is correct.

*Proof.* Similar to the proof of Lemma 8.

To show that FL is correct we need the following proposition.

**Proposition 4** Let X be an ordered list and let x be an ancestor of X[i] for some  $i \in \{1, ..., k\}$ . If x is an ancestor of some node in X other than X[i] then x is an ancestor of X[i-1] or X[i+1].

*Proof.* Recall that  $u \triangleleft v$  iff  $\operatorname{pre}(u) < \operatorname{pre}(v)$  and  $\operatorname{post}(u) < \operatorname{post}(v)$ . Since  $x \triangleleft X[i]$  we have  $\operatorname{pre}(x) < \operatorname{pre}(X[i])$  and  $\operatorname{post}(X[i]) < \operatorname{post}(x)$ . Assume there exists a descendant X[j] of x such that  $j \notin \{i-1, i, i+1\}$ . If j < i-1 we have

$$\operatorname{pre}(x) \le \operatorname{pre}(X[j]) < \operatorname{pre}(X[i-1]),$$

where the first inequality follows from  $x \prec X[j]$  and the second from X being ordered. And

$$post(X[i-1]) < post(X[i]) \le post(x),$$

where the first inequality follows from X being ordered and the second from  $x \prec X[i]$ . Thus  $x \prec X[i-1]$ . Similarly, for j > i+1, we have  $\operatorname{pre}(x) \leq \operatorname{pre}(X[i]) < \operatorname{pre}(X[i+1])$  and  $\operatorname{post}(X[i+1]) < \operatorname{post}(X[j]) \leq \operatorname{post}(x)$  implying that  $x \prec X[i+1]$ .

Proposition 4 shows that the doubly linked list implementation of DEEP\* is correct. Since all changes to the list are either deletions or insertions of a parent in the place of its child, the list L (and thus also Z) is ordered at the beginning of each iteration of the outer loop.

**Lemma 10** Procedure  $FL(X, \alpha)$  is correct.

*Proof.* Let  $F = \{f(x, \alpha) \mid x \in X\}$ . We first show that  $FL(X, \alpha) \subseteq F$ . Consider a node  $x \in FL(X, \alpha)$ . Since x is in L after the final iteration, x was deleted from Z during some iteration. Thus label $(x) = \alpha$ . For any  $y \in X$  we follow the path from y to the root and stop the first time we meet a node with label  $\alpha$  or even earlier since we keep the list deep. Thus  $x \in F$ .

The set  $\operatorname{FL}(X,\alpha)$  is a deep set, and therefore  $\operatorname{DEEP}(F) \subseteq \operatorname{FL}(X,\alpha) \subseteq F \Rightarrow \operatorname{DEEP}(F) = \operatorname{FL}(X,\alpha)$ . Hence, it remains to show that  $\operatorname{DEEP}(F) \subseteq \operatorname{FL}(X,\alpha)$ . Let x be a node in  $\operatorname{DEEP}(F)$ , let  $z \in X$  be a node such that  $x = \operatorname{fl}(z,\alpha)$ , and let  $z = x_1, x_2, \ldots, x_k = x$  be the nodes on the path from z to x. We will argue that after each iteration of the algorithm we have  $x_i \in L$  for some i. Since  $\operatorname{label}(x_i) \neq \alpha$  for i < k this is the same as  $x_i \in Z$  for i < k. Before the first iteration we have  $x_1 \in X = Z$ . As long as i < k we replace  $x_i$  with  $x_{i+1}$  in case 2 of the for loop, since  $\operatorname{label}(x_i) \neq \alpha$ . When i = k we remove  $x_k$  from Z but keep it in L. It remains to show that we do not delete  $x_i$  in the computation of  $\operatorname{DEEP}^*(Z,L)$  in any iteration. If  $x_i$  is removed then there is a node  $y \in L$  that is a descendant of  $x_i$  and thus also a descendant of x. We argued above that  $L \setminus Z \subseteq F$  and thus  $y \in Z$  since  $x \in \operatorname{DEEP}(F)$ . But since  $x \in \operatorname{DEEP}(F)$  no node on the path from y to x can have label  $\alpha$  and therefore  $x_i$  will eventually be reinserted in Z.

#### 4.4 Complexity of the Set Procedures

For the running time of the node list implementation observe that, given the data structure described in Section 4.1, all set procedures, except FL, perform a single pass over the input using constant time at each step. Hence we have,

**Lemma 11** For any tree T there is a data structure using  $O(n_T)$  space and preprocessing which supports each of the procedures PARENT, DEEP, MOPRIGHT, MOPLEFT, and NCA in linear time (in the size of their input).

The running time of a single call to FL might take time  $O(n_T)$ . Instead we will divide the calls to FL into groups and analyze the total time used on such a group of calls. The intuition behind the division is that for a path in P the calls made to FL by EMB are done bottom up on disjoint lists of nodes in T.

**Lemma 12** For disjoint ordered node lists  $X_1, \ldots, X_k$  and labels  $\alpha_1, \ldots, \alpha_k$ , such that any node in  $X_{i+1}$  is an ancestor of some node in  $\operatorname{FL}_T(X_i, \alpha_i)$ ,  $1 \leq i < k$ , all of  $\operatorname{FL}_T(X_1, \alpha_1), \ldots, \operatorname{FL}_T(X_k, \alpha_k)$  can be computed in  $O(n_T)$  time.

*Proof.* Let Z and L be as in the implementation of the procedure. Since DEEP\* takes time O(|Z|) and each of the steps in the for loop takes constant time, we only need to show that the total length of the lists Z—summed over all the calls—is  $O(n_T)$  to analyze the total time usage. We will show that any node in T can be in Z at the beginning of the while loop at most twice during all calls to FL. The size of Z cannot increase in the iterations of the for loop (line 3–10), and thus the size of Z when DEEP\* is called (line 11) is at most the size of Z at the beginning of this iteration of the while loop.

Consider a single call to FL. Except for the first iteration, a node can be in Z only if one of its children were in Z in the last iteration. Note that Z is ordered at the beginning of each for loop. Thus if a node is in Z at the beginning of the while loop none of its children are in Z and thus in one call to FL a node can be in Z only once.

Look at a node z the first time it appears in Z at the beginning of an execution of the while loop. Assume that this is in the call  $FL(X_i, \alpha_i)$ .

- If  $z \in X_i$  then z cannot be in Z in any later calls, since no node in  $X_j$  where j > i can be a descendant of a node in  $X_i$ .
- If  $label(z) \neq \alpha_i$  then z is removed from Z in case 2 or case 3 of the procedure and cannot be in Z in any of the later calls. To see this consider the time when z is removed from Z (case 2 or case 3). Since the set L is deep at the beginning of the while loop and  $Z \subseteq L$ , no descendant of z will appear in Z later in this call to FL, and no node in the output from  $FL(X_i, \alpha_i)$  can be a descendant of z. Since any node in  $X_j$ , j > i, is an ancestor of some node in  $FL(X_i, \alpha_i)$  neither z or any descendant of z can be in any  $X_j$ , j > i. Thus z cannot appear in Z in any later calls to FL.
- Now if label(z) =  $\alpha_i$  then we might have  $z \in X_{i+1}$ . In that case, z will appear in Z in the first iteration of the procedure call  $FL(X_{i+1}, \alpha_i)$ , but not in any later calls since the lists are disjoint, and since no node in  $X_j$  where j > i+1 can be a descendant of a node in  $X_{i+1}$ . If label(z) =  $\alpha_i$  and  $z \notin X_{i+1}$  then clearly z cannot appear in Z in any later call.

Thus a node in T is in Z at the beginning of an execution of the while loop at most twice during all the calls.  $\Box$ 

#### 4.5 Complexity of the Tree Inclusion Algorithm

Using the node list implementation of the set procedures we get:

**Lemma 13** For trees P and T the tree inclusion problem can be solved in  $O(l_P n_T)$  time.

Proof. By Lemma 11 we can preprocess T in  $O(n_T)$  time and space. Let g(n) denote the time used by FL on a list of length n. Consider the time used by EMB(root(P)). We bound the contribution for each node  $v \in V(P)$ . If v is a leaf we are in case 1 of EMB. The cost of computing  $\operatorname{FL}(L(T), \operatorname{label}(v))$  is  $O(g(l_T))$ , and by Lemma 12 (with k=1) we get  $O(g(l_T)) = O(n_T)$ . Hence, the total cost of all leaves is  $O(l_P n_T)$ . If v has a single child w we are in case 2 of EMB, and by Lemma 11 the cost is  $O(g(|\operatorname{EMB}(w)|))$ . If v has more than one child the cost of Mopright, NCA, and Deep is bounded by  $\sum_{w \in \operatorname{child}(v)} O(|\operatorname{EMB}(w)|)$ . Furthermore, since the length of the output of Mopright (and thus NCA) is at most  $z = \min_{w \in \operatorname{child}(v)} |\operatorname{EMB}(w)|$  the cost of FL is O(g(z)). Hence, the total cost for internal nodes is,

$$\sum_{v \in V(P) \setminus L(P)} O\left(g(\min_{w \in \text{child}(v)} |\text{EMB}(w)|) + \sum_{w \in \text{child}(v)} |\text{EMB}(w)|\right) = \sum_{v \in V(P)} O(g(|\text{EMB}(v)|)). \tag{3}$$

Next we bound (3). For any  $w \in \text{child}(v)$  we have that Emb(w) and Emb(v) are disjoint ordered lists. Furthermore we have that any node in Emb(v) must be an ancestor of some node in FL(Emb(w), label(v)).

Hence, by Lemma 12, for any leaf to root path  $\delta = v_1, \ldots, v_k$  in P, we have that  $\sum_{u \in \delta} g(|\text{EmB}(u)|) = O(n_T)$ . Let  $\Delta$  denote the set of all root to leaf paths in P. It follows that,

$$\sum_{v \in V(T)} g(|\mathrm{Emb}(v)|) \le \sum_{p \in \Delta} \sum_{u \in p} g(|\mathrm{Emb}(u)|) = O(l_P n_T) .$$

Since this time is the same as the time spent at the leaves the time bound follows.

To analyze the space used by the algorithm we first bound the size of Emb(v) for each node  $v \in V(P)$ . We then use this to bound the total the size of embeddings stored in the recursion stack in the computation of Emb(root(P)), i.e., the total size of embeddings stored by recursive calls during the computation.

**Lemma 14** For any tree P we have  $\forall v \in V(P)$ :

$$|\mathrm{Emb}_T(v)| \le \frac{l_T}{l_{P(v)}}$$
.

*Proof.* By Lemma 5 EMB(v) is the set of deep occurrences of P(v) in T. By the definition of deep the occurrences are disjoint and no node in one occurrence can be an ancestor of a node in another occurrence. Each occurrence has at least  $l_{P(v)}$  descendant leaves and each of these leaves is an ancestor of at least one distinct leaf in T (see also Figure 4(b)). Thus the number of occurrences is bounded by  $l_T/l_{P(v)}$ .

**Lemma 15** The total size of saved embeddings on the recursion stack at any time during the computation of EmB(root(P)) is at most  $O(l_T)$ .

Proof. Let node v be the node for which we are currently computing EMB. Let p be the path from the root to v and let  $w_0, \ldots, w_\ell$  be the light nodes on this path. Let  $\ell = \operatorname{ldepth}(v)$ . There is one embedding on the stack for each light node on the path (see Figure 6): For the heavy nodes on the path there can be no saved embeddings in the recursion as the algorithm always recurses on the heavy child first. For each light node  $w_i$  on the path p except the root  $w_0$  the stack will contain either EMB(heavy(parent( $w_i$ ))), or  $U_j = \operatorname{MopRight}(U_{j-1}, R_j)$ , where  $v_j$  is  $w_i$ 's right sibling. The computation of  $U_j$  is a series of MopRight (or MopLeft) computations that started with the pair of node lists (EMB(heavy(parent( $w_i$ ))), EMB(heavy(parent( $w_i$ )))) as the first argument to MopRight (or MopLeft). As the output of MopRight (or MopLeft) can be no larger than the input to the procedure we have  $|U_j| = O(|\operatorname{EMB}(\operatorname{heavy}(\operatorname{parent}(w_i))))|$  and thus the total space used at any time during the recursion is

$$O\left(\sum_{i=1}^{\text{ldepth}(v)} |\text{EMB}(\text{heavy}(\text{parent}(w_i)))|\right).$$

By Lemma 14 we have

$$|\text{EmB}(\text{heavy}(\text{parent}(w_i)))| \le \frac{l_T}{l_{P(\text{heavy}(\text{parent}(w_i)))}},$$

and thus

$$\sum_{i=1}^{\text{ldepth}(v)} |\text{EMB}(\text{heavy}(\text{parent}(w_i)))| \le l_T \sum_{i=1}^{\text{ldepth}(v)} \frac{1}{l_{P(\text{heavy}(\text{parent}(w_i)))}}.$$
 (4)

By the definition of heavy the node heavy (parent $(w_i)$ ) has more leaves in its subtree than  $w_i$ , i.e.,

$$l_{P(w_i)} \le l_{P(\text{heavy}(\text{parent}(w_i)))}$$
 (5)

Obviously, heavy(parent( $w_i$ )) has no more leaves in its subtree than its parent, i.e.,

$$l_{P(\text{heavy}(\text{parent}(w_i)))} \le l_{P(\text{parent}(w_i))}$$
 (6)

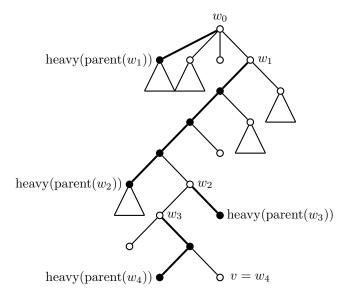


Figure 6: Path from root to v. The heavy nodes are black and the light nodes are white. The heavy edges are the thick edges and the light edges are thin.

Since  $w_i$  is light it has at most half the number of leaves in its subtree as its parent, that is

$$l_{P(w_i)} \le l_{P(\text{parent}(w_i))}/2. \tag{7}$$

Combining this with the fact that  $w_i$  is an ancestor of  $w_{i+1}$  and heavy(parent( $w_{i+1}$ )) we get,

$$\begin{split} l_{P(\text{heavy}(\text{parent}(w_{j}))} &\leq l_{P(\text{parent}(w_{j}))} & \text{by (6)} \\ &\leq l_{P(w_{j-1})} & \text{since } w_{j-1} \text{ is an ancestor of } w_{j} \\ &\leq l_{P(\text{parent}(w_{j-1}))}/2 & \text{by (7)} \\ &\leq l_{P(w_{j-2})}/2 & \text{since } w_{j-2} \text{ is an ancestor of parent}(w_{j-1}) \\ &\leq l_{P(\text{heavy}(\text{parent}(w_{j-2})))}/2, & \text{by (5)} \end{split}$$

for any  $2 < j \le \text{ldepth}(v)$ . Let  $l_i = l_{P(\text{heavy}(\text{parent}(w_i)))}$  for all i. To bound the sum in (4) we will use that  $l_i \le l_{i-2}/2$ ,  $l_i < l_{i-1}$ , and  $l_{\text{ldepth}(v)} \ge 1$ . We have

$$\sum_{i=1}^{\operatorname{ldepth}(v)} \frac{1}{l_i} \leq 2 \sum_{i=2, i \text{ odd}}^{\operatorname{ldepth}(v)} \frac{1}{l_i} \leq 2 \cdot 2 = 4,$$

since the  $l_i$ 's in the last sum are decreasing with a factor of 2. Combining this with Equation (4) we get

$$\sum_{i=1}^{\operatorname{depth}(v)} |\operatorname{Emb}(\operatorname{heavy}(\operatorname{parent}(w_i)))| \leq l_T \sum_{i=1}^{\operatorname{ldepth}(v)} \frac{1}{l_{P(\operatorname{heavy}(\operatorname{parent}(w_i)))}} \leq 4l_T.$$

**Theorem 2** For trees P and T the tree inclusion problem can be solved in  $O(l_P n_T)$  time and  $O(n_T)$  space.

*Proof.* The time bound follows from Lemma 13. Next consider the space used by EMB(root(P)). The preprocessing of Section 4.1 uses only  $O(n_T)$  space. By Lemma 15 the space used for the saved embeddings is  $O(l_T) = O(n_T)$ .

## 4.6 An Alternative Algorithm

In this section we present an alternative algorithm. Since the time complexity of the algorithm in the previous section is dominated by the time used by FL, we present an implementation of this procedure which leads to a different complexity. Define a firstlabel data structure as a data structure supporting queries of the form  $fl(v,\alpha)$ ,  $v \in V(T)$ ,  $\alpha \in \Sigma$ . Maintaining such a data structure is known as the tree color problem. This is a well-studied problem, see e.g. [5, 16, 18, 31]. With such a data structure available we can compute FL as follows.

 $FL(X, \alpha)$ : Return the list  $DEEP([fl(X[1], \alpha), ..., fl(X[|X|], \alpha)])$ .

**Theorem 3** Let P and T be trees. Given a firstlabel data structure using  $s(n_T)$  space,  $p(n_T)$  preprocessing time, and  $q(n_T)$  time for queries, the tree inclusion problem can be solved in  $O(p(n_T) + l_P l_T \cdot q(n_T))$  time and  $O(s(n_T) + n_T)$  space.

Proof. Constructing the firstlabel data structures uses  $O(s(n_T))$  space and  $O(p(n_T))$  time. The total cost of the leaves is bounded by  $O(l_p l_T \cdot q(n_T))$ , since the cost of a single leaf is  $O(l_T \cdot q(n_T))$ . As in the proof of Theorem 2 we have that the total time used by the internal nodes is bounded by  $\sum_{v \in V(P)} g(|\text{EMB}(v)|)$ , where g(n) is the time used by FL on a list of length n, that is,  $g(n) \leq n \cdot q(n_T)$ . By Lemma 11 and Lemma 15 for any leaf to root path  $\delta = v_1, \ldots, v_k$  in P, we have that  $\sum_{u \in \delta} |\text{EMB}(u)| \leq O(l_T)$ . Let  $\Delta$  denote the set of all root to leaf paths in P. It follows that,

$$\sum_{v \in V(P)} g(|\mathrm{Emb}(v)|) \leq \sum_{p \in \Delta} \sum_{u \in p} g(|\mathrm{Emb}(u)|) \leq \sum_{p \in \Delta} O(l_T \cdot q(n_T)) \leq O(l_P l_T \cdot q(n_T)).$$

Since this time is the same as the time spent at the leaves the time bound follows.

Several firstlabel data structures are available, for instance, if we want to maintain linear space we have,

**Lemma 16 (Dietz [16])** For any tree T there is a data structure using  $O(n_T)$  space,  $O(n_T)$  expected preprocessing time which supports firstlabel queries in  $O(\log \log n_T)$  time.

The expectation in the preprocessing time is due to perfect hashing. Since our data structure does not need to support efficient updates we can remove the expectation by using the deterministic dictionary of Hagerup et al. [20]. This gives a worst-case preprocessing time of  $O(n_T \log n_T)$ . However, using a simple two-level approach this can be reduced to  $O(n_T)$  (see e.g. [38]). Plugging in this data structure we obtain,

**Corollary 2** For trees P and T the tree inclusion problem can be solved in  $O(l_P l_T \log \log n_T + n_T)$  time and  $O(n_T)$  space.

# 5 A Faster Tree Inclusion Algorithm

In this section we present a new tree inclusion algorithm which has a worst-case subquadratic running time. As discussed in the introduction, the general idea is to divide T into clusters of logarithmic size which we can efficiently preprocess and then use this to speed up the computation with a logarithmic factor.

## 5.1 Clustering

In this section we describe how to divide T into clusters and how the macro tree is created. For simplicity in the presentation we assume that T is a binary tree. If this is not the case it is straightforward to construct a binary tree B, where  $n_B \leq 2n_T$ , and a mapping  $g: V(T) \to V(B)$  such that for any pair of nodes  $v, w \in V(T)$ , label $(v) = \text{label}(g(v)), v \prec w$  iff  $g(v) \prec g(w)$ , and  $v \prec w$  iff  $g(v) \prec g(w)$ . The nodes in the set  $U = V(B) \setminus \{g(v) \mid v \in V(T)\}$  are assigned a special label  $\beta \notin \Sigma$ . It follows that for any tree  $P, P \sqsubseteq T$  iff  $P \sqsubseteq B$ .

Let C be a connected subgraph of T. A node in V(C) adjacent to a node in  $V(T)\backslash V(C)$  is a boundary node. The boundary nodes of C are denoted by  $\delta C$ . We have  $\mathrm{root}(T) \in \delta C$  if  $\mathrm{root}(T) \in V(C)$ . A cluster of C is a connected subgraph of C with at most two boundary nodes. A set of clusters CS is a cluster partition of T iff  $V(T) = \bigcup_{C \in CS} V(C)$ ,  $E(T) = \bigcup_{C \in CS} E(C)$ , and for any  $C_1, C_2 \in CS$ ,  $E(C_1) \cap E(C_2) = \emptyset$ ,  $|E(C_1)| \geq 1$ . If  $|\delta C| = 1$  we call C a leaf cluster and otherwise an internal cluster.

We use the following recursive procedure  $\text{CLUSTER}_T(v, s)$ , adopted from [6], which creates a cluster partition CS of the tree T(v) with the property that |CS| = O(s) and  $|V(C)| \leq \lceil n_T/s \rceil$  for each  $C \in CS$ . A similar cluster partitioning achieving the same result follows from [3, 4, 19].

Cluster T(v, s): For each child u of v there are two cases:

- 1.  $|V(T(u))| + 1 \le \lceil n_T/s \rceil$ . Let the nodes  $\{v\} \cup V(T(u))$  be a leaf cluster with boundary node v.
- 2.  $|V(T(u))| \ge \lceil n_T/s \rceil$ . Pick a node  $w \in V(T(u))$  of maximum depth such that  $|V(T(u))| + 2 |V(T(w))| \le \lceil n_T/s \rceil$ . Let the nodes  $V(T(u)) \setminus V(T(w)) \cup \{v, w\}$  be an internal cluster with boundary nodes v and w. Recursively, compute CLUSTER $_T(w, s)$ .

**Lemma 17** Given a tree T with  $n_T > 1$  nodes, and a parameter s, where  $\lceil n_T/s \rceil \ge 2$ , we can build a cluster partition CS in  $O(n_T)$  time, such that |CS| = O(s) and  $|V(C)| \le \lceil n_T/s \rceil$  for any  $C \in CS$ .

Proof. The procedure CLUSTER<sub>T</sub>(root(T), s) clearly creates a cluster partition of T and it is straightforward to implement in  $O(n_T)$  time. Consider the size of the clusters created. There are two cases for u. In case  $1, |V(T(u))| + 1 \le \lceil n_T/s \rceil$  and hence the cluster  $C = \{v\} \cup V(T(u))$  has size  $|V(C)| \le \lceil n_T/s \rceil$ . In case  $2, |V(T(u))| + 2 - |V(T(w))| \le \lceil n_T/s \rceil$  and hence the cluster  $C = V(T(u)) \setminus V(T(w)) \cup \{v, w\}$  has size  $|V(C)| \le \lceil n_T/s \rceil$ .

Next consider the size of the cluster partition. Let  $c = \lceil n_T/s \rceil$ . We say that a cluster C is bad if  $|V(C)| \le c/2$  and good otherwise. We will show that at least a constant fraction of the clusters in the cluster partition are good. It is easy to verify that the cluster partition created by procedure Cluster has the following properties:

- (i) Let C be a bad internal cluster with boundary nodes v and w ( $v \prec w$ ). Then w has two children with at least c/2 descendants each.
- (ii) Let C be a bad leaf cluster with boundary node v. Then the boundary node v is contained in a good cluster.

By (ii) the number of bad leaf clusters is at most twice the number of good internal clusters and by (i) each bad internal cluster has two child clusters. Therefore, the number of bad internal clusters is bounded by the number of leaf clusters. Let  $b_i$  and  $g_i$  denote the number of bad and good internal clusters, respectively, and let  $b_l$  and  $g_l$  denote the number of bad and good leaf clusters, respectively. We have

$$b_i \le b_l + g_l \le 2g_i + g_l,$$

and therefore the number of bad clusters is bounded by

$$b_l + b_i \le 2g_i + g_l + 2g_i = 4g_i + g_l$$
.

Thus the number of bad clusters is at most 4 times the number of good clusters, and therefore at most a constant fraction of the total number of clusters. Since a good cluster is of size more than c/2, there can be at most 2s good clusters and thus |CS| = O(s).

Let  $C \in CS$  be an internal cluster with  $v, w \in \delta C$ . The *spine path* of C is the path between v, w excluding v and w. A node on the spine path is a *spine node*. A node to the left and right of v or of any node on

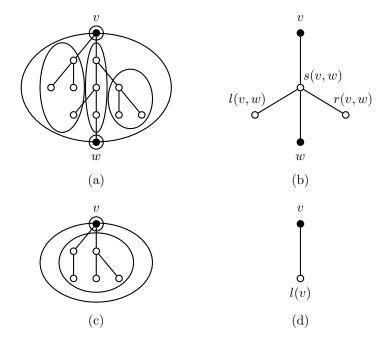


Figure 7: The clustering and the macro tree. (a) An internal cluster. The black nodes are the boundary nodes and the internal ellipses correspond to the boundary nodes, the right and left nodes, and spine path. (b) The macro tree corresponding to the cluster in (a). (c) A leaf cluster. The internal ellipses are the boundary node and the leaf nodes. (d) The macro tree corresponding to the cluster in (c).

the spine path is a *left node* and *right node*, respectively. If C is a leaf cluster with  $v \in \delta C$  then any proper descendant of v is a *leaf node*.

Let CS be a cluster partition of T as described in Lemma 17. We define an ordered macro tree M. Our definition of M may be viewed as an "ordered" version of the macro tree defined in [6]. The node set V(M) consists of the boundary nodes in CS. Additionally, for each internal cluster  $C \in CS$  with  $v, w \in \delta C$ ,  $v \prec w$ , we have the nodes s(v, w), l(v, w) and r(v, w) and edges (v, s(v, w)), (s(v, w), l(v, w)), (s(v, w), w), and (s(v, w), r(v, w)). That is, the nodes l(v, w), r(v, w) and w are all children of s(v, w). The nodes are ordered so that  $l(v, w) \vartriangleleft w \vartriangleleft r(v, w)$ . For each leaf cluster C,  $v \in \delta C$ , we have the node l(v) and edge (v, l(v)). Since root(T) is a boundary node, M is rooted at root(T). Figure 7 illustrates these definitions.

With each node  $v \in V(T)$  we associate a unique macro node denoted c(v). Let  $u \in V(C)$ , where  $C \in CS$ .

$$c(u) = \begin{cases} u & \text{if } u \text{ is boundary node,} \\ l(v) & \text{if } u \text{ is a leaf node and } v \in \delta C, \\ s(v,w) & \text{if } u \text{ is a spine node, } v,w \in \delta C, \text{ and } v \prec w, \\ l(v,w) & \text{if } u \text{ is a left node, } v,w \in \delta C, \text{ and } v \prec w, \\ r(v,w) & \text{if } u \text{ is a right node, } v,w \in \delta C, \text{ and } v \prec w. \end{cases}$$

Conversely, for any macro node  $i \in V(M)$  define the *micro forest*, denoted C(i), as the induced subgraph of T of the set of nodes  $\{v \mid v \in V(T), i = c(v)\}$ . We also assign a *set* of labels to i given by label $(i) = \{label(v) \mid v \in V(C(i))\}$ . If i is a spine node or a boundary node the unique node in V(C(i)) of greatest depth is denoted by first(i). Finally, for any set of nodes  $\{i_1, \ldots, i_k\} \subseteq V(M)$  we define  $C(i_1, \ldots, i_k)$  as the induced subgraph of the set of nodes  $V(C(i)) \cup \cdots \cup V(C(i))$ .

The following propositions state useful properties of ancestors, nearest common ancestor, and the left-to-right ordering in the micro forests and in T. The propositions follow directly from the definition of the clustering. See also Figure 8.

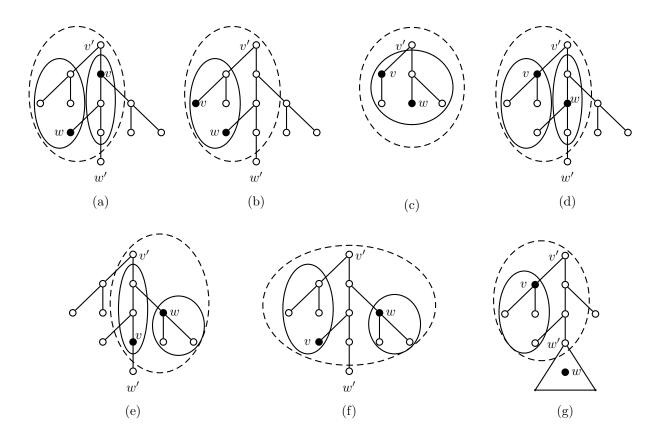


Figure 8: Examples from the propositions. In all cases v' and w' are top and bottom boundary nodes of the cluster, respectively. (a) Proposition 5(ii). Here c(v) = s(v', w') and c(w) = l(v', w') (solid ellipses). The dashed ellipse corresponds to C(c(w), s(v', w'), v'). (b) Proposition 6(i) and 7(ii). Here c(v) = c(w) = l(v', w') (solid ellipse). The dashed ellipse corresponds to C(c(w), s(v', w'), v'). (c) Proposition 6(ii) and 7(i). Here c(v) = c(w) = l(v') (solid ellipse). The dashed ellipse corresponds to C(c(v), v'). (d) Proposition 6(iii). Here c(v) = l(v', w') and c(w) = s(v', w') (solid ellipses). The dashed ellipse corresponds to C(c(v), c(w), v'). (e) Proposition 6(iv). Here c(v) = s(v', w') and c(w) = r(v', w') (solid ellipses). The dashed ellipse corresponds to C(c(v), c(w), v'). (f) Proposition 7(iv). Here c(v) = l(v', w') and c(w) = r(v', w') (solid ellipses). The dashed ellipse corresponds to C(c(v), c(w), s(v', w'), v'). (g) Proposition 7(v). Here c(v) = l(v', w') (solid ellipse) and  $w' \leq_M c(w)$ . The dashed ellipse corresponds to C(c(v), s(v', w'), v', w').)

**Proposition 5 (Ancestor relations)** For any pair of nodes  $v, w \in V(T)$ , the following hold

- (i) If c(v) = c(w) then  $v \prec_T w$  iff  $v \prec_{C(c(v))} w$ .
- (ii) If  $c(v) \neq c(w)$ , and for some boundary nodes v', w' we have c(v) = s(v', w'), and  $c(w) \in \{l(v', w'), r(v', w')\}$ , then  $v \prec_T w$  iff  $v \prec_{C(c(w), s(v', w'), v')} w$ .
- (iii) In all other cases,  $v \prec_T w$  iff  $c(v) \prec_M c(w)$ .

Case (i) says that if v and w belong to the same macro node then v is an ancestor of w iff v is an ancestor of w in the micro forest for that macro node. Case (ii) says that if v is a spine node and w is a left or right node in the same cluster then v is an ancestor of w iff v is an ancestor of w in the micro tree induced by that cluster (Figure 8(a)). Case (iii) says that in all other cases v is an ancestor of w iff the macro node v belongs to is an ancestor of the macro node w belongs to in the macro tree.

**Proposition 6 (Left-of relations)** For any pair of nodes  $v, w \in V(T)$ , the following hold

- (i) If  $c(v) = c(w) \in \{r(v', w'), l(v', w')\}$  for some boundary nodes v', w', then  $v \triangleleft w$  iff  $v \triangleleft_{C(c(v), v', s(v', w'))} w$ .
- (ii) If c(v) = c(w) = l(v') for some boundary node v', then  $v \triangleleft w$  iff  $v \triangleleft_{C(c(v),v')} w$ .
- (iii) If c(v) = l(v', w') and c(w) = s(v', w') for some boundary nodes v', w', then  $v \triangleleft w$  iff  $v \triangleleft_{C(c(v), c(w), v')} w$ .
- (iv) If c(v) = s(v', w') and c(w) = r(v', w') for some boundary nodes v', w', then  $v \triangleleft w$  iff  $v \triangleleft_{C(c(v), c(w), v')} w$ .
- (v) In all other cases,  $v \triangleleft w$  iff  $c(v) \triangleleft_M c(w)$ .

Case (i) says that if v and w are both either left or right nodes in the same cluster then v is to the left of w in the micro tree induced by their macro node together with the spine and top boundary node of the cluster (Figure 8(b)). Case (ii) says that if v and w are both leaf nodes in the same cluster then v is to the left of w iff v is to the left of w in the micro tree induced by that leaf cluster (Figure 8(c)). Case (iii) says that if v is a left node and w is a spine node in the same cluster then v is to the left of w iff v is to the left of w in the micro tree induced by their two macro nodes and the top boundary node of the cluster (Figure 8(d)). Case (iv) says that if v is a spine node and w is a right node in the same cluster then v is to the left of w iff v is to the left of w in the micro tree induced by their two macro nodes and the top boundary node of the cluster (Figure 8(e)). In all other cases v is to the left of w if the macro node v belongs to is to the left of the macro node of w in the macro tree (Case (v)).

### **Proposition 7 (Nca relations)** For any pair of nodes $v, w \in V(T)$ , the following hold

- (i) If c(v) = c(w) = l(v') for some boundary node v', then  $\operatorname{nca}_T(v, w) = \operatorname{nca}_{C(c(v), v')}(v, w)$ .
- (ii) If  $c(v) = c(w) \in \{l(v', w'), r(v', w')\}$  for some boundary nodes v', w', then  $nca_T(v, w) = nca_{C(c(v), s(v', w'), v')}(v, w)$ .
- (iii) If c(v) = c(w) = s(v', w') for some boundary nodes v', w', then  $\operatorname{nca}_T(v, w) = \operatorname{nca}_{C(c(v))}(v, w)$ .
- (iv) If  $c(v) \neq c(w)$  and  $c(v), c(w) \in \{l(v', w'), r(v', w'), s(v', w')\}$  for some boundary nodes v', w', then  $nca_T(v, w) = nca_{C(c(v), c(w), s(v', w'), v')}(v, w)$ .
- (v) If  $c(v) \neq c(w)$ ,  $c(v) \in \{l(v', w'), r(v', w'), s(v', w')\}$ , and  $w' \leq_M c(w)$  for some boundary nodes v', w', then  $\text{nca}_T(v, w) = \text{nca}_{C(c(v), s(v', w'), v', w')}(v, w')$ .
- (vi) In all other cases,  $nca_T(v, w) = nca_M(c(v), c(w))$ .

Case (i) says that if v and w are leaf nodes in the same cluster then the nearest common ancestor of v and w is the nearest common ancestor of v and w in the micro tree induced by that leaf cluster (Figure 8(c)). Case (ii) says that if v and w are both either left nodes or right nodes then the nearest common ancestor of v and w is the nearest common ancestor in the micro tree induced by their macro node together with the spine and top boundary node of the cluster (Figure 8(b)). Case (iii) says that if v and w are both spine nodes in the same cluster then the nearest common ancestor of v and w is the nearest common ancestor of v and w in the micro tree induced by their macro node. Case (iv) says that if v and w are in different macro nodes but are right, left, or spine nodes in the same cluster then the nearest common ancestor of v and w is the nearest common ancestor of v and w in the micro tree induced by that cluster (we can omit the bottom boundary node) (Figure 8(f)). Case (v) says that if v is a left, right, or spine node, and the bottom boundary node w of v is cluster is an ancestor in the macro tree of the macro node containing w, then the nearest common ancestor of v and w is the nearest common ancestor of v and w in the micro tree induced by the macro node of v, the spine node, and the top and bottom boundary nodes of v is cluster (Figure 8(g)). In all other cases the nearest common ancestor of v and w is the nearest common ancestor of their macro nodes in the macro tree (Case (vi)).

## 5.2 Preprocessing

In this section we describe how to preprocess T. First build a cluster partition CS of the tree T with clusters of size s, to be fixed later, and the corresponding macro tree M in  $O(n_T)$  time. The macro tree is preprocessed as in Section 4.1. However, since nodes in M contain a set of labels, we now store a dictionary for label(v) for each node  $v \in V(M)$ . Using the deterministic dictionary of Hagerup et al. [20] all these dictionaries can be constructed in  $O(n_T \log n_T)$  time and  $O(n_T)$  space. Furthermore, we extend the definition of fl such that  $fl_M(v,\alpha)$  is the nearest ancestor w of v such that  $\alpha \in label(w)$ .

Next we show how to preprocess the micro forests. For any cluster  $C \in CS$ , deep sets  $X, Y, Z \subseteq V(C)$ ,  $i \in \mathbb{N}$ , and  $\alpha \in \Sigma$  define the following procedures.

SIZE(X): Return the number of nodes in X.

LEFT(i, X): Return the leftmost i nodes in X.

RIGHT(i, X): Return the rightmost i nodes in X.

LEFTOF(X,Y): Return all nodes of X to the left of the leftmost node in Y.

RIGHTOF(X,Y): Return all nodes of X to the right of the rightmost node in Y.

MATCH(X, Y, Z), where  $X = \{m_1 \triangleleft \cdots \triangleleft m_k\}, Y = \{v_1 \triangleleft \cdots \triangleleft v_k\}, \text{ and } Z \subseteq Y. \text{ Return } R := \{m_j \mid v_j \in Z\}.$ 

MOP(X,Y) Return the pair  $(R_1, R_2)$ , where  $R_1 = \text{mop}(X,Y)|_1$  and  $R_2 = \text{mop}(X,Y)|_2$ .

If we want to specify that a procedure applies to a certain cluster C we add the subscript C. In addition to these procedures we also define the set procedures on clusters, that is, PARENT, NCA, DEEP, and FL, as in Section 3. Collectively, we will call these the *cluster procedures*. We represent the input and output sets in the procedures as bit strings indexed by preorder numbers. Specifically, a subset X in a cluster C is given by a bit string  $b_1 \ldots b_s$ , such that  $b_i = 1$  iff the ith node in a preorder traversal of C is in X. If C contains fewer than s nodes we set the remaining bits to 0.

The procedure  $\operatorname{SIZE}(X)$  is the number of ones in the bit string. The procedure  $\operatorname{LEFT}(i,X)$  corresponds to setting all bits in X larger than the ith set bit to zero. Similarly,  $\operatorname{RIGHT}(i,X)$  corresponds to setting all bits smaller than the ith largest set bit to zero. Similarly, the procedures  $\operatorname{LEFTOF}(X,Y)$ ,  $\operatorname{RIGHTOF}(X,Y)$ ,  $\operatorname{MOP}(X,Y)$ , and  $\operatorname{MATCH}(X,Y,Z)$  only depend on the preorder of the nodes and thus only on the bit string and not any other information about the cluster.

Next we show how to implement the cluster procedures efficiently. We precompute the value of all procedures, except FL, for all possible inputs and clusters. By definition, these procedures do not depend on any specific labeling of the nodes in the cluster. Hence, it suffices to precompute the value for all rooted, ordered trees with at most s nodes. The total number of these is less than  $2^{2s}$  (consider e.g. an encoding using balanced parenthesis). Furthermore, the number of possible input sets is at most  $2^s$ . Since at most 3 sets are given as input to a cluster procedure, it follows that we can tabulate all solutions using less than  $2^{3s} \cdot 2^{2s} = 2^{5s}$  bits of memory. Hence, choosing  $s \leq 1/10 \log n$  we use  $O(2^{\frac{1}{2} \log n}) = O(\sqrt{n})$  bits. Using standard bit wise operations each solution is easily implemented in O(s) time giving a total time of  $O(\sqrt{n} \log n)$ .

Since the procedure FL depends on the alphabet, which may be of size  $n_T$ , we cannot efficiently apply the same trick as above. Instead define for any cluster  $C \in CS$ ,  $X \subseteq V(C)$ , and  $\alpha \in \Sigma$ :

ANCESTOR(X): Return the set  $\{x \mid x \text{ is an ancestor of a node in } X\}$ .

EQ<sub>C</sub>( $\alpha$ ): Return the set  $\{x \mid x \in V(C), \text{label}(x) = \alpha\}.$ 

Clearly, ANCESTOR can be implemented as above. For EQ<sub>C</sub> note that the total number of distinct labels in C is at most s. Hence, EQ<sub>C</sub> can be stored in a dictionary with at most s entries each of which is a bit string of length s. Thus, (using again the result of [20]) the total time to build all such dictionaries is  $O(n_T \log n_T)$ .

By the definition of FL we have that,

$$\operatorname{FL}_C(X, \alpha) = \operatorname{DEEP}_C(\operatorname{ANCESTOR}_C(X) \cap \operatorname{EQ}_C(\alpha)).$$

Since intersection can be implemented using a binary and-operation,  $FL_C(X, \alpha)$  can be computed in constant time. Later, we will also need to compute union of sets represented as bit strings and we note that this can be done using a binary or-operation.

To implement the set procedures in the following section we often need to "restrict" the cluster procedures to work on a subtree of a cluster. Specifically, for any set of macro nodes  $\{i_1,\ldots,i_k\}$  in the *same* cluster C (hence,  $k \leq 5$ ), we will replace the subscript C with  $C(i_1,\ldots,i_k)$ . For instance, PARENT $_{C(s(v,w),l(v,w))}(X) = \{parent(x) \mid x \in X \cap V(C(s(v,w),l(v,w)))\} \cap V(C(s(v,w),l(v,w))$ . To implement all restricted versions of the cluster procedures, we compute for each cluster  $C \in CS$  a bit string representing the set of nodes in each micro forest. Clearly, this can be done in  $O(n_T)$  time. Since there are at most 5 micro forests in each cluster it follows that we can compute any restricted version using an additional constant number of and-operations.

Note that the total preprocessing time and space is dominated by the construction of deterministic dictionaries which use  $O(n_T \log n_T)$  time and  $O(n_T)$  space.

## 5.3 Implementation of the Set Procedures

Using the preprocessing from the previous section we show how to implement the set procedures in sublinear time. First we define a compact representation of node sets. Let T be a tree with macro tree M. For simplicity, we identify nodes in M with a number almost equal to their preorder number, which we denote their macro tree number: All nodes nodes except spine and left nodes are identified with their preorder number + 1 if they have a left node as a child and with their preorder number otherwise, and left nodes are identified with their preorder number - 1. Hence, we swap the order of left and spine nodes in the macro tree numbering. We will explain the reason for using macro tree numbers below. Note that the macro tree numbers are the same as the preorder numbers would be if we had let l(v, w) and r(v, w) be children of v instead of children of s(v, w) in the definition of the macro tree.

Let  $S \subseteq V(T)$  be any subset of nodes of T. A micro-macro node array (abbreviated node array) X representing S is an array of size  $n_M$ . The *i*th entry, denoted X[i], represents the subset of nodes in C(i), that is,  $X[i] = V(C(i)) \cap S$ . The set X[i] is encoded using the same bit representation as in Section 5.2. By our choice of parameter in the clustering the space used for this representation is  $O(n_T/\log n_T)$ .

We can now explain the reason for using macro tree numbers to identify the nodes instead of preorder numbers. Consider a node array representing a deep set. If a left node and the corresponding spine node are both non-empty, then all nodes in the left node are to the left of the node in the spine node. Formally,

**Proposition 8** Consider a node array X representing a deep set X. For any pair of nodes  $v, w \in X$ , such that  $v \in X[i]$  and  $w \in X[j]$ ,  $i \neq j$ , we have

$$v \triangleleft w \Leftrightarrow i < j$$
.

*Proof.* By Proposition 6(v) the claim is true for  $i \triangleleft j$ . The remaining cases are i = l(v', w') and j = s(v', w') (Proposition 6 (iii)) and i = s(v', w') and j = r(v', w') (Proposition 6(iv)). In both cases i < j and it follows immediately that  $v \triangleleft w \Rightarrow i < j$ . For the other direction, it follows from the structure of the macro tree that in both cases either  $v \triangleleft w$  or  $w \triangleleft v$ . But  $\mathcal{X}$  is deep and thus  $v \triangleleft w$ .

Thus, by using macro tree numbers we encounter the nodes in X according to their preorder number in the original tree T. This simplifies the implementation of all the procedures except DEEP, since they all get deep sets as input.

We now present the detailed implementation of the set procedures on node arrays. As in Section 4 we assume that the input to all of the procedures, except DEEP, represent a deep set. Let X be a node array.

#### **Procedure** Parent(X)

```
1 Initialize an empty node array R of size n_M (R[i] := \emptyset \text{ for } i = 1, \dots n_M) and set i := 1.
 2 while i \leq n_M do
       while X[i] = \emptyset do i := i + 1.
 3
       There are three cases depending on the type of i:
 4
       case 1. i \in \{l(v, w), r(v, w)\}
 5
           Compute N := PARENT_{C(i,s(v,w),v)}(X[i]).
 6
           foreach j \in \{i, s(v, w), v\} do
 7
              R[j] := R[j] \cup (N \cap V(C(j))).
 8
           end
 9
10
       case 2. i = l(v)
11
           Compute N := PARENT_{C(i,v)}(X[i]).
12
           foreach j \in \{i, v\} do
13
              R[j] := R[j] \cup (N \cap V(C(j))).
14
           end
15
16
       case 3. i \notin \{l(v, w), r(v, w), l(v)\}
17
           Compute N := PARENT_{C(i)}(X[i]).
18
           if N \neq \emptyset then
19
               set R[i] := R[i] \cup N.
20
           else if j := parent_M(i) \neq \bot then
21
            set R[j] := R[j] \cup \{first(j)\}.
22
           end
23
       Set i := i + 1.
24
25 end
26 Return R.
```

Procedure Parent has three cases. Case 1 handles the fact that left or right nodes may have a node on a spine or the top boundary node as parent. Since no left or right nodes can have their parent outside their cluster there is no need to compute parents in the macro tree. Case 2 handles the fact that a leaf node may have the boundary node as parent. Since no leaf node can have its parent outside its cluster there is no need to compute parents in the macro tree. Case 3 handles boundary and spine nodes. In this case there is either a parent within the micro forest or we can use the macro tree to compute the parent of the root of the micro tree. Since the input to Parent is deep we only need to do one of the two things. If the computation of parent in the micro tree returns a nonempty set, this set is added to the output (line 18). Otherwise (the returned set is empty), we compute parent of i in the macro tree (line 19). If the computation of parent in the macro tree returns a node j, this will either be a spine node or a boundary node. To take care of the case where j is a spine node, we add the lowest node (first(j)) in j to the output (line 20). If j is a boundary node this is just j itself.

Implementation of NCA We now give the implementation of procedure NCA. The input to procedure NCA is two node arrays X and Y representing two subsets  $\mathcal{X}, \mathcal{Y} \subseteq V(T)$ ,  $|\mathcal{X}| = |\mathcal{Y}| = k$ . The output is a node array R representing the set  $\text{Deep}(\{\text{nca}(\mathcal{X}_i, \mathcal{Y}_i) \mid 1 \leq i \leq k\})$ , where  $\mathcal{X}_i$  and  $\mathcal{Y}_i$  is the ith element of  $\mathcal{X}$  and  $\mathcal{Y}$ , w.r.t. their preorder number in the tree, respectively. We also assume that we have  $\mathcal{X}_i \triangleleft \mathcal{Y}_i$  for all i (since NCA is always called on a set of minimum ordered pairs). Note, that  $\mathcal{X}_l$  and  $\mathcal{Y}_l$  can belong to different clusters/nodes in the macro tree, i.e., we might have  $\mathcal{X}_l \in X[i]$  and  $\mathcal{Y}_l \in Y[j]$  where  $i \neq j$ .

### Procedure NCA(X,Y)

```
1 Initialize an empty node array R of size n_M, set i := 1 and j := 1.
 2 while i \leq n_M and j \leq n_M do
        while X[i] = \emptyset do i := i + 1.
 3
        while Y[j] = \emptyset do j := j + 1.
 4
        Set n := \min(\text{SIZE}(X[i]), \text{SIZE}(Y[j])), X_i := \text{LEFT}(n, X[i]), \text{ and } Y_j := \text{LEFT}(n, Y[j]).
 5
        Compare i and j. There are two cases:
 6
        case 1. i = j.
 7
            Set
 8
                                         S := \begin{cases} C(i, v), & \text{if } i = l(v), \\ C(i, s(v, w), v), & \text{if } i \in \{l(v, w), r(v, w)\}. \end{cases}
            Compute N := \text{NCA}_S(X_i, Y_i).
 9
            foreach macro node h = c(s) where s \in V(S) do
10
             set R[h] := R[h] \cup (N \cap V(C(h))).
11
            \quad \text{end} \quad
12
13
        case 2. i \neq j.
14
            Compute h := NCA_M(i, j). There are two subcases:
15
            case (a) h is a boundary node
16
                Set R[h] := 1.
17
            case (b) h is a spine node s(v, w)
18
                 There are three subcases:
19
                 case i. i \in \{l(v, w), s(v, w)\}\ and j \in \{s(v, w), r(v, w)\}\ 
\mathbf{20}
                 Compute N := \text{NCA}_{C(i,j,s(v,w),v)}(X_i, Y_j).
21
22
                 case ii. i = l(v, w) and w \leq j
23
                 Compute N := \text{NCA}_{C(i,s(v,w),v,w)}(\text{RIGHT}(1,X_i),w).
24
                 case iii. j = r(v, w) and w \leq i
25
                 Compute N := NCA_{C(j,s(v,w),w,v)}(w, LEFT(1,Y_j)).
26
27
                 Set R[h] := R[h] \cup (N \cap V(C(h))) and R[v] := R[v] \cup (N \cap V(C(v))).
28
29
        Set X[i] := X[i] \setminus X_i and Y[j] := Y[j] \setminus Y_i.
30
31 end
32 Return Deep(R).
```

In the main loop of procedure NCA (line 2–27) we first find the next non-empty entries in the node arrays X[i] and Y[j] (line 3 and 4). We then compare the sizes of X[i] and Y[j] and construct two sets of equal sizes  $X_i$  and  $Y_j$  consisting of the min(SIZE(X[i]), SIZE(Y[j])) leftmost nodes from X[i] and Y[j] (line 5). In Section 5.4 we prove the following invariant on  $X_i$  and  $Y_j$ 

LEFT
$$(1, X_i) = \mathcal{X}_l$$
 and LEFT $(1, Y_i) = \mathcal{Y}_l$  for some  $l$ .

The procedure has two main cases.

- If i = j (Case 1) then i is either a leaf, left, or right node due to the invariant and the assumption on the input that  $\mathcal{X}_l \triangleleft \mathcal{Y}_l$  (for a formal proof see Section 5.4). If i is a leaf node the nearest common ancestors of all pairs in  $X_i$  and  $Y_j$  are in the leaf node or the boundary node. If i is a left or right node the nearest common ancestors of all the pairs are in i, on the spine, or in the top boundary node. In line 9 we compute NCA in the appropriate cluster depending on the type of i.
- If  $i \neq j$  (Case 2) we first compute the nearest common ancestor h of i and j in the macro tree (line

- 14). Due to the structure of the macro tree h is either a spine node or a boundary node (left, right, and leaf nodes have no descendants). If h is a boundary node all pairs in  $X_i$  and  $Y_j$  have the same nearest common ancestor, namely h (Case 2(a)). If h is a spine node there are three cases depending on the types of i and j.
  - In Case 2(b)i we have i = l(v, w) and  $j \in \{s(v, w), r(v, w)\}$  (see Figure 8(d) and (f)), or i = s(v, w) and j = r(v, w) (see Figure 8(e)). In this case we compute NCA in the cluster containing i, j, s(v, w), v.
  - In case 2(b)ii i is a left node l(v, w) and j is a (not necessarily proper) descendant of w (see Figure 8(g)). In this case we compute NCA on the rightmost node in  $X_i$  and w in the cluster containing i, v, w, s(v, w). We can restrict the computation to RIGHT(1,  $X_i$ ) because we always run DEEP on the output from NCA before using it in any other computation and all nearest common ancestors of the pairs in  $X_i$  and  $Y_j$  will be on the spine, and the deepest one will be the nearest common ancestor of the rightmost nodes in  $X_i$  and  $Y_j$  (see Section 5.4 for a formal proof).
  - Case 2(b)iii is similar to Case 2(b)ii.

In the end of the iteration we have computed the nearest common ancestors of all the pairs in  $X_i$  and  $Y_j$  and the nodes from these pairs are removed from X[i] and Y[j].

**Implementation of** DEEP The implementation of DEEP resembles the previous implementation, but takes advantage of the fact that the input list is in macro tree order.

#### Procedure DEEP(X)

```
1 Initialize an empty node array R of size n_M.
 2 Find the smallest j such that X[j] \neq \emptyset. If no such j exists stop. Set i := j + 1.
   while i \leq n_M do
 4
        while X[i] = \emptyset do i := i + 1.
        Compare i and i. There are three cases:
 5
        case 1. i \triangleleft i.
 6
             Set
                                           S := \begin{cases} C(j,v), & \text{if } j = l(v), \\ C(j,s(v,w),v), & \text{if } j \in \{l(v,w),r(v,w)\}, \\ C(j), & \text{otherwise.} \end{cases}
            Set R[j] := \text{DEEP}_S(X[j]).
 8
        case 2. j \prec i.
 9
             if j = s(v, w) and i = r(v, w) then
10
                 \text{compute } N := \mathsf{DEEP}_{C(r(v,w),s(v,w),v)}(X[i] \cup X[j]).
11
                 Set R[j] := X[j] \cap N.
12
             end
13
        case 3. i \prec j (can happen if i = s(v, w) and j = l(v, w)).
14
             Compute N := \text{DEEP}_{C(l(v,w),s(v,w),v)}(X[i] \cup X[j]).
15
             Set R[j] := X[j] \cap N, X[i] := X[i] \cap N.
16
        Set i := i and i := i + 1.
17
19 Set R[j] := \text{DEEP}_S(X[j]), where S is set as in Case 1.
20 Return R.
```

The procedure DEEP has three cases. In case 1 node i is to the right of our "potential output node" j. Since any node l that is a descendant of j must be to the left of i (l < i) it cannot appear later in the list

X than i. We can thus safely add  $DEEP_S(X[j])$  to R at this point. To ensure that the cluster we compute DEEP on is a tree we include the top boundary node if j is a leaf node and the top and spine node if j is a left or right node. We add the result to R and set i to be our new potential output node.

In case 2 node j is an ancestor of i and therefore no node from C(j) can be in the output list unless j is a spine node and i is the corresponding right node. If this is the case we compute DEEP of X[j] and X[i] in the cluster containing i and j and add the result for j to the output and set i to be our new potential output node.

In case 3 node i is an ancestor of j. This can only happen if j is a left node and i the corresponding spine node. We compute DEEP of X[j] and X[i] in the cluster containing i and j and add the result for j to the output. We restrict X[i] to the nodes both in X[i] and the result N of the DEEP computation, and let i be our potential output node. The results for X[i] cannot be added directly to the input since there might be nodes later in the input that are descendants of i. Since a left node has no children we can safely add the result for j to the output.

After iterating through the whole node array X we add the last potential node j to the output after computing DEEP of it as in Case 1.

Implementation of Mopricht We now give the implementation of procedure Mopricht. Procedure Mopricht takes a pair of node arrays (X, Y) and another node array Z as input. The pair (X, Y) represents a set of minimum ordered pairs, where the first coordinates are in X and the second coordinates are in Y. To simplify the implementation of procedure Mopricht it calls two auxiliary procedures Mopricht and Match defined below. Procedure Mopricht computes mop of Y and Z, and procedure Match computes the first coordinates from X corresponding to the first coordinates from the minimum ordered pairs of Y and Z computed by Mopricht.

## **Procedure** MopRight((X,Y),Z)

- 1 Compute M := MopSim(Y, Z).
- **2** Compute  $R := MATCH(X, Y, M|_1)$ .
- **3** Return  $(R, M|_2)$ .

Procedure MopSim takes two node arrays as input and computes mop of these.

### **Procedure** MopSim(X,Y)

```
1 Initialize two empty node arrays R and S of size n_M.
 2 Set i := 1, j := 1, (r_1, r_2) := (0, \emptyset), (s_1, s_2) := (0, \emptyset).
 з repeat
        while X[i] = \emptyset do set i := i + 1.
 4
        There are four cases:
 5
        case I. i = l(v, w) for some v, w.
 6
           Until Y[j] \neq \emptyset and either i \triangleleft j, i = j, or j = s(v, w) do set j := j + 1.
 7
        case II. i = s(v, w) for some v, w.
 8
           Until Y[j] \neq \emptyset and either i \triangleleft j or j = r(v, w) do set j := j + 1.
 9
        case III. i \in \{r(v, w), l(v)\}\ for some v, w.
10
        Until Y[j] \neq \emptyset and either i \triangleleft j or i = j do set j := j + 1.
11
        case IV. i is a boundary node.
12
           Until Y[j] \neq \emptyset and i \triangleleft j do set j := j + 1.
13
14
        Compare i and j. There are two cases:
15
        case 1. i \triangleleft j.
16
            if s_1 < j then
17
            set R[r_1] := R[r_1] \cup r_2, S[s_1] := S[s_1] \cup s_2, and (s_1, s_2) := (j, \text{LEFT}_{C(j)}(1, Y[j])).
18
19
            Set (r_1, r_2) := (i, RIGHT_{C(i)}(1, X[i])) and i = i + 1.
20
21
        otherwise // case 2.
22
            Compute (r,s) := MOP_{C(i,j,v)}(X[i],Y[j]), where v is the top boundary node in the cluster i
23
            and j belong to.
            if r \neq \emptyset then
24
                if s_1 < j or if s_1 = j and LEFTOF_{C(i,j)}(X[i], s_2) = \emptyset then
25
                 set R[r_1] := R[r_1] \cup r_2, S[s_1] := S[s_1] \cup s_2.
26
27
                Set (r_1, r_2) := (i, r) and (s_1, s_2) := (j, s).
28
            end
29
            There are two subcases:
30
            case (a) i = j, or i = l(v, w) and j = s(v, w).
31
             Set X[i] := RIGHT_{C(i)}(1, RIGHTOF_{C(i)}(X[i], r)) and j := j + 1.
32
            case (b) i = s(v, w) \ and \ j = r(v, w).
33
               if r = \emptyset then set j := j + 1 else set i := j.
34
35
       endsw
36
37 until i > n_M or j > n_M;
38 Set R[r_1] := R[r_1] \cup r_2 and S[s_1] := S[s_1] \cup s_2.
39 Return (R, S).
```

Procedure MopSim is somewhat similar to the previous implementation of the procedure MopRight from Section 4.2. As in the previous implementation we have a "potential pair"  $((r_1, r_2), (s_1, s_2))$ , where  $r_1$  and  $s_1$  are macro nodes,  $r_2 \subseteq X[r_1]$ ,  $s_2 \subseteq Y[s_1]$ , where  $r_2 = \{r^1 \lhd \cdots \lhd r^k\}$  and  $s_2 = \{s^1 \lhd \cdots \lhd s^k\}$  such that  $r^l \lhd s^l$  for  $l = 1, \ldots k$ . Furthermore, for any l there exists no node  $y \in Y[j]$ , for  $j < s_1$ , such that  $r^l \lhd y \lhd s^l$  and no node  $x \in X[i]$ , for  $i < r_1$ , such that  $r^l \lhd x \lhd s^l$ .

We have the following invariant at the beginning of each iteration:

$$\nexists x \in X[i], \text{ such that } x \le x', \text{ for any } x' \in r_2.$$
(8)

We first find the next non-empty macro node i. We then have 4 cases depending on which kind of node

i is.

- In Case I i is a left node. Due to Proposition 6 we can have mop in i (case (i), see Fig. 8(b)), in the spine (case (iii), see Fig. 8(d)), or in a node to the right of i (case(v)).
- In Case II i is a spine node. Due to Proposition 6 we can have mop in the right node (case (iv), see Fig. 8(e)) or in a node to the right of i (case(v)).
- In Case III i is a right node or a leaf node. Due to Proposition 6 we can have mop in i (case (i) and (ii), see Fig. 8(b)-(c)) or in a node to the right of i (case(v)).
- In the last case (Case IV) i must be a boundary node and mop must be in a node to the right of i.

We then compare i and j. The case where  $i \triangleleft j$  is similar to the previous implementation of the procedure. We compare j with our potential pair (line 16). If  $s_1 < j$  then  $s_1 \triangleleft j$  since the input is deep, and we can insert  $r_2$  and  $s_2$  into our output node arrays R and S, respectively. We also set  $s_1$  to j and  $s_2$  to the leftmost node in Y[j] (if  $s_1 = j$  we already have  $(s_1, s_2) = (j, \text{LEFT}_{C(j)}(1, Y[j]))$ ). Then—both if  $s_1 \triangleleft j$  or  $s_1 = j$ —we set  $r_1$  to i and  $r_2$  to the rightmost node in X[i] (line 19). That we only need the rightmost node in X[i] and the leftmost node in Y[j] follows from the definition of mop and the structure of the macro tree.

Case 2  $(i \not\preceq j)$  is more complicated. In this case we first compute mop in the cluster i and j belong to (line 21). If this results in any minimum ordered pairs  $(r \neq \emptyset)$  we must update our potential pair (line 22–27). Otherwise we leave the potential pair as it is and only update i and j. If  $r \neq \emptyset$  we compare  $s_1$  and j (line 23). As in Case 1 of the procedure we add our potential pair to the output and update the potential pair with r and s if  $s_1 < j$ , since this implies  $s_1 \lhd j$ . If  $s_1 = j$  and no nodes in X[i] are to the left of the leftmost node in  $s_2$  we also add the potential pair to the output and update it. We show in the next section that in this case  $|s_2| = 1$ . Therefore we can safely add the potential pair to the output. In all other cases the pair  $(r, s) \neq (\emptyset, \emptyset)$  shows a contradiction to our potential pair and we update the potential pair without adding anything to the output.

Finally, in Case 2, we update X[i], i, and j (line 28–32). There are two cases depending on i and j. In Case (a) either i=j or i is a left node and j is the corresponding spine node. In both cases we can have nodes in X[i] that are not to the left of any node in Y[j]. These nodes could be in a minimum ordered pair with nodes from another macro node. We show in the next section that this can only be true for the rightmost node in X[i]. X[i] is updated accordingly. After this update all nodes in Y[j] are to the left of all nodes in X[i] in the next iteration and therefore j is incremented. In Case (b) i is a spine node and j is the corresponding right node. Since the input lists are deep, there is only one node in X[i]. If  $r = \emptyset$  then no node in Y[j] is to the right of the single node in X[i]. Since the input arrays are deep, no node later in the array X can be to the left of any node in Y[j] and we therefore increment j. If  $r \neq \emptyset$  then  $(r_1, r_2) = (i, X[i])$  and we update i. Instead of incrementing i by one we set i := j, this is correct since all macro nodes with macro node number between i and j are descendants of i, and thus contains no nodes from X, since X is deep.

When reaching the end of one of the arrays we add our potential pair to the output and return (line 35–36).

As in Section 4.2 we can implement MopLeft similarly to MopRight.

Recall that procedure Moprish calls Match to find the first coordinates from X corresponding to the first coordinates from the minimum ordered pairs computed by Moprish. Procedure Match takes three node arrays X, Y, and Y' representing deep sets  $\mathcal{X}$ ,  $\mathcal{Y}$ , and  $\mathcal{Y}'$ , where  $|\mathcal{X}| = |\mathcal{Y}|$ , and  $\mathcal{Y}' \subseteq \mathcal{Y}$ . The output is a node array representing the set  $\{\mathcal{X}_j \mid \mathcal{Y}_j \in \mathcal{Y}'\}$ .

## **Procedure** MATCH(X,Y,Y')

```
1 Initialize an empty node array R of size n_M.
 2 Set X_L := \emptyset, Y_L := \emptyset, Y_L' := \emptyset, x := 0, y := 0, i := 1 and j := 1.
 з repeat
        while X[i] = \emptyset do set i := i + 1.
 4
        while Y[j] = \emptyset do set j := j + 1.
 5
        Set x := SIZE(X[i]) and y := SIZE(Y[j]).
 6
        Compare Y[j] and Y'[j]. There are two cases:
 7
        case 1. Y[j] = Y'[j]
 8
             Compare x and y. There are three subcases:
 9
             case (a) x = y.
10
             Set R[i] := R[i] \cup X[i], i := i + 1, and j := j + 1.
11
12
             case (b) x < y.
13
             Set R[i] := R[i] \cup X[i], Y[j] := RIGHT(y - x, Y[j]), Y'[j] := Y[j], and i := i + 1.
14
15
             case (c) x > y.
16
              Set X_L := \text{LEFT}(y, X[i]), R[i] := R[i] \cup X_L, X[i] := X[i] \setminus X_L, \text{ and } j := j + 1.
17
18
        case 2. Y[j] \neq Y'[j]
19
            Compare x and y. There are three subcases:
20
             case (a) x = y.
21
             Set R[i] := R[i] \cup MATCH(X[i], Y[j], Y'[j]), i := i + 1, and j := j + 1.
22
23
             case (b) x < y.
24
                Set Y_L := \text{LEFT}(x, Y[j]), Y'_L := Y'[j] \cap Y_L, R[i] := R[i] \cup \text{MATCH}(X[i], Y_L, Y'_L), Y[j] := Y[j] \setminus Y_L, Y'[j] := Y'[j] \setminus Y'_L, \text{ and } i := i + 1.
25
26
27
             case (c) x > y.
28
                Set X_L := \text{LEFT}(y, X[i]), R[i] := R[i] \cup \text{MATCH}(X_L, Y[j], Y'[j]), X[i] := X[i] \setminus X_L, and j := j + 1.
29
30
31
32
зз until i > n_M or j > n_M;
34 Return R.
```

Procedure MATCH proceeds as follows. First we find the first non-empty entries in the two node arrays X[i] and Y[j] (line 4–5). We then compare Y[j] and Y'[j] (line 7).

If they are equal we keep all nodes in X with the same rank as the nodes in Y[j] (case 1). We do this by splitting into three cases. If there are the same number of nodes X[i] and Y[j] we add all nodes in X[i] to the output and increment i and j (case 1(a)). If there are more nodes in Y[j] than in X[i] we add all nodes in X[i] to the output and update Y[j] and Y'[j] to contain only the y-x leftmost nodes in Y[j] (case 1(b)). We then increment i and iterate. If there are more nodes in X[i] than in Y[j] we add the first y nodes in X[i] to the output, increment j, and update X[i] to contain only the nodes we did not add to the output (case 1(c)).

If  $Y[j] \neq Y'[j]$  we call the cluster procedure MATCH (case 2). Again we split into three cases depending on the number of nodes in X[i] and Y[j]. If they have the same number of nodes we can just call MATCH on X[i], Y[j], and Y'[j] and increment i and j (case 2(a)). If SIZE(Y[j]) > SIZE(X[i]) we call match with X[i] the leftmost SIZE(X[i]) nodes of Y[j] and with the part of Y'[j] that are a subset of these leftmost SIZE(X[i]) nodes of Y[j] (case 2(b)). We then update Y[j] and Y'[j] to contain only the nodes we did not use in the call to MATCH and increment i. If SIZE(Y[j]) < SIZE(X[i]) we call MATCH with the leftmost SIZE(Y[j]) nodes

of X[i], Y[j], and Y'[j] (case 2(c)). We then update X[i] to contain only the nodes we did not use in the call to MATCH and increment j.

Implementation of FL Procedure FL takes as input a node array X representing a deep set and a label  $\alpha$ .

### **Procedure** $FL(X, \alpha)$

```
1 q Initialize an empty node array R of size n_M and two node lists L and S.
         while X[i] = \emptyset do set i := i + 1.
 3
         There are three cases depending on the type of i:
 4
         case 1. i \in \{l(v, w), r(v, w)\}
 5
             Compute N := \operatorname{FL}_{C(i,s(v,w),v)}(X[i],\alpha).
 6
             if N \neq \emptyset then
 7
                 foreach j \in \{i, s(v, w), v\} do set R[j] = R[j] \cup (N \cap V(C(j))).
 8
             else set L := L \circ \operatorname{parent}_{M}(v).
 9
         case 2. i = l(v)
10
             Compute N := \operatorname{FL}_{C(i,v)}(X[i], \alpha).
11
             if N \neq \emptyset then
12
                 foreach j \in \{i, v\} do set R[j] := R[j] \cup (N \cap V(C(j))).
13
             else set L := L \circ \operatorname{parent}_{M}(v).
14
         case 3. i \notin \{l(v, w), r(v, w), l(v)\}
15
             Compute N := \operatorname{FL}_{C(i)}(X[i], \alpha).
16
             if N \neq \emptyset then
17
                 set R[i] := R[i] \cup N.
18
             else set L := L \circ \operatorname{parent}_{M}(i).
19
20
21 until i > n_M;
22 Compute the list S := \operatorname{FL}_M(L, \alpha).
23 foreach node i \in S do set R[i] := R[i] \cup FL_{C(i)}(first(i), \alpha).
24 Return Deep(R)
```

The FL procedure is similar to PARENT. The cases 1, 2 and 3 compute FL on a micro forest. If the result is within the micro tree we add it to R and otherwise we store in a node list L the node in the macro tree which contains the parent of the root of the micro forest. Since we always call DEEP on the output from  $FL(X,\alpha)$  there is no need to compute FL in the macro tree if N is nonempty. We then compute FL in the macro tree on the list L, store the results in a list S, and use this to compute the final result.

Consider the cases of procedure FL. In case 1 i is a left or right node. Due to Proposition 5 case (i) and (ii) fl of a node in i can be in i or on the spine or in the top boundary node. If this is not the case it can be found by a computation of FL of the parent of the top boundary node of i's cluster in the macro tree (Proposition 5 case (iii)). In case 2 i is a leaf node. Then fl of a node in i must either be in i, in the top boundary node, or can be found by a computation of FL of the parent of the top boundary node of i's cluster in the macro tree. If i is a spine node or a boundary node (case 3), then fl of a node in i is either in i or can be found by a computation of FL of the parent of i in the macro tree.

#### 5.4 Correctness of the Set Procedures

The following lemmas show that the set procedures are correctly implemented.

Lemma 18 Procedure Parent is correctly implemented.

*Proof.* We will prove that in iteration i the procedure correctly computes the parents of all nodes in the macro node i. There are four cases depending on the type of i.

- Consider the case  $i \in \{l(v, w), r(v, w)\}$ , i.e., i is a left or right node. For all nodes x in C(i), parent(x) is either in C(i), on the spine s(v, w), or is the top boundary node v. The parents of all input nodes in C(i) is thus in N computed in Case 1 in the procedure. The last line in Case 1 ("For each  $j \in \{i, s(v, w), v\}, \ldots$ ") adds the set of parents to the appropriate macro node in the output array.
- If i is a leaf node l(v) then for any node  $x \in C(i)$ , parent(x) is either in C(i) or is the boundary node v. The parents of all input nodes in C(i) is thus in N computed in Case 2 in the procedure. The last line in Case 2 ("For each  $j \in \{i, v\}, \ldots$ ") adds the set of parents to the appropriate macro node in the output array.
- If i is a spine node s(v, w) then the input contains at most one node in C(i), since the input to the procedure is deep. For any  $x \in C(i)$ , parent(x) is either a node on the spine or the top boundary node v. This is handled by Case 3 in the procedure. Let x be the node in X[i]. If parent(x) = v, then  $N = \emptyset$ , and we compute j, which is the parent v of i in the macro tree, and add j to the output array (since j = v is a boundary node first(j) = v). If parent(x) is another node y on the spine, then  $N = \{y\} \neq \emptyset$  and y is added to the output array.
- If i is a boundary node v, then parent(v) is either another boundary node v', the bottom node on a spine, or  $\bot$  if v is the root. This is handled by Case 3 in the procedure. In all three cases  $N = \emptyset$  and we compute the parent j of i in the macro tree. If i is the root, then  $j = \bot$  and we do nothing. Otherwise, we add first(j) to the output. If parent(v) is a boundary node then first(j) is the bottom node on j.

In each iteration of the procedure we might add nodes to the output, but we never delete anything written to the output in earlier iterations. Procedure Parent thus correctly computes the parents of all nodes in X.

Before proving the correctness of procedure NCA we will prove the following invariant on the variables  $X_i$  and  $Y_j$  in the procedure.

**Lemma 19** In procedure NCA we have the following invariant of  $X_i$  and  $Y_j$ :

LEFT
$$(1, X_i) = \mathcal{X}_l$$
 and LEFT $(1, Y_i) = \mathcal{Y}_l$  for some  $l$ .

*Proof.* The proof is by induction on the number of iterations of the outer loop. After the while loop on X in the first iteration (line 3), i is the smallest integer such that  $X[i] \neq \emptyset$ . Due to the macro tree order of the array X, X[i] contains the first nodes from X w.r.t. the preorder of the original tree (Proposition 8). Similarly, Y[j] contains the leftmost node in  $\mathcal{Y}$ . The invariant now follows immediately from the assignment of  $X_i$  and  $Y_j$ .

For the induction step consider iteration m and let i' and j' be the values of i and j after the while loops in the previous iteration, i.e, after line 4. By the induction hypothesis  $\text{LEFT}(1, X_{i'}) = \mathcal{X}_l$  and  $\text{LEFT}(1, Y_{j'}) = \mathcal{Y}_l$  for some l. Let  $n' = \min(\text{SIZE}(X[i']), \text{SIZE}(Y[j']))$ . Then  $X_{i'}$  contains  $\mathcal{X}_l, \ldots, \mathcal{X}_{l+n'}$  and  $Y_{j'}$  contains  $\mathcal{Y}_l, \ldots, \mathcal{Y}_{l+n'}$ . We will show that  $\text{LEFT}(1, X_i) = \mathcal{X}_{l+n'+1}$ . In the end of the previous iteration we removed  $X_{i'}$  from X[i'] (line 26). There are two cases depending on wether X[i'] is empty or not at the beginning of iteration m.

- If  $X[i'] \neq \emptyset$  then it clearly contains  $\mathcal{X}_{l+n'+1}$  as its leftmost node. Since a spine node can only contain one node from  $\mathcal{X}$ , i' cannot be a spine node. Thus i = i', when we get to line 5 in the current iteration It follows that LEFT $(1, X_i) = \mathcal{X}_{l+n'+1}$ .
- $X[i'] = \emptyset$ . It follows from the macro tree order of X that X[i] contains  $\mathcal{X}_{l+n'+1}$  as its leftmost node.

It follows by a similar argument that LEFT $(1, Y_j) = \mathcal{Y}_{l+n'+1}$ .

**Lemma 20** Let X and Y be two node arrays representing the deep sets X and Y, |X| = |Y| = k, and let  $X_i$  and  $Y_i$  denote the ith element of X and Y, w.r.t. their preorder number in the tree, respectively. For all i = 1, ..., k, assume  $X_i \triangleleft Y_i$ . Procedure NCA(X, Y) correctly computes  $DEEP(\{nca(X_i, Y_i) | 1 \le i \le k\})$ .

*Proof.* We are now ready to show that the procedure correctly takes care of all possible cases from Proposition 7. The proof is split into two parts. First we will argue that some of the cases from the proposition cannot occur during an iteration of the outer loop of NCA. Afterwards we prove that the procedure takes care of all the cases that can occur.

Case (iii) cannot happen since if i = j is a spine node then  $\mathcal{X}_l$  is either a descendant or an ancestor of  $\mathcal{Y}_l$  contradicting the assumption on the input that  $\mathcal{X}_l \lhd \mathcal{Y}_l$ . Case (vi) can only happen if  $i \neq j$ : If i = j and we are in case (vi) then i = j is a boundary node, and this would imply that C(i) only consists of one node, i.e.,  $\mathcal{X}_l = X[i] = Y[j] = \mathcal{Y}_l$  contradicting the assumption on the input that  $\mathcal{X}_l \lhd \mathcal{Y}_l$ . Due to this assumption on the input we also have that in case (iv) of the proposition i is either a left node or a spine node and j is a spine node or a right node. For case (v) either i is a left node and j is a descendant of the bottom boundary node of j's cluster. All the other cases from case (v) would contradict the assumption that  $\mathcal{X}_l \lhd \mathcal{Y}_l$ .

The procedure first constructs two sets  $X_i$  and  $Y_j$  containing the elements  $\mathcal{X}_l, \ldots, \mathcal{X}_{l+n}$  and  $\mathcal{Y}_l, \ldots, \mathcal{Y}_{l+n}$  for some l, respectively, where  $n = \min(\text{SIZE}(X[i]), \text{SIZE}(Y[j]))$ . The procedure NCA has two main cases depending on whether i = j or not. Case 1 (i = j) takes care of cases (i)–(ii) from Proposition 7. Case 2  $(i \neq j)$  takes care of the remaining cases from Proposition 7 (iv)–(vi) that can occur.

First consider Case 1: We compute nearest common ancestors N of the n nodes in  $X_i$  and  $Y_j$  in a cluster S depending on what kind of node i is. We need to show that Case 1 handles Case (i) and (ii) from the Proposition correctly.

- Case (i). i = j is a leaf node. By the Proposition the nearest common ancestors of the pairs in  $(\mathcal{X}_l, \mathcal{Y}_l), \ldots, (\mathcal{X}_{l+n}, \mathcal{Y}_{l+n})$  from  $X_i$  and  $Y_j$  is either in c(i) or in the boundary node, i.e., in C(i, v).
- Case (ii). i = j is a left or right node. By the Proposition the nearest common ancestors of the pairs in  $\{(\mathcal{X}_l, \mathcal{Y}_l), \dots, (\mathcal{X}_{l+n}, \mathcal{Y}_{l+n})\}$  from  $X_i$  and  $Y_j$  is either in c(i), on the spine, or in the top boundary node, i.e., in C(i, s(v, w), v).

Thus S is correctly set in both cases. After the computation of N in line 9 the output is then added to the entries in the output array R for each of the macro nodes belonging to nodes in V(S) (line 10–12). Case 1 thus handles Case (i)-(ii) (and only these two cases) from Proposition 7.

Next consider Case 2  $(i \neq j)$ . We first compute the nearest common ancestor h of i and j in the macro tree. The macro node h is either a boundary node or a spine node due to the structure of the macro tree (see also Proposition 7). We will show that Case 2 takes care of the remaining cases.

- Case (iv). From the above discussion it follows that we have one of the three following cases. i = l(v, w) and j = s(v, w), i = l(v, w) and j = r(v, w), or i = s(v, w) and j = r(v, w). All three cases are handled in Case 2(b)i of the procedure. It follows from the proposition that NCA is computed in the correct cluster.
- Case (v). It follows from the discussion above that either i = l(v, w) and  $w \leq j$ , or j = r(v, w) and  $w \leq i$ . These two cases are handled by Case 2(b)ii and 2(b)iii of the procedure. It follows from the Proposition that NCA is computed in the correct cluster. We need to argue that we can restrict the computation of NCA to the pair (RIGHT(1,  $X_i$ ), w) instead of computing NCA for all nodes in  $\{\mathcal{X}_l, \ldots, \mathcal{X}_{l+n}\}$ . Consider the case where i = l(v, w) and  $w \leq j$  (Case 2(b)ii of the procedure). Since  $w \leq \mathcal{Y}_r$  for all  $r = l, \ldots l + n$ , and  $\mathcal{X}_l \lhd \mathcal{X}_{l+1} \lhd \ldots \lhd \mathcal{X}_{l+n}$ , then  $\operatorname{nca}(\mathcal{X}_r, \mathcal{Y}_r) \leq \operatorname{nca}(\mathcal{X}_{l+n}, \mathcal{Y}_{l+n})$  for all  $r = l, \ldots l + n$ . Thus we do not need to compute  $\operatorname{nca}(\mathcal{X}_r, \mathcal{Y}_r)$  for  $r \neq n + l$ , since the output of the procedure is  $\operatorname{DEEP}(\{\operatorname{nca}(\mathcal{X}_i, \mathcal{Y}_i) | 1 \leq i \leq k\})$ . A similar argument shows that we can restrict the computation to  $(w, \operatorname{LEFT}(1, Y_i))$  in Case 2(b)iii.

• Case (vi). It follows from the discussion above and the proposition that  $i \neq j$  and i and j are in different clusters, and we are not in any of the cases from (iv) and (v). Thus h must be a boundary node and all the pairs  $\{(\mathcal{X}_l, \mathcal{Y}_l), \ldots, (\mathcal{X}_{l+n}, \mathcal{Y}_{l+n})\}$  have the same nearest common ancestor, namely h. This is handled by Case 2(a).

We have now argued that the procedure correctly takes care of all possible cases from Proposition 7. It remains to show that all pairs from  $\{\operatorname{nca}(\mathcal{X}_i,\mathcal{Y}_i)|1\leq i\leq k\}$  are considered during the computation. It follows from the invariant that we only consider pairs from the input. In the last lines we remove the nodes from the input that we have computed the notes of in this iteration. It follows from the proof of the invariant that no entry in the input arrays is left nonempty. Thus all pairs are taken care of.

To prove that procedure DEEP is correctly implemented we will use the following fact about preorder and postorder numbers in the macro tree.

**Proposition 9** Let i and j be nodes in the macro tree identified by their macro tree number such that i < j. For all  $x \in C(i), y \in C(j)$  we have

```
1. \operatorname{pre}(x) < \operatorname{pre}(y) unless i = l(v, w) and j = s(v, w).
```

2. post(y) > post(x) unless i = s(v, w) and j = r(v, w).

**Proposition 10** Let  $x_1, ..., x_n$  be nodes from the macro tree associated with their macro tree number such that  $x_1 < x_2 < \cdots < x_n$ . If  $x_i \triangleleft x_j$  for some i and j then  $x_i \triangleleft x_k$  for all  $x_k > x_j$ .

*Proof.* From  $x_i \triangleleft x_j$  we have  $\operatorname{pre}(x_i) \triangleleft \operatorname{pre}(x_j)$  and  $\operatorname{post}(x_i) \triangleleft \operatorname{post}(x_j)$ . Since  $x_k > x_j$  we have  $\operatorname{pre}(x_j) \triangleleft \operatorname{pre}(x_k)$  unless  $x_k = s(v, w)$  and  $x_j = l(v, w)$ . In that case,  $\operatorname{pre}(x_k) + 1 = \operatorname{pre}(x_j) > \operatorname{pre}(x_i)$ . Since  $x_i \triangleleft x_j$  we have  $x_i \neq x_j$  and thus  $\operatorname{pre}(x_k) > \operatorname{pre}(x_i)$ .

It remains to show that  $post(x_i) < post(x_k)$ . Assume for the sake of contradiction that  $post(x_k) < post(x_i) < post(x_i)$ . This implies  $x_i \prec x_k$  and  $x_j \prec x_k$  contradicting  $x_i \lhd x_j$ .

We will first prove the following invariants on i and j in procedure DEEP.

**Lemma 21** In line 5 of procedure DEEP we have the following invariant on i and j: For all l such that j < l < i we have  $X[l] = \emptyset$ .

*Proof.* Let i' be the value of i in line 5 of the previous iteration of the outer loop (line 3–18). Then i is the smallest index greater than i' such that the corresponding entry in X is nonempty. This is true since i was set to i' + 1 in the end of the previous iteration (line 17), and in line 4 of this iteration i was incremented until we found a nonempty entry. Since j = i' (this was also set in line 17 of the previous iteration), i is the first nonempty entry greater than j and the claim follows.

**Lemma 22** At the beginning of each iteration of the main loop of procedure DEEP (line 3) we have the following invariant on j: For all nodes  $x \in X[j]$  and  $y \in X[l]$ , where  $1 \le l < j$ , we have  $x \not\prec y$ .

Proof. Recall that  $x \prec y \Leftrightarrow \operatorname{pre}(x) < \operatorname{pre}(y)$  and  $\operatorname{post}(y) < \operatorname{post}(x)$ . By Proposition 9 the only case where we can have  $\operatorname{pre}(x) < \operatorname{pre}(y)$  is if l = l(v, w) and j = s(v, w) for some v, w. Assume this is the case. If  $X[l] = \emptyset$  the claim follows trivially. Otherwise, let i' and j' be the values of i and j in the previous iteration, respectively (since l < j and  $X[l] \neq \emptyset$  there must be such an iteration). We have j = l + 1, i' = j = s(v, w) and j' = l = l(v, w). Thus in the previous iteration the procedure entered case 3, where X[i'] was set to  $X[i'] \cap \operatorname{DEEP}_{C(l(v,w),s(v,w),v)}(X[i'] \cup X[j'])$ , and thus X[j] contains no nodes that are ancestors of nodes in X[j'] = X[l].

*Proof.* We will prove that  $x \in \text{DEEP}(X)$  iff  $x \in X$  and there exists no  $y \in X$  such that  $x \prec y$ .

Assume  $x \in \text{DEEP}(X)$ . Consider the iteration when x is assigned to the output. There are three cases depending on which case we are in when x is added to the input. If  $j \triangleleft i$  (Case 1 of the procedure) then  $x \in \text{DEEP}_S(X[j])$  and it follows from the invariant on j (Lemma 22) that x has no descendants in any nodes  $y \in X[l]$ , l < j. For j < l < i the claim follows directly from Lemma 21. It remains to show that x has no descendants in X[l] for  $l \ge i$ . By Proposition 10 we have  $j \triangleleft l$  for all l > i and the claim follows from Proposition 5.

If  $j \prec i$  (Case 2 of the procedure) then j is a spine node s(v,w) and i is the corresponding right node r(v,w), and we compute  $N := \text{DEEP}_{C(r(v,w),s(v,w),v)}(X[i] \cup X[j])$ . Since  $x \in \text{DEEP}(X)$  we have  $x \in R[j] = X[j] \cap N$ . It follows from the invariant (Lemma 22) and the computation of N that x has no descendants in X[l] for any  $l \leq j$ . For l > j it follows from the structure of the macro tree that for any l > i we have  $j \lhd l$ . For j < l < i the claim follows directly from Lemma 21. The claim follows from Proposition 5. For j < l < i the claim follows directly from Lemma 21.

If  $i \prec j$  (Case 3 of the procedure) then i is a spine node s(v, w) and j is the corresponding left node l(v, w), and we compute  $N := \text{DEEP}_{C(l(v,w),s(v,w),v)}(X[i] \cup X[j])$ . Since  $x \in \text{DEEP}(X)$  we have  $x \in R[j] = X[j] \cap N$ . It follows from the computation of N that x has no descendants in  $X[i] \cup X[j]$ . Since l(v, w) has no descendants in the macro tree it follows from Proposition 5 that x has no descendants in X[l] for any  $l \neq j$ .

If x is assigned to the output in line 19 then it follows from the invariant on j (Lemma 22) and the computation of  $\text{DEEP}_S(X[j])$  that x has no descendants in X.

For the other direction let  $x \in X$  be a node such that  $X \cap V(T(x)) = \{x\}$ . Let l be the index such that  $x \in X[l]$ . All nonempty entries in X are i in line 5 at some iteration. Consider the iteration when i = l. Unless i = l(v, w) and j = s(v, w) (Case 3 of the procedure) X[i] is not changed in this iteration. If we are in Case 3, then N is computed and X[i] is set to  $X[i] \cap N$ . Since x has no descendants in X we have  $x \in N$  and thus  $x \in X[i]$  after the assignment. At the end of this iteration j is set to i. Consider the next iteration when j = l. If  $j \triangleleft i$  or  $i > n_M$  then  $x \in DEEP_S(X[j]) = R[j]$ . If  $j \triangleleft i$  we have j = s(v, w) and i = r(v, w) since x has no descendants in x. For the same reason we have  $x \in N$  and thus  $x \in X[j] \cap N = R[j]$ . If  $i \triangleleft j$  we have i = s(v, w) and j = l(v, w). Again  $x \in N$  and thus  $x \in X[j] \cap N = R[j]$ .

We now consider procedures MopSim and Match.

**Lemma 24** Let  $((r_1, r_2), (s_1, s_2))$  be as defined in procedure MOPSIM. Then  $r_1$  and  $s_1$  are macro nodes,  $r_2 \subseteq X[r_1]$ ,  $s_2 \subseteq Y[s_1]$ , where  $r_2 = \{r^1 \triangleleft \cdots \triangleleft r^k\}$  and  $s_2 = \{s^1 \triangleleft \cdots \triangleleft s^k\}$ . For any  $l = 1, \ldots, k$  we have

- 1.  $r^l \triangleleft s^l$ .
- 2. for all  $j \leq s_1$  there exists no node  $y \in Y[j]$  such that  $r^l \triangleleft y \triangleleft s^l$ ,
- 3. for all  $i \leq r_1$  there exists no node  $x \in X[i]$  such that  $r^l \triangleleft x \triangleleft s^l$ .

Proof. It follows immediately from the code that  $r_1$  and  $s_1$  are macro nodes and that  $r_2 \subseteq X[r_1]$ ,  $s_2 \subseteq Y[s_1]$ , where  $r_2 = \{r^1 \lhd \cdots \lhd r^{k_1}\}$  and  $s_2 = \{s^1 \lhd \cdots \lhd s^{k_2}\}$ . Due to the macro tree order of the tree and the fact that X represents a deep set, no node in X[i] can be to the right of any node in  $X[r_1]$  for  $i < r_1$ . To prove condition 3 it is thus enough to prove it for  $i = r_1$ . We proceed by induction on the number k of iterations of the outer loop (line 3–34). We consider the time right after the kth iteration of the loop, i.e., right before the (k+1)th iteration. The base case (k=0) is trivially satisfied.

For the induction step let  $i^*$  and  $j^*$  be the values of i and j at line 14 in iteration k. Let  $r'_i$  and  $s'_i$  for i = 1, 2 be the values of  $r_i$  and  $s_i$ , respectively, after the (k-1)th round. There are 3 cases:

- 1.  $r'_2 = r_2$  and  $s'_2 = s_2$ : the claim follows directly from the induction hypothesis.
- 2.  $r_2' \neq r_2$  and  $s_2' = s_2$ : condition 2 from the lemma follows directly from the induction hypothesis. Since  $s_2$  and thus also  $s_1$  were not changed,  $r_2$  was set in case 1 of the procedure and  $j^* = s_1$ . Therefore,  $i^* \triangleleft j^*$ ,  $r_1 = i^*$ , and  $|r_2| = 1$ . Let  $r_2 = \{r^1\}$  and  $s_2 = \{s^1\}$ . We have  $r_1 = i^* \triangleleft j^* = s_1$  and thus  $r^1 \triangleleft s^1$

satisfying condition 1 from the lemma. To prove condition 3 is satisfied we only have to consider the case  $i = r_1$ . Since  $r_2$  was set in case 1 of the procedure,  $r^1$  is the rightmost node in  $X[r_1]$  and it follows immediately that there exists no node  $x \in X[r_1]$  such that  $r^1 \triangleleft x \triangleleft s^1$ .

3.  $r_2' \neq r_2$  and  $s_2' \neq s_2$ : We first prove condition 1 and 3. If the potential pair was set in case 1 (line 15–19) of the procedure then  $r_1 = i^* \lhd j^* = s_1$  and  $|r_2| = 1$  implying  $r^1 \lhd s^1$  (condition 1). The node  $r^1$  is the rightmost node in  $X[r_1]$  (line 19) and it follows that there exists no node  $x \in X[r_1]$  such that  $r^1 \lhd x \lhd s^1$  proving condition 3. If the potential pair was set in case 2 then both condition 1 and 3 follows from the correctness of the implementation of mop and the computation  $(r,s) = \text{mop}_{C(i,j,v)}(X[i^*],Y[j^*]) = \text{mop}_{C(i,j,v)}(X[i^*],Y[j^*])$  in line 21.

Let  $y \in Y[j]$ , for  $j \leq s_1$ , be a node such that  $y \notin s_2$ . To prove condition 2 is satisfied we will show that  $r^l$  is not to the left of y. There are two cases

- $j = s_1$ . Since  $s'_2 \neq s_2$  there are two cases depending on which case of the procedure the potential pair was set in. If the potential pair was set in case 2 of the procedure the claim follows from the correctness of the implementation of mop and the computation  $(r, s) = \text{mop}_{C(i,j,v)}(X[i^*], Y[j^*]) = \text{mop}_{C(i,j,v)}(X[i^*], Y[s_1])$  in line 21.
  - If the potential pair was set in case 1, then  $r_1 = i^* \triangleleft j^* = s_1$ . Since  $s_2 \neq s_2'$ ,  $s_2$  was changed in the kth iteration and is therefore the leftmost node in  $Y[s_1]$  (line 17). The claim follows.
- $j < s_1$ . We will use that we just proved the claim for  $j = s_1$ . Assume for the sake of contradiction that there exists a  $y \in Y[j]$  such that  $r^l \triangleleft y$ . Since Y is representing a deep set and due to the macro tree order of Y this implies  $r^l \triangleleft y \triangleleft y'$  for all  $y' \in Y[s_1]$  contradicting that the claim is true for  $j = s_1$ .

**Lemma 25** We have the following invariant at the beginning of each iteration of the main loop (line 3) of MOPSIM:

$$\nexists x \in X[i]$$
, such that  $x \leq x'$ , for any  $x' \in r_2$ .

Proof. By induction on the number of iterations of the outer loop. In the base case  $r_2 = \emptyset$  and the condition is trivially satisfied. Note that X is representing a deep set and thus either  $x \triangleleft x'$  or  $x' \triangleleft x$  for all  $x \in X[i]$ . For the induction step let i', j', and  $r'_2$  be the values of i, j, and  $r_2$  respectively in the iteration before this. By the induction hypothesis  $x' \triangleleft x$  for all  $x \in X[i']$  and  $x' \in r'_2$ . Due to the macro tree order of X and the fact that X represents a deep set, all nodes in X[i'] are to the left of all nodes in X[i]. Thus, if  $r_2 = r'_2$  it follows from the induction hypotheses that  $x' \triangleleft x$  for all  $x \in X[i]$  and  $x' \in r'_2 = r_2$ . For  $r'_2 \neq r_2$  there are two cases: If  $i' \triangleleft j'$  then  $r_2 = \text{RIGHT}_{C(i')}(1, X[i'])$  and i > i' and thus the condition is satisfied. Otherwise  $r_2$  was set in case 2 of the procedure. Since  $r_2 \neq r'_2$  we have  $r_2 = r \subseteq X[i']$  and  $r \neq \emptyset$ . There are two subcases: If i = j or i = l(v, w) and j = s(v, w) (Case 2(a) of the procedure) then X[i] either contains a single node, which is the rightmost of the nodes in X[i'] that are to the right of all nodes in  $r_2$  or if there are no such nodes  $X[i] = \emptyset$ . In both cases the condition is satisfied. If i = s(v, w) and j = r(v, w) then i > i' and the condition is satisfied.

#### **Lemma 26** Procedure MOPSIM is correctly implemented.

*Proof.* Let  $\mathcal{X}$  and  $\mathcal{Y}$  be the sets represented by X and Y, respectively. Let  $R = \text{MopSim}(X,Y)|_1$  and  $S = \text{MopSim}(X,Y)|_2$ . For simplicity we will slightly abuse the notation and write  $(x,y) \in \text{MopSim}(X,Y)$  iff there exists an i such that  $x \in R[i]$  and  $y \in S[i]$ . We want to show that

$$(x,y) \in \text{MopSim}(X,Y) \Leftrightarrow (x,y) \in \text{mop}(\mathcal{X},\mathcal{Y}).$$

Assume  $(x,y) \in \text{MOPSIM}(X,Y)$ . Consider the round where x and y were added to R and S, respectively. We have  $x = r^l \in r_2$  and  $y = s^l \in s_2$ . We want to show that there is no node  $x' \in X[i]$  for any i such that  $x \triangleleft x' \triangleleft y$  and no node  $y' \in Y[j]$  for any j such that  $x \triangleleft y' \triangleleft y$ . By Lemma 24 this is true for  $i \leq r_1$  and  $j \leq s_1$ . By the macro tree order of Y we have that  $y \triangleleft y'$  for any  $y' \in Y[j]$  when  $j > s_1$ . Let i' be the value of i in the round where x and y is added to the output. We will show that no node in X[i'] is to the left of any node in  $s_2$ . Due to the macro tree order of X this implies that no node in X[i] is to the left of any node in  $s_2$  for any  $i \geq i'$ . If  $i' = r_1$  then it follows directly from Lemma 24. If  $i' > r_1$  it follows from the implementation of the procedure that i' is the first non-empty entry in X greater than  $r_1$ . Thus the claim follows for any j. We now return to show that no node in X[i'] is to the left of any node in  $s_2$ . There are two cases depending on whether  $j = s_1$  or  $j > s_1$ . If  $j > s_1$  then j was changed either in one of the four cases I–IV, or in the previous iteration in case 2. If j was equal to  $s_1$  at the beginning of this iteration then j was incremented in one of the four cases I–IV. Thus none of the cases applied to  $s_1$ . By Proposition 6 no node in X[i'] can be to the left of a node in  $X[s_1]$ . Since  $s_2 \subseteq X[s_1]$  the claim follows. If  $j = s_1$  it follows from case 2 that LEFT $(X[i'], s_2) = \emptyset$  (otherwise the potential pairs would not have been added to the output in this iteration) and the claim follows immediately.

Now assume  $(x,y) \in \text{mop}(\mathcal{X},\mathcal{Y})$ . We will deal with each of the cases from Proposition 6 separately.

```
1. Case (i): c(x) = c(y) = r(v, w).
```

- 2. Case (i): c(x) = c(y) = l(v, w).
- 3. Case (ii): c(x) = c(y) = l(v).
- 4. Case (iii): c(x) = l(v, w) and c(y) = s(v, w).
- 5. Case (iv): c(x) = s(v, w) and c(y) = r(v, w).
- 6. Case (v): c(x) = l(v, w) and c(y) = r(v, w).
- 7. Case (v):  $c(x) \triangleleft c(y)$  and c(x) and c(y) belong to different clusters.

Note that if  $c(x) \triangleleft c(y)$  then x is the rightmost node in X[c(x)] and y is the leftmost node in Y[c(y)]. We first show that in all cases we will have  $x = r^l \in r_2$  and  $y = s^l \in s_2$  for some l at some iteration. Consider the first iteration where either  $x \in X[i]$  or  $y \in Y[j]$ . Let i' and j' be the values of i and j, respectively, in this iteration. There are three cases:

- (a)  $x \in X[i']$  and  $y \in Y[j']$ . For case 1–5 the procedure goes into case 2. From the correctness of MOP we get  $x \in r$  and  $y \in s$ . Thus  $r \neq \emptyset$  and we set  $(r_1, r_2) = (i', r)$  and  $(s_1, s_2) = (j', s)$  and the claim follows. For case 6–7 the procedure goes into case 1. Since this iteration is the first where  $y \in Y[j]$  we have  $j' > s_1$  and we set  $(r_1, r_2) = (i', \text{RIGHT}_{C(i')}(1, X[i']))$  and  $(s_1, s_2) = (j', \text{LEFT}_{C(j')}(1, Y[j']))$ . Since x is the rightmost node in X[i'] and y is the leftmost node in Y[c(j')] the claim follows.
- (b)  $x \in X[i']$  and  $y \notin Y[j']$ . Since  $(x,y) \in \text{mop}(\mathcal{X},\mathcal{Y})$  this implies j' < c(y) and there exists no node  $y' \in Y[j']$  such that  $x \triangleleft y'$ . Assume that there existed such a y'. Then  $x \triangleleft y' \triangleleft y$  due to the macro tree order of Y contradicting  $(x,y) \in \text{mop}(\mathcal{X},\mathcal{Y})$ . Thus  $i' \not \vartriangleleft j'$ . From case I–IV of the procedure it follows that either i' = j', i' = l(v, w) and j' = s(v, w), or i' = s(v, w) and j' = r(v, w). From this and j' < c(y) it follows that we are in case 4 or 7 from above.

The procedure enters case 2 in this iteration. If we are in case 4 then i' = l(v, w) = j' and c(y) = s(v, w). If  $r = \emptyset$  then i = i', X[i'] is unchanged, and j = j' + 1 = s(v, w) = c(y) at the end of this iteration. If  $r \neq \emptyset$  then x must be to the right of all nodes in  $x' \in r$ . Assume that there is a  $x' \in r$  such that  $x \triangleleft x'$ . Since  $x' \in r$  there exists a node  $y' \in s$  such that  $x' \triangleleft y' \triangleleft y$ . That  $y' \triangleleft y$  follows from  $y' \in l(v, w)$  and  $y \in s(v, w)$  and the assumption that  $\mathcal{Y}$  is deep. Thus  $x \triangleleft x' \triangleleft y' \triangleleft y$  contradicting that  $(x, y) \in \text{mop}(\mathcal{X}, \mathcal{Y})$ . Therefore, i = i',  $x \in X[i']$  and j = j' + 1 = s(v, w) = c(y) at the end of this iteration. From case I of the procedure and the analysis of case (a) it follows that  $x = r^l \in r_2$  and  $y = s^l \in s_2$  for some l.

Now assume we are in case 7. By the same argument as before  $i=i', x \in X[i]$ , and j>j' at the end of this iteration. Unless i'=l(v,w)=j' this implies that  $i \triangleleft j$  at line 14 ("Compare i and j") in the next iteration. If i'=l(v,w)=j' then either  $i \triangleleft j$  after the first loop in the next iteration (line 7), and the claim follows as before, or i=l(v,w) and j=s(v,w). In the last case we get into case (b) again, but it follows from the analysis that in the iteration after the next we will have  $i \triangleleft j=c(y)$ . The claim follows from the analysis of case (a).

(c)  $x \notin X[i']$  and  $y \in Y[j']$ . It follows by inspection of the cases that unless we are in case 1 we have  $i' \lhd j'$ . If we are in case 1 (j' = c(x) = r(v, w) = c(y)) we have either  $i' \lhd j'$  or i' = s(v, w). First we consider the cases 2–7. Since  $i' \lhd j'$  the procedure enters case 1 in this iteration. Thus i is incremented and j stays the same. This happens until i = c(x). Now consider case 1. If  $i' \lhd j'$  the procedure enters case 1 in this iteration. Thus i is incremented and j stays the same. In the next iteration either the same happens or i' = s(v, w). If i' = s(v, w) the procedure enters case 2. Since i' is a spine node and  $\mathcal{X}$  is deep, X[i] contains only one node x'. By the structure of the macro tree and the assumption that  $\mathcal{X}$  is deep  $x' \lhd x$ . Since  $x \lhd y \in Y[j']$  this implies  $r \ne \emptyset$ . It follows from case 2(b) of the procedure that i is incremented while j stays the same. At line 14 ("Compare i and j") in the next iteration we will have j = i = r(v, w) since all entries in X between j' and r(v, w) are empty due to the assumption that  $\mathcal{X}$  is deep. The claim follows from the analysis in case (a).

It remains to show that once  $x = r^l \in r_2$  and  $y = s^l \in s_2$  they will stay this way until added to the output. Consider the iteration where x and y are assigned to  $r_2$  and  $s_2$ . At the end of this iteration either i or j or both are incremented. Assume j is incremented while the potential pairs are still unchanged. Since j is incremented we have  $s_1 < j$  until  $s_1$  is changed. It follows from case 1 and 2 of the procedure that in this case  $(r_1, r_2)$  is only changed if at the same time  $(s_1, s_2)$  are changed and right before that  $(r_1, r_2)$  and  $(s_1, s_2)$  are added to the output.

Consider first case 1–3. If i is incremented then j is incremented in one the cases I–V in the next iteration since i'=j'. By the above argument x and y are added to the output. For case 4 j is incremented (case 2(a) of the procedure) and the claim follows as before. For case 5–7 first note that  $r_2$  and  $s_2$  contain only one node each, i.e.,  $x=r_2$  and  $y=s_2$ . For case 5 i is incremented (case 2(b) of the procedure). Since  $\mathcal{X}$  is deep we have  $i \geq r(v,w) = j'$  at line 14 ("Compare i and j") in the next iteration. If i > r(v,w) then j > j' and the claim follows. If i = r(v,w) the procedure enters case 2. If  $r = \emptyset$  then j is incremented and the claim follows. If  $r \neq \emptyset$  then  $s_1 = j$  and  $(x,y) \in \text{mop}(\mathcal{X},\mathcal{Y})$  implies  $\text{LEFT}_{C(i,j)}(X[i],s_2) = \text{LEFTOF}_{C(i,j)}(X[i],y) = \emptyset$ . Thus  $(r_1,r_2)$  and  $(s_1,s_2)$  are added to the output. If we are in case 6 and 7, i is incremented. Consider case 6. Since  $(x,y) \in \text{mop}(\mathcal{X},\mathcal{Y})$  all entries in X between l(v,w) and r(v,w) are empty. Thus at line 14 ("Compare i and j") in the next iteration  $i \geq r(v,w)$ . The proof is equivalent to the one for case 5. Consider case 7. If j is a boundary node then all entries in X between c(x) and j are empty. Thus j is incremented in the second loop of the next iteration. For all other cases for j the proof is similar to the proof of case 5.

**Lemma 27** In procedure MATCH we have the following invariant of X[i] and Y[j] in line 6:

LEFT
$$(1, X[i]) = \mathcal{X}_l$$
 and LEFT $(1, Y[j]) = \mathcal{Y}_l$  for some  $l$ .

Proof. Induction on the number of iterations of the outer loop. Base case: In the first iteration X[i] and Y[j] are the first nonempty entries in X and Y and thus LEFT $(1, X[i]) = \mathcal{X}_1$  and LEFT $(1, Y[j]) = \mathcal{Y}_1$ . For the induction step let i' and j' be the values of i and j in the previous iteration. By the induction hypothesis LEFT $(1, X[i']) = \mathcal{X}_{l'}$  and LEFT $(1, Y[j']) = \mathcal{Y}_{l'}$ . If x = |X[i']| = |Y[j']| both i and j were incremented and LEFT $(1, X[i]) = \mathcal{X}_{l'+x}$  and LEFT $(1, Y[j]) = \mathcal{Y}_{l'+x}$ . If x = |X[i']| < |Y[j']| then i was incremented implying LEFT $(1, X[i]) = \mathcal{X}_{l'+x}$ . In that case j = j' and Y[j] = LEFT(x, Y[j']) implying LEFT $(1, Y[j]) = \mathcal{Y}_{l'+x}$ . Similarly, if |X[i']| > |Y[j']| = y we have LEFT $(1, X[i]) = \mathcal{X}_{l'+y}$ . In that case j = j' and LEFT $(1, Y[j]) = \mathcal{Y}_{l'+y}$ .  $\square$ 

**Lemma 28** Procedure MATCH is correctly implemented.

Proof. We need to show that for all  $1 \leq k \leq |\mathcal{X}|$ :  $\mathcal{X}_k \in \text{MATCH}(X,Y,Y') \Leftrightarrow \mathcal{X}_k \in \{\mathcal{X}_j|\mathcal{Y}_j \in \mathcal{Y}\}$ . Consider the iteration where  $\mathcal{X}_k \in X[i]$  and  $\mathcal{Y}_k \in Y[j]$ . By Lemma 27 such an iteration exists. If Y[j] = Y[j'] then  $\mathcal{Y}_k \in \mathcal{Y}'$  implying  $\mathcal{X}_k \in \{\mathcal{X}_j|\mathcal{Y}_j \in \mathcal{Y}\}$ . It follows from the implementation of case 1(a) and 1(b) that if  $x \leq y$  all nodes in X[i] are added to the output and thus  $\mathcal{X}_k \in \text{MATCH}(X,Y,Y')$ . If x > y then  $\mathcal{X}_k \in \text{LEFT}(y,X[i])$  since  $\mathcal{Y}_k \in Y[j]$  and thus  $\mathcal{X}_k \in \text{MATCH}(X,Y,Y')$ .

If  $Y[j] \neq Y'[j]$  the procedure calls MATCH with some subset of X[i], Y[j], and Y'[j] depending on the size of x and y. By Lemma 27 and the correctness of MATCH it follows that  $\mathcal{X}_k \in \text{MATCH}(X, Y, Y') \Leftrightarrow \mathcal{X}_k \in \{\mathcal{X}_j | \mathcal{Y}_j \in \mathcal{Y}\}.$ 

## Lemma 29 Procedure MopRight is correctly implemented.

*Proof.* Follows from the correctness of MopSim (Lemma 26) and Match (Lemma 28).  $\hfill\Box$ 

Finally, we consider correctness of the FL procedure.

### Lemma 30 Procedure FL is correctly implemented.

Proof. Let  $\mathcal{X}$  denote the set represented by X and let  $F = \{ \mathrm{fl}(x, \alpha) | x \in \mathcal{X} \}$ . To show  $\mathrm{FL}(X, \alpha) \subseteq F$  we will first show that for any node x added to R during the computation  $x \in F$ . Consider a node  $x \in R[i]$  for some i. Either x was added directly to R after a computation of N in one of the three cases of the procedure or it was added after the computation of S. In the first case  $x \in F$  follows from the correctness of  $\mathrm{FL}_C$ . If x was added after the computation of S it follows from the correctness of  $\mathrm{FL}_M$  that  $x \in C(i)$  for some  $i \in S$ . Due to the correctness of  $\mathrm{FL}_C$  we have  $x \in F$ .

To show  $\mathrm{DEEP}(F) \subseteq \mathrm{FL}(X,\alpha)$  we use Proposition 5. Let x be a node in  $\mathrm{DEEP}(F)$  and let x' be a node in X such that  $\mathrm{fl}(x',\alpha)=x$ . We have  $x'\in X[i]$  for some i. If i is a left or right node then according to Proposition 5 x can be in i (case (i)), on the spine (case (ii)), in the top boundary node (case (ii)), or in an ancestor of i in the macro tree (case (iii)). If x is in the same cluster as x' then it follows from the correctness of  $\mathrm{FL}_C$  that  $x\in N$ . Thus x is added to R and due to the correctness of  $\mathrm{DEEP}$  we have  $x\in \mathrm{FL}(X,\alpha)$ . If c(x) is in a different cluster than c(x') then c(x) is an ancestor of c(x') in the macro tree due to Proposition 5. Since  $x\in \mathrm{DEEP}(F)$  we have  $N=\emptyset$  and thus parent $(v)\prec_M c(x')$  is added to L. It follows from the correctness of  $\mathrm{FL}_M$  that  $c(x)\in S$ . Due to the structure of the macro tree c(x) is either a boundary node or a spine node and thus  $x=\mathrm{fl}_{C(c(x))}(\mathrm{first}(c(x)),\alpha)=\mathrm{FL}_{C(c(x))}(\mathrm{first}(c(x)),\alpha)$ . The last equality follows from the correctness of  $\mathrm{FL}_C$ . That  $x\in \mathrm{FL}(X,\alpha)$  now follows from the above analysis showing that only nodes from F are added to R and the correctness of  $\mathrm{DEEP}$ .

If i is a leaf node then x can be in i (case (i)), in the top boundary node (case (iii)), or in an ancestor of i in the macro tree (case (iii)). The correctness follows by an analysis similar to the one for the previous case. If i is a spine node or a boundary node, then x is either in i (case (i)) or in an ancestor of i in the macro tree (case (iii)). The correctness follows by an analysis similar to the one for the first case.

## 5.5 Complexity of the Tree Inclusion Algorithm

To analyze the complexity of the node array implementation we first bound the running time of the above implementation of the set procedures. All procedures scan the input from left-to-right while gradually producing the output. In addition to this procedure FL needs a call to a node list implementation of FL on the macro tree. Given the data structure described in Section 5.2 it is easy to check that each step in the scan can be performed in O(1) time giving a total of  $O(n_T/\log n_T)$  time. Since the number of nodes in the macro tree is  $O(n_T/\log n_T)$ , the call to the node list implementation of FL is easily done within the same time. Hence, we have the following lemma.

**Lemma 31** For any tree T there is a data structure using  $O(n_T)$  space and  $O(n_T \log n_T)$  preprocessing time which supports all of the set procedures in  $O(n_T/\log n_T)$  time.

Next consider computing the deep occurrences of P in T using the procedure Emb of Section 3 and Lemma 31. The following lemma bounds the space usage.

**Lemma 32** The total size of the saved embeddings at any time during the computation of Emb(root(P)) is  $O(n_T)$ .

Proof. Let v be the node for which we are currently computing EMB. Let p be the path from the root to v and let  $w_0, \ldots, w_l$  be the light nodes on this path. We have l = ldepth(v). As in the proof of Lemma 15 it suffices to bound  $|\text{EMB}(\text{heavy}(\text{parent}(w_i)))|$  for all i. Assume that  $l_P \leq l_T$  (otherwise we can check this in linear time and conclude that P cannot be included in T). Each of the node arrays use  $O(n_T/\log n_T)$  space and therefore by Corollary 1 we have that  $\sum_{i=1}^{l} |\text{EMB}(\text{heavy}(\text{parent}(w_i)))| = O(n/\log n_T \cdot \log l_P) = O(n_T)$ .

For the time complexity note that during the computation of EMB(root(P)) each node  $v \in V(P)$  contributes a constant number of calls to the set procedures. Hence, the total time used by the algorithm is  $O(n_P n_T / \log n_T + n_T \log n_T)$ . Thus we have shown the following.

**Theorem 4** For trees P and T the tree inclusion problem can be solved in  $O(n_P n_T / \log n_T + n_T \log n_T)$  time and  $O(n_T)$  space.

Combining the results in Theorems 2, 4 and Corollary 2 we have the main result of Theorem 1.

## 6 Conclusion

We have presented three algorithms for the tree inclusion problem, which match or improve the best known time complexities while using only linear space. We believe that some of the new ideas are likely to be of both practical and theoretical value in future work. From a practical perspective, space is a common bottleneck for processing large data sets and hence reducing the space can significantly improve performance in practice. From a theoretical perspective, we have introduced several non-trivial algorithms to manipulate sets of nodes in trees that may have applications to other problems. For instance, the NCA procedure from Section 5 computes multiple nearest common ancestor queries in time *sublinear* in the size of input sets.

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