Fully Dynamic Maintenance of *k*-Connectivity in Parallel

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Abstract—Given a graph G = (V, E) with n vertices and m edges, the k-connectivity of G denotes either the k-edge connectivity or the k-vertex connectivity of G. In this paper, we deal with the fully dynamic maintenance of k-connectivity of G in the parallel setting for k = 2, 3. We study the problem of maintaining k-edge/vertex connected components of a graph undergoing repeatedly dynamic updates, such as edge insertions and deletions, and answering the query of whether two vertices are included in the same k-edge/vertex connected component. Our major results are the following: 1) An NC algorithm for the 2-edge connectivity problem is proposed, which runs in $O(\log n \log(m/n))$ time using $O(n^{3/4})$ processors per update and query. 2) It is shown that the biconnectivity problem can be solved in $O(\log^2 n) \tan O(n\alpha(2n, n)/\log n)$ processors per update and O(1) time with a single processor per query or in $O(\log n \log \frac{m}{n})$ time using $O(n\alpha(2n, n)/\log n)$ processors per update and $O(\log n)$ time using $O(n\alpha(2n, n)/\log n)$ processors per query. 4) An NC algorithm for the 3-edge connectivity problem. To the best of our knowledge, the proposed algorithms are the first NC algorithms for the proposed NC algorithm for the 2-edge connectivity problem. To the best of our knowledge, the proposed algorithms run on a CRCW PRAM.

Index Terms—NC algorithms, 2-edge/vertex connectivity, 3-edge/vertex connectivity, dynamic data structures, parallel algorithm design and analysis, graph problems.

1 INTRODUCTION

A fully dynamic graph algorithm is one that allows edge insertions and deletions and recomputes a desired graph property quickly after each such update. A partially dynamic graph algorithm is one that allows edge insertion updates only. In this paper, we focus on the design of fully dynamic NC algorithms for maintaining the k-connectivity of graphs. Given a graph G = (V, E)with n vertices and m edges, the k-connectivity of G refers to either the k-edge connectivity or k-vertex connectivity [18], which will be formally defined in Section 2.1.

1.1 Previous Related Results

k-connectivity of *G* is a basic property in graph theory with wide applications in robust routing, distributed computing, reliable communications networks design, etc. There have been extensive studies of algorithms for dynamically maintenance of the *k*-connectivity of graphs with $k \leq 4$. For example, Even and Shiloach [12] studied the connected component problem (k = 1) in the early 1980s. Frederickson [15] studied the minimum spanning

• H. Shen is with the School of Computing and Information Technology, Griffith University, Nathan, QLD 4111, Australia. E-mail: hong@cit.gu.edu.au. tree problem by giving a fully dynamic algorithm with $O(\sqrt{m})$ time per update and O(1) time per query. Eppstein et al. [10] improved Frederickson's algorithm by presenting an $O(\sqrt{n}\log(m/n))$ algorithm, using a sparsification technique. Later, Eppstein et al. [11] improved their own algorithm further by an $O(\log(m/n))$ factor. Henzinger and King [25] gave an algorithm for fully dynamic maintenance of minimum spanning trees with $O(\sqrt[3]{n}\log n)$ time per update and constant time per query. Randomization has also been used to improve the time complexity of fully dynamic maintenance of connected components. Henzinger and King [24], [23] showed that a spanning forest can be maintained in $O(\log^3 n)$ expected amortized time per update, and the query can be answered in $O(\log n / \log \log n)$ time. The update time was further improved to $O(\log^2 n)$ by Henzinger and Thorup [27]. Since the initial submission of this paper in 1995, there has been further development in this topic. In particular, Holm et al. [28] recently gave an improved algorithm for the fully dynamic maintenance of connected components with $O(\log^2 n)$ amortized time per update and $O(\log n / \log \log n)$ time per query and an improved algorithm for the fully dynamic maintenance of minimum spanning forests with $O(\log^4 n)$ amortized time per update and $O(\log n / \log \log n)$ time per query, thus matching the previously best randomized bound and improving substantially the previously best deterministic bounds of $O(\sqrt[3]{n}\log n)$. Note that Holm et al.'s algorithm assumes that the initial graph is empty and the update time is the amortized time. Thus, a single insertion or deletion operation in their algorithm takes $\Omega(n)$ time in the worst case. In the parallel setting, Ferragina [13] studied fully

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dynamic maintenance of connected components on an EREW PRAM provided only a small number of processors are used, and the proposed parallel algorithm requires $O((\frac{n \log n}{p} + p^{1+\epsilon})(\frac{\log f}{\log \log n}))$ time using O(p) processors per update, where $p \leq (n \log n)^{1/(2+\epsilon)}$, $f = \frac{m}{n \log n}$, and $0 \leq \epsilon < 1$. Das and Ferragina [8] and Ferragina [14] developed parallel algorithms for the fully dynamic maintenance of minimum spanning trees and connected components which requires $O(\log n)$ time and the total of $O(n^{2/3} \log \frac{m}{n})$ work. Independently, Liang and McKay [35] also proposed parallel algorithms for the same problem which requires $O(\log n \log \frac{m}{n})$ time and $O(n^{2/3})$ processors.

As Galil and Italiano [19] pointed out, the fully dynamic maintenance of 2-connectivity, however, is much harder than the fully dynamic maintenance of connected components because the maintenance of connected components only involves merging two connected components into one connected component or splitting a connected component into two connected components per update, whereas the maintenance of 2-edge/vertex connected components may need merging O(n) 2-edge/vertex connected components into one or splitting a 2-edge/vertex connected component into O(n) 2-edge/vertex connected components per update. Westbrook and Tarjan [45] presented the first partially dynamic algorithms for the maintenance of both 2-edge connected components and biconnected (2-vertex connected) components, which require $O(m\alpha(m, n))$ time for the maintenance of 2-connectivity of graphs when there are m queries and edge insertions. In other words, the algorithm takes $O(\alpha(m, n))$ amortized time per update or per query, where $\alpha(m,n)$ is the inverse of Ackermann's function. Westbrook and Tarjan left an open problem as to whether there exists a sublinear time, fully dynamic algorithm for the 2-connectivity problem. Since then, finding a sublinear time algorithm for the problem has been a challenging task. Galil and Italiano [19] presented the first sublinear time algorithm for 2-edge connectivity, by using Frederickson's clustering technique and the other techniques. Their algorithm requires $O(m^{2/3})$ time. Later, Frederickson [16] improved this algorithm to $O(\sqrt{m})$ using an ambivalent data structure. Eppstein et al. [10], [11] further improved Frederickson's algorithm by developing $O(\sqrt{n}\log(m/n))$ and $O(\sqrt{n})$ time algorithms, based on their sparsification technique. They also derived algorithms for the fully dynamic maintenance of biconnectivity and triconnectivity, which take $O(\alpha(q, n))$ amortized time per insertion or per query, and O(n) time per deletion, where q is the total number of queries made. Furthermore, Henzinger and King [24] generalized their randomization technique for connected components by giving an $O(\log^5 n)$ expected amortized time bound per update. Recently, Holm et al. [28] extended their results for connected components by presenting an algorithm for the fully dynamic maintenance of 2-edge connected components. Their algorithm requires $O(\log^4 n)$ amortized time per update.

The fully dynamic maintenance of biconnected components seems much harder than that of 2-edge connected components. There is a reduction technique [18] which can reduce a k-edge connectivity problem to a k-vertex connectivity problem, but the reverse reduction has not yet been found. In 1992, Rauch [40] presented the first fully dynamic algorithm for the maintenance of biconnected components. Her algorithm takes $O(m^{2/3})$ amortized time per update and O(1) time per query. She later improved her algorithm to $O(\min\{\sqrt{m}\log n, n\})$ using new data structures and the sparse certificate technique [41]. In 1995, Henzinger and Poutré [26] further improved her result to $O(\sqrt{n \log n} \log \frac{m}{n})$. Later, Henzinger and King [23], [24] presented polylogarithmic time randomized algorithms for the fully dynamic maintenance of biconnectivity [23] using a novel decomposition of graphs and the randomization technique. Their algorithm requires $O(\Delta \log^4 n)$ expected amortized time per update, where Δ is the maximum degree of the graph. Holm et al. [28] further improved the above results by giving a deterministic algorithm, which takes $O(\log^4 n)$ amortized time per update.

When k = 3, Galil and Italiano [20] presented a partially dynamic algorithm for the maintenance of 3-edge connected components. Their algorithm takes $O((n+q)\alpha(q,n))$ time for a sequence of q queries and updates on an n-vertex graph. Independently, La Poutré et al. [34] presented a partially dynamic algorithm which has the same time bound as Galil Italiano [20]. Later, Galil and Italiano [21] claimed that they developed a fully dynamic algorithm for the problem which requires $O(m^{2/3})$ time per update and per query by minor modifications to their 2-edge connectivity algorithm [19]. Moreover, La Poutré [33] also proposed a partially dynamic algorithm for the maintenance of triconnected components, which needs $O((n+q)\alpha(q,n))$ time for a sequence of q queries and updates. As for k > 3, there are only a few results available. Kanevsky et al. [30] presented a partially dynamic algorithm for the maintenance of 4-vertex connected components which requires $O(\alpha(q, n))$ amortized time per update and per query provided that the graph is triconnected, where q is the number of operations performed. Dinitz [9] presented a partially dynamic algorithm for the maintenance of 4-edge connected components. Note that using Eppstein et al.'s [10], [11] sparsification technique and the above algorithms, the following results then follow easily: 1) fully dynamic maintenance of 3-edge connected components can be done in $O(n^{2/3})$ time per update and per query, 2) fully dynamic maintenance of triconnected components can be done in $O(\alpha(q, n))$ time per insertion or per query, and in O(n) time per deletion, and 3) fully dynamic maintenance of 4-vertex connected components can be done in $O(\log n)$ time per insertion, in $O(n \log n)$ time per deletion, and in O(1) time per query.

Surprisingly, we have not seen any NC algorithms for the fully dynamic maintenance of *k*-connectivity with $k \ge 2$ in the parallel environment. In particular, we have not seen any NC algorithms for the addressed problems using O(n) processors only. In this paper, we will focus on developing such NC algorithms for the problems for the first time. It must be mentioned that the previously known sequential algorithms except the algorithms by Henzinger and King [23], [24] and Holm et al. [28] are hardly parallelizable since they use the restricted clustering decomposition and depth-first search technique which are inherently sequential. The algorithms by Henzinger and King [23], [24] are Las Vegas type of randomized algorithms, which are incomparable to the deterministic algorithms, because the deterministic simulation of the randomized algorithms may require $\Omega(m)$ amount of work. The algorithm due to Holm et al. [28] does guarantee that each update can be done in $O(\log^4 n)$ amortized time, but the time complexity is amortized, which means that all the data structures must be rebuilt after undergoing $\Omega(m)$ updates. The rebuild takes $\Omega(n)$ time. In other words, it takes $\Omega(n)$ time per update in the worst case. Therefore, even if their algorithm can be parallelized, $\Omega(n)$ processors must be required in order to achieve a polylogarithmic time complexity per update in the worst case.

1.2 Our Contributions

In this paper, we deal with the fully dynamic maintenance of *k*-edge/vertex connectivity property of a graph in the parallel environment with k = 2,3 after undergoing an intermixed sequence of insertion, deletion, and query operations, which are described as follows:

- SameConnComp(x,y): return true if vertices x and y are in the same k-edge/vertex connected component; otherwise, return false.
- *InsertEdge(x,y):* insert a new edge between vertex *x* and vertex *y.*
- *DeleteEdge(x,y):* delete an existent edge between vertex *x* and vertex *y*.

We present NC algorithms for the fully dynamic maintenance of *k*-edge/vertex connected components in *G* for k = 2, 3 by using O(n) processors in contrast to the usual $\Omega(m)$ processors needed for computing these properties from scratch. Our major results are as follows:

1. An NC algorithm for the 2-edge connectivity problem is presented, which requires

$$O(\log n \log(m/n))$$

time using $O(n^{3/4})$ processors per update and per query;

- 2. An NC algorithm for the biconnectivity problem is also given, which requires either $O(\log^2 n)$ time using $O(n\alpha(2n, n)/\log n)$ processors per update and O(1) time with a single processor per query or $O(\log n \log \frac{m}{n})$ time using $O(n\alpha(2n, n)/\log n)$ processors per update and $O(\log n)$ time with $O(n\alpha(2n, n)/\log n)$ processors per query;
- 3. An NC algorithm for the triconnectivity problem is derived, which requires

$$O(\log n \log \frac{m}{n} + \log n \log \log n / \alpha(3n, n))$$

time using $O(n\alpha(3n, n)/\log n)$ processors per update and O(1) time with a single processor per query;

4. An NC algorithm for the 3-edge connectivity problem is obtained, which has the same time and

processor complexities as the algorithm for the fully dynamic maintenance of triconnectivity.

To the best of our knowledge, the proposed algorithms for the above problems are the first NC algorithms that use only O(n) processors in contrast to $\Omega(m + n)$ processors for solving them from scratch. At the same time, it should be mentioned that the basic idea of the proposed algorithm for the fully maintenance of 2-edge connectivity comes from Galil et al.'s sequential algorithm, although, their algorithm by itself is not easily parallelizable. Our NC algorithm for this problem is achieved by replacing the sequential part of their algorithm with a highly parallelizable approach.

In this paper, the following three parallel computational models will be used:

- 1. An Exclusive Read and Exclusive Write (EREW for short) PRAM in which concurrent read and concurrent write are forbidden.
- 2. A CREW PRAM in which Concurrent Read is allowed but only Exclusive Write is permitted.
- 3. A CRCW PRAM in which both Concurrent Read and Concurrent Write are allowed, but the write conflicts are resolved arbitrarily.

The remainder of this paper is organized as follows: In Section 2, we introduce some basic concepts with respect to the *k*-edge/vertex connectivity, such as the sparse *k*-edge/ vertex certificates, Galil et al.'s reduction technique [18], and Eppstein et al.'s sparsification technique [10]. In Section 3, for the sake of simplicity, we consider the fully dynamic maintenance of 2-edge connected components in a connected graph. The data structures used for the maintenance are described and the algorithms for connectivity queries, edge insertions, and edge deletions are presented. In Section 4, we first extend the results in connected graphs to disconnected graphs and then present an improved NC algorithm for fully dynamic maintenance of 2-edge connectivity in general graphs. We finally discuss the fully dynamic maintenance of 3-edge connected components. In Section 5, we deal with the fully dynamic maintenance of biconnected components and triconnected components and we conclude our discussions in Section 6.

2 Preliminaries

2.1 Definitions

G(V, E) is k-edge connected if it is still connected after deleting any k - 1 edges from it. Assuming that G contains no less than k + 1 vertices, G is k-vertex connected if it is still connected after deleting any k - 1 vertices from it. If G is not k-edge/vertex connected, it contains k-edge/vertex connected components. Given a graph G and an integer k, a pair of vertices u and v in G is said to be k-edge connected if the removal of any k - 1 edges from G leaves u and vconnected. This is an equivalence relationship and written as \equiv_k , i.e., if a pair of vertices x and y is k-edge connected, we write $x \equiv_k y$. The vertices in G are partitioned by this relationship into equivalence classes and each equivalence class is a k-edge connected component in G.

It seems much more complicated to define the *k*-vertex *connected component* in G with k > 3. Here, we only define this concept for $1 \le k \le 3$, which will be used in this paper. Given a graph G(V, E), let E_1, E_2, \ldots, E_h be an edge partition of E into equivalence classes such that two edges e' and e'' are in the same equivalence class if and only if either 1) e' = e'', or 2) there is a simple cycle in G containing both e' and e''. For all i, $1 \le i \le h$, let V_i be the set of endpoints of the edges in E_i . Then, the subgraph $G_i = (V_i, E_i)$ is a biconnected component (or 2vertex connected component) in G. An edge contained in no cycle is a biconnected component by itself and is referred to as a bridge in G. A graph G without including any bridge is called 2-edge connected. If G contains bridges, then the removal of a bridge disconnects the graph. Usually, 2-vertex connectivity is called biconnectivity and a biconnected component is called a block. If, for every pair of vertices, there are two vertex disjoint paths in G between them, then G is called a biconnected graph. It is well known [1] that G is not biconnected if and only if there is a vertex $v \in V$ whose removal disconnects G. Such a *v* is called an *articulation point* of *G*.

Let $\{a, b\}$ be a pair of vertices in a biconnected graph G. Suppose the edges in G are partitioned into equivalence classes E_1, E_2, \ldots, E_p such that two edges, which lie on a common path not containing any vertex of $\{a, b\}$ except as an endpoint, are in the same class. The classes E_i are called the separation classes of G with respect to $\{a, b\}$. If there are at least two separation classes, then $\{a, b\}$ is a *separation pair* of G unless 1) there are exactly two separation classes, one class consists of a single edge, or 2) there are exactly three classes, each consisting of a single edge. If G is a biconnected graph that does not contain a separation pair, then G is triconnected. Let $\{a, b\}$ be a separation pair of G and E_1, E_2, \ldots, E_p be the separation classes of G with respect to $\{a, b\}$. Let $E' = \bigcup_{i=1}^{k} E_i$ and $E'' = \bigcup_{i=k+1}^{p} E_i$ and $|E'| \ge 2$ and $|E''| \ge 2$. Let $G_1 = (V(E'), E' \cup \{(a, b)\})$ and $G_2 = (V(E''), E'' \cup \{(a, b)\})$, where V(E') and V(E'') are the sets of endpoints of edges in E' and E''. G_1 and G_2 are called the *split graphs* of G with respect to $\{a, b\}$. Replacing G by the two split graphs is called *splitting* G. The new edges (a, b) added to both G_1 and G_2 are called *virtual edges*. If G is biconnected, then any split graph of G is also biconnected. Suppose G is split, the split graphs are split and so on, until no more splits are possible. The graphs constructed in this way are called *split components* of G. The split components of G are of three types: triple bonds of form $(\{a, b\}, \{(a, b), (a, b), (a, b)\})$, triangles of form $(\{a, b, c\}, \{(a, b), (a, c), (b, c)\})$, and triconnected graphs. The split components of G are called *triconnected* components in G. A triconnected component usually is also called 3-vertex connected component and 3-vertex connectivity is called triconnectivity.

An inverted tree $T(V, E_T)$ is a directed tree rooted at a distinguished vertex $r, r \in V$, such that for every other vertex v ($v \neq r$) there is a parent pointer pointing to the parent $F_T(v)$ of v, an edge $\langle v, F_T(v) \rangle \in E_T$, and $F_T(r) = r$. Given an inverted tree T and any two vertices v and u in it, the *lowest common ancestor* LCA(u, v) of u and v is the first common vertex on the paths from u (v) to the root. It is known that every LCA query in T can be answered in O(1) time using one processor on an EREW PRAM

provided that *T* is properly preprocessed. The preprocessing takes $O(\log n)$ time and uses O(n) processors on an EREW PRAM if *T* contains *n* vertices [42]. Note that all the trees used in this paper are maintained as inverted trees.

2.2 Transformation Technique of Graphs

We now introduce a technique to transform a general graph with arbitrary degree into a graph with bounded degree. It is well known [22] that a graph G(V, E) can be transformed to another graph G'(V', E') such that every vertex in G' has no degree greater than three. The transformation proceeds as follows: For each vertex u of degree $d \ge 4$ in G, assume that $v_0, v_1, \ldots, v_{d-1}$ are the adjacent vertices of u in G. Then, G' is generated by replacing u with new vertices $u_0, u_1, \ldots, u_{d-1}$, adding edges $(u_i, u_{(i+1) \mod d})$ and edges (u_i, v_i) to G', $0 \le i \le d - 1$. One important property of the transformation is that it preserves the 2-edge connectivity of G [19]. That is, if two vertices are in a 2-edge connected component in G, then they will also be in the same 2-edge connected component in G'. Otherwise, they are still not in the same 2-edge connected component in G'.

It is clear that G' contains O(m + n) edges and vertices and each vertex has degree no greater than three. Without loss of generality, in Section 3, we assume G(V, E) is a graph with O(m + n) vertices and each vertex has a degree no greater than three when dealing with the 2-edge connectivity. Otherwise, the transformation can be applied, which takes $O(\log n)$ time using O(m + n) processors on an EREW PRAM.

2.3 The Sparse *k*-Edge (*k*-Vertex) Certificate

Let G(V, E) be an arbitrary graph. An *edge separator* in *G* is such a subset of edges that the removal of the edges in it will disconnect *G*. A *sparse k-edge certificate* of *G* is a subgraph *H* in *G* containing O(kn) edges and every edge separator in *H* of size less than *k* is also an edge separator in *G*. A *sparse k-vertex certificate* of *G* can be defined similarly. In the following, we show how to find a sparse *k*-edge/ *k*-vertex certificate of *G* [38], [4].

Let T_1 be a maximal spanning forest in G and T_i be a maximal spanning forest of graph $G_i = G - \bigcup_{j=1}^{i-1} T_j$ for i > 1. Denote by $U_i = \bigcup_{j=1}^{i} T_j$, the union of the *i* maximal spanning forests T_1, T_2, \ldots, T_i . The graph U_k is a *sparse k-edge certificate* of G and has the following property:

Lemma 1 [38], [39]. The graph U_k is *l*-edge connected if and only if G is *l*-edge connected for any integer *l* with $1 \le l \le k$.

Lemma 1 always holds no matter whether G is a simple graph or not. We now introduce the *scan-first search technique* [4] on G which will be used to find a sparse k-vertex certificate of G.

Definition 1. A scan-first search in a connected undirected graph G(V, E) starting from a specified vertex r is a systematic way of visiting the vertices in G. To scan a vertex is to visit all previously unvisited adjacent vertices of that vertex. At the beginning of the search, only r is visited. Then, the search iteratively scans an already visited but unscanned vertex until all vertices are scanned. If *G* is disconnected, a scan-first search spanning forest of *G* is then obtained by applying the scan-first search on each connected component in *G*. Cheriyan et al. [4] presented an efficient parallel algorithm for finding a scan-first search forest in *G*, which proceeds as follows: It first finds an arbitrary spanning tree *T* in *G*. It then traverses *T* by labeling the vertices with their preorder numberings in *T*. It finally constructs the following scan-first tree ST: Let b(v) be the parent of v in ST and N(v) be the set of adjacent vertices of v in *G*; b(v) is such a vertex in N(v) that it has the smallest preorder numbering for every $v \in V$, $v \neq r$.

Lemma 2 [4]. For a graph with n vertices and m edges, a scanfirst search spanning forest can be found in $O(\log n)$ time using C(n,m) processors on a CRCW PRAM, where C(n,m)is the number of processors used to find spanning trees in all connected components in $O(\log n)$ time.

Note that, currently, the best result of C(n,m) is $O((m+n)\alpha(m,n)/\log n)$ on a CRCW PRAM [6]. With the above preparation, the sparse *k*-vertex certificate of *G* is obtained as follows: For i = 1, 2, ..., k, let E_i be the edge set of a scan-first search spanning forest in graph $G_{i-1} = (V, E - (E_1 \cup ... \cup E_{i-1}))$ and $G_0 = (V, E)$. Then, the graph $C_k = E_1 \cup ... \cup E_k$ is a sparse *k*-vertex certificate of *G*, which contains *n* vertices and no more than k(n-1) edges.

Lemma 3 [4]. The graph C_k defined is *l*-vertex connected if and only if G is *l*-vertex connected for any integer *l* with $1 \le l \le k$.

2.4 Sparsification Technique

Eppstein et al.'s sparsification technique [10] is based on a *certificate* concept. Before introducing the technique, we reproduce some of the definitions which are fundamental to the technique.

- **Definition 2.** For any graph property \mathcal{P} and graph G(V, E), a certificate for G is a graph G' such that G has the property \mathcal{P} if and only if G' has the property.
- **Definition 3.** For any graph property \mathcal{P} and a graph G(V, E), a strong certificate for G is a graph G' on the same vertex set such that, for any $H, G \cup H$ has the property \mathcal{P} if and only if $G' \cup H$ has the property.
- **Definition 4.** Let \mathcal{A} be a function mapping graphs to strong certificates. Then, \mathcal{A} is stable if it has the following two properties: 1) for any graphs G and H, $\mathcal{A}(G \cup H) = \mathcal{A}(\mathcal{A}(G) \cup H)$; 2) for any graph G and edge $e \in E$, $\mathcal{A}(G - \{e\})$ differs from $\mathcal{A}(G)$ by O(1) edges.

The sparsification technique is described as follows: We partition the edges in G(V, E) into $\lceil m/n \rceil$ groups, all except one contain exactly n edges. The remaining group, called the *small* group may contain between 1 and n edges. When inserting an edge into G, we place it into the small group. When deleting an edge from a group, we move another edge from the small group into the group, to keep the group size invariant (i.e., n edges exactly). If deleting the last edge in the small group, we remove the small group entirely and, if inserting an edge to the small group while there are already n edges in it, we start a new small group. Having finished edge grouping, we now establish a complete

binary tree \mathcal{BT} with $\lceil m/n \rceil$ leaf nodes, which correspond to the $\lceil m/n \rceil$ groups. \mathcal{BT} is called the *sparsification tree* in which every *node* corresponds to a subgraph formed by the edges in the graphs at the leaf nodes of \mathcal{BT} that are the descendants of the node. For every node in \mathcal{BT} , we maintain a sparse certificate \mathcal{P} for it. The sparse certificate at a leaf node is obtained by applying the procedure of finding a sparse certificate to the subgraph consisting of the edges in the corresponding group. The sparse certificate at a nonleaf node is the union of the sparse certificates of its two children nodes.

Each update in *G* will cause either a single leaf node of \mathcal{BT} to split or two leaf nodes to merge in the worst case. Since only the edge insertions and deletions are allowed, the number of edges in every group except the small group, will never be changed. When an edge is inserted or deleted, only O(1) groups are involved. For each of these groups and its $O(\log \frac{m}{n})$ ancestor nodes in \mathcal{BT} , we maintain the sparse certificates of these nodes dynamically. As results, the sparse certificate of *G* is the sparse certificate at the root of \mathcal{BT} .

2.5 The Reduction Technique

A reduction technique due to Galil and Italiano [18] can reduce a *k*-edge connectivity problem in a graph G = (V, E)to a *k*-vertex connectivity problem in another graph $\varphi_k(G) =$ $(\varphi(V), \varphi(E))$ and $\varphi_k(G)$ is constructed as follows: For every vertex *v* in *G*, there are k - 2 vertices

$$\varphi(v_1), \varphi(v_2), \ldots, \varphi(v_{k-2})$$

in $\varphi_k(G)$, referred to as *node-vertices* of $\varphi_k(G)$. For every edge e in G, there is a vertex $\varphi(e)$ in $\varphi_k(G)$, referred to as *arc-vertex* of $\varphi_k(G)$. Let v be any vertex of G and $u_0, u_1, \ldots, u_{d-1}$ be the adjacent vertices of v in G, and $e_i = (v, u_i)$ be an edge in G, $0 \le i \le d-1$. Then, edges $(\varphi(e_i), \varphi(e_{(i+1) \mod d}))$ and $(\varphi(e_i), \varphi(v_j))$ are in $\varphi_k(G)$, $0 \le i \le d-1$, $0 \le j \le k-2$. Notice that $\varphi_k(G)$ is obtained from G by replacing v and its incident edges $e_0, e_1, \ldots, e_{d-1}$ by a wheel with k-2 hubs

$$\varphi(v_1), \varphi(v_2), \ldots, \varphi(v_{k-2})$$

and vertices $\varphi(e_0), \varphi(e_1), \ldots, \varphi(e_{d-1})$. This implies that $\varphi_k(G)$ has O(m + kn) vertices and O(km) edges. Clearly, $\varphi_k(G)$ can be constructed in $O(\log n)$ time using O(k(m + n)) processors on an EREW PRAM if G contains O(n) vertices and O(m) edges. The graph $\varphi_k(G)$ has the following property:

Theorem 1 [18]. G(V, E) is k-edge connected if and only if $\varphi_k(G)(\varphi(V), \varphi(E))$ is k-vertex connected.

Following Theorem 1, to check whether two vertices, x and y in G, are k-edge connected can be done through checking whether their corresponding node-vertices $\varphi(x_1)$ and $\varphi(y_1)$ in $\varphi_k(G)$ are k-vertex connected.

3 MAINTAINING 2-EDGE CONNECTED COMPONENTS IN CONNECTED GRAPHS

For the sake of simplicity, in this section, we assume that G(V, E) is connected and contains O(m + n) vertices and edges. Every vertex in *G* has a degree no greater than three.



Fig. 1. A graph, its spanning tree, and the vertex cluster partition. (a) A graph G(V, E). (b) A spanning tree and a vertex cluster partition of G.

Otherwise, the transformation technique will be applied in order to obtain a graph with the degree no more than three. Fig. 1a shows such a graph.

3.1 Overview

We present our results in the following order: First, we decompose G into a number of disjoint connected subgraphs and build data structures for each of them to keep the information on the subgraph. Also, we build a data structure to keep the information among the subgraphs, which is referred to as the *compact representation* $\hat{\mathcal{G}}$ of G. $\hat{\mathcal{G}}$ is the super graph of G, defined as follows: Each vertex in \mathcal{G} corresponds to a subgraph of G. There is an edge between two vertices in \mathcal{G} if there are edges in G between the vertices in the two corresponding subgraphs. For any two subgraphs in G, we use two data structures to keep the edge information between them. Then, we assume that all the data structures for 2-edge connected components in each subgraph have been already built. Now, consider a query which asks whether two given vertices x and y are in the same 2-edge connected component in G. To answer this query, we proceed as follows: Let *x* be in C_x and *y* be in C_y . We check whether C_x and C_y are in the same 2-edge connected component in $\hat{\mathcal{G}}$. If yes, then x and y are in the same 2-edge connected component in G. Otherwise, for every articulation point C in G separating C_x from C_y , we examine the induced subgraph of G by the vertices in C to see whether there is a potential bridge internal to the subgraph separating x from y in G. If yes, then x and y are not in the same 2-edge connected component in G. Finally, we deal with the fully dynamic maintenance of the data structures when there is an edge insertion or deletion. Note

that each insertion/deletion only involves updating a constant number of subgraphs and $\hat{\mathcal{G}}$. Accordingly, we only update those involved subgraphs and $\hat{\mathcal{G}}$ instead of the entire graph *G*, which is much cheaper in terms of the maintenance cost because of the smaller sizes of these subgraphs and $\hat{\mathcal{G}}$ compared to *G*. Therefore, it is possible to reduce the maintenance cost to respond the updates and queries.

The rest of this section is organized as follows: Section 3.2 introduces the data structures that the proposed algorithm uses and Section 3.3 presents algorithms for constructing the data structures. Section 3.4 proposes an algorithm for query processing, using the proposed data structures. Section 3.5 shows how to maintain the data structures efficiently when there are edge insertions and deletions.

3.2 Data Structures for 2-Edge Connected Components

The proposed algorithm, basically, is a parallel implementation of Galil and Italiano's algorithm [19]. Therefore, most data structures used in their algorithm are adopted. Let G(V, E) be a connected graph with the maximum degree three. The *vertex cluster partition* of *G* is described as follows: Following the vertex clustering idea by Frederickson [15], the vertex set *V* is partitioned into V_1, V_2, \ldots, V_p disjoint subsets such that 1) $\lfloor K/2 \rfloor \leq |V_i| \leq \lceil 3K/2 \rceil$ for all $i, 1 \leq i \leq p$ and 2) the induced subgraph of *G* by the vertices in V_i is connected, where *K* is a parameter to be determined later. $V = \bigcup_{i=1}^p V_i$ and p = O(m/K) because *G* contains O(m + n)vertices and each vertex in it has a degree no greater than three. Each vertex subset V_i is a *vertex cluster* (or cluster for short) of *G*. The size of a cluster is the number of vertices in



Fig. 2. The induced subgraph $\mathcal{G}(\mathcal{C}_1)$ and the full representation $\mathcal{F}(\mathcal{C}_1)$ of \mathcal{C}_1 . (a) The induced subgraph $\mathcal{G}(\mathcal{C}_\infty)$ of G by the vertices in cluster \mathcal{C}_1 . (b) The full presentation $\mathcal{F}(\mathcal{C}_1)$ of a vertex cluster \mathcal{C}_1 where $\mathcal{C}_1 = \{1, 2, 3, 4, 5, 11\}$.

it. However, the size of each cluster is O(K). For example, a spanning tree in *G* is shown in Fig. 1b. The vertex set of *G* is partitioned into two vertex clusters $C_1 = \{1, 2, 3, 4, 5, 11\}$ and $C_2 = \{6, 7, 8, 9, 10\}$ with K = 4.

3.2.1 Data Structures Pertinent to a Vertex Cluster C

Induced Subgraph $\mathcal{G}(\mathcal{C})$ of \mathcal{C} . Let \mathcal{C} be a vertex cluster. An edge in G with both its endpoints in \mathcal{C} is said to be *internal* to \mathcal{C} while an edge in G with only one endpoint in \mathcal{C} is said to be *incident* to \mathcal{C} . The induced subgraph of G by the vertices in \mathcal{C} is denoted by $\mathcal{G}(\mathcal{C})$. For the above example, $\mathcal{G}(\mathcal{C}_1)$ is shown in Fig. 2a. Edge (2, 4) in G is internal to \mathcal{C}_1 while edge (4, 6) in G is incident to \mathcal{C}_1 .

Full Representation $\mathcal{F}(\mathcal{C})$ of \mathcal{C} . The *full representation* $\mathcal{F}(\mathcal{C})$ of a vertex cluster \mathcal{C} is a simple graph, which consists of $\mathcal{G}(\mathcal{C})$ a vertex for each vertex cluster adjacent to \mathcal{C} and the edges incident to \mathcal{C} . In other words, the vertex set of $\mathcal{F}(\mathcal{C})$ consists of the vertices internal to \mathcal{C} and the corresponding vertices of the vertex clusters which are incident to \mathcal{C} . The edge set of $\mathcal{F}(\mathcal{C})$ is composed of all the edges in G internal and incident to \mathcal{C} . Fig. 2b shows $\mathcal{F}(\mathcal{C}_1)$, where \mathcal{C}_2 is the cluster adjacent to \mathcal{C}_1 .

Tree Representation T(C) of C. The *tree representation* T(C) of a vertex cluster C is a tree in which the edges are the bridges in $\mathcal{G}(C)$ and the vertices are the 2-edge connected components in $\mathcal{G}(C)$. For a given $\mathcal{G}(C_1)$, in Fig. 2a, the tree representation $T(C_1)$ of C_1 is shown in Fig. 3a, where vertices 1, 2, and 3 are in a 2-edge connected component

called *a*, vertices 4, 5, and 11 are in the 2-edge connected components called *b*, *c*, and *d*, respectively.

3.2.2 Edge Set Presentation between Two Vertex Clusters

Given two clusters C_i and C_j , let $E_{i,j} = \{e_1, e_2, \dots, e_h\}$ be a subset of *E* in which the two endpoints of every edge are in C_i and C_{i} , respectively. The edges in $E_{i,i}$ are represented by two data structures: One is for C_i , named $E(C_i, C_j)$, and the other is for C_j , named $E(C_j, C_i)$. Let X be an endpoint set of the edges in $E_{i,j}$, which is a subset of C_j , then $E(\mathcal{C}_i, \mathcal{C}_j) = \{y \mid y \text{ is a 2-edge connected component vertex}\}$ in $\mathcal{T}(\mathcal{C}_i)$, including x and $x \in X$ }. Therefore, $E(\mathcal{C}_i, \mathcal{C}_i)$ is a vertex subset of $\mathcal{T}(\mathcal{C}_i)$. In other words, $E(\mathcal{C}_i, \mathcal{C}_i)$ contains those vertices of $\mathcal{T}(\mathcal{C}_i)$ that, for each such vertex *u*, there is at least an edge $(x, y) \in E_{i,j}$ incident to u, where $x \in C_i$, $y \in u$, and u is a 2-edge connected component in $\mathcal{G}(\mathcal{C}_i)$. $E(\mathcal{C}_i, \mathcal{C}_i)$ is represented by a balanced binary search tree and the key to access a vertex in it is the preorder numbering of the vertex in $\mathcal{T}(\mathcal{C}_i)$. For the example, in Fig. 1b, there are two vertex clusters C_1 and C_2 . $E_{2,1} = \{(4,6), (5,6), (11,10)\}.$ $E(\mathcal{C}_2, \mathcal{C}_1) = \{b, c, d\}$. Fig. 3b shows $E(\mathcal{C}_2, \mathcal{C}_1)$, where *preorder*(*x*) is the preorder numbering of *x* in $\mathcal{T}(\mathcal{C}_1)$.

3.2.3 Super Graphs and Equivalent Graphs

The super graph $\hat{\mathcal{G}}$ of *G* is a multigraph defined as follows: Each vertex in $\hat{\mathcal{G}}$ is a vertex cluster. There is an edge



Fig. 3. (a) The tree representation $\mathcal{T}(\mathcal{C}_1)$ of \mathcal{C}_1 . (b) $E(\mathcal{C}_2, \mathcal{C}_1)$ and its binary search tree representation.



Fig. 4. The super graph $\hat{\mathcal{G}}$ of a graph G.

between two vertex clusters C_i and C_j if $|E_{i,j}| = 1$ and there are two edges between them if $|E_{i,j}| \ge 2$. For example, the graph in Fig. 4 is a super graph of a graph G.

A simple graph $\overline{\mathcal{G}}(\mathcal{V}, \mathcal{E})$, derived from $\hat{\mathcal{G}}$, is defined as follows:

 $\mathcal{V} = \{\mathcal{C}_i \mid \mathcal{C}_i \text{ is a vertex cluster}\}$

 $\cup \{v_{i,j,1}, v_{i,j,2} | \text{there are two edges in } \hat{\mathcal{G}} \text{ between } \mathcal{C}_i \text{ and } \mathcal{C}_j \},\$

where $v_{i,j,k}$ is an *artificial vertex*, k = 1, 2.

 $\mathcal{E} = \{ (\mathcal{C}_i, \mathcal{C}_j) | \text{there is only one edge in } \hat{\mathcal{G}} \text{ between } \mathcal{C}_i \text{ and } \mathcal{C}_j \}$ $\cup \{ (\mathcal{C}_i, v_{i,j,k}), (v_{i,j,k}, \mathcal{C}_j) | \text{there are two edges between}$ $\mathcal{C}_i \text{ and } \mathcal{C}_i \text{ in } \hat{\mathcal{G}} \text{ and } k = 1, 2 \}.$

Thus, the vertex set of $\hat{\mathcal{G}}$ is a subset of the vertex set of $\overline{\mathcal{G}}$. Graph $\overline{\mathcal{G}}(\mathcal{V}, \mathcal{E})$ is called the *equivalent* graph of $\hat{\mathcal{G}}$ in the sense that they both have the same sets of articulation points and bridges. Assume that $\hat{\mathcal{G}}$ contains m' edges and n' vertices. Then, $\overline{\mathcal{G}}(\mathcal{V}, \mathcal{E})$ contains no more than 2m' edges and n' + m' vertices. For example, the graph in Fig. 5a is the equivalent graph of a super graph shown in Fig. 4.

From now on, we assume that there is no big difference between $\hat{\mathcal{G}}$ and $\overline{\mathcal{G}}$ with respect to the articulation points and bridges. The purpose that we here introduce the equivalent graph $\overline{\mathcal{G}}$ for $\hat{\mathcal{G}}$ is to make Tarjan and Vishkin's parallel algorithm for finding biconnected components can be applied, which is only applicable for simple graphs.

3.3 Construction of Data Structures

We now show how to construct the defined data structures in parallel. In order to obtain a vertex cluster partition of G, a spanning tree $T(V, E_T)$ of G is generated, using any efficient parallel algorithm such as the algorithm by Awerbuch and Shiloach [2]. Assume T has already been built and stored as an inverted tree. Our approach for the vertex cluster partition of G is based on the following lemma due to Lipton and Tarjan [37].

Lemma 4 [37]. Given an n-vertex tree with the maximum degree three, there is a vertex partitioning by removing exactly one edge from the tree. As the result, the tree is separated into two subtrees and the number of vertices in each subtree is between |n/3| and [2n/3].

The vertex cluster partition of G then can be obtained as follows:

- **Lemma 5.** Given an inverted tree T with m vertices and the maximum degree three, partition its vertices into cm/K vertex clusters such that every vertex cluster contains O(K) vertices. This can be done in $O(\log(m/K))$ time using O(m) processors on an EREW PRAM, where $m \ge 3K/2$, cm/K is the number of clusters, and c is constant.
- **Proof.** Let \mathcal{F} be a forest of inverted trees. Initially, $\mathcal{F} = \{T\}$. For each tree $T' \in \mathcal{F}$, we proceed as follows: If the number of vertices n_1 in T' is between K/2 and 3K/2, we do nothing about it. Otherwise, traverse T' by a Eulerian tour and make it into a linked list L. Delete such an edge in L that the two resulting subtrees have almost the same



Fig. 5. The equivalent graph $\overline{\mathcal{G}}$ of $\hat{\mathcal{G}}$ and the resulting graph through replacing a vertex cluster \mathcal{C}_2 in $\overline{\mathcal{G}}$ by $\mathcal{F}(\mathcal{C}_2)$. (a) The equivalent graph $\overline{\mathcal{G}}$ of a super graph $\hat{\mathcal{G}}$. (b) Replace \mathcal{C}_2 and its artificial adjacent vertices by $\mathcal{F}(\mathcal{C}_2)$ in $\overline{\mathcal{G}}$.

number of vertices (one has $n_1/3$ vertices and the other has $2n_1/3$ vertices in the worst case). The existence of such a partition is guaranteed by Lemma 4. Remove T'from \mathcal{F} and add the two new trees to \mathcal{F} . Repeat the above procedure on \mathcal{F} until the size of every tree in \mathcal{F} is between K/2 and 3K/2. There are $O(\log(m/K))$ calls for the procedure. As a result, every tree in \mathcal{F} is a vertex cluster of G because its size is between K/2 and 3K/2and the tree is a connected subgraph in G. Therefore, the vertex cluster partition can be done in $O(\log m)$ time using O(m) processors on an EREW PRAM [31] because the size of the resulting tree is reduced by one third at least for each procedure call, compared with the size of its immediate predecessor. \Box

Note that the restricted vertex cluster partition by Frederickson cannot be applied here because it uses the depth-first-search technique which seems highly sequential. Furthermore, although the algorithms for finding a spanning tree and a vertex cluster partition take $O(\log n)$ time using O(m + n) processors and $O(\log(m/K))$ time using O(m) processors, respectively, they are only done once. In the design of dynamic algorithms, the time spent for this part is called *preprocessing time*, which is not taken into account as part of the dynamic maintenance cost for the operations including insertions, deletions, and queries.

Given a vertex cluster partition of G, the defined data structures are constructed as follows: For a vertex cluster C, the graph $\mathcal{G}(\mathcal{C})$ can be constructed in O(1) time using O(K) processors on a CRCW PRAM because $\mathcal{G}(\mathcal{C})$ contains O(K) vertices and edges and the degree of each vertex in it is constant. Since each vertex cluster C contains O(K) vertices and each vertex in G has a degree no greater than three, then the full representation $\mathcal{F}(\mathcal{C})$ of C contains O(K) vertices and edges. $\mathcal{F}(\mathcal{C})$ can be constructed in $O(\log K)$ time with O(K) processors using the adjacency lists of those vertices in G that are also in C and the vertex cluster partition of G. Note that the maximum degree of the vertices in $\mathcal{F}(\mathcal{C})$ is O(K), not a constant. The tree representation $\mathcal{T}(\mathcal{C})$ of C can be obtained by the following lemma:

- **Lemma 6.** The full tree representation T(C) of a vertex cluster C can be constructed in $O(\log K)$ time using O(K) processors on a CRCW PRAM.
- **Proof.** We here provide a constructive proof. We know that $\mathcal{G}(\mathcal{C})$ is a simple graph. First, find all biconnected components in $\mathcal{G}(\mathcal{C})$ by applying Tarjan and Vishkin's algorithm [44]. This takes $O(\log K)$ time with O(K) processors on a CRCW PRAM because $\mathcal{G}(\mathcal{C})$ contains O(K) vertices and edges. Then, find all bridges in $\mathcal{G}(\mathcal{C})$. The edge in a biconnected component consisting of one edge only is a bridge in $\mathcal{G}(\mathcal{C})$, which can be found in O(1) time with O(K) processors after all biconnected components in $\mathcal{G}(\mathcal{C})$ have been found. Let *B* be the set of bridges in $\mathcal{G}(\mathcal{C})$. Let $\mathcal{G}'(\mathcal{C})$ be the resulting graph after removing the edges in *B* from $\mathcal{G}(\mathcal{C})$, which can be done in O(1) time with O(K) processors because of the constant degree of each vertex in $\mathcal{G}(\mathcal{C})$. Finding connected components in $\mathcal{G}'(\mathcal{C})$, which in fact are the 2-edge connected components (*BCs*) in $\mathcal{G}(\mathcal{C})$ takes $O(\log K)$ time using O(K) processors on a CRCW PRAM [43]. Finally,

use the vertex in *BC* with the smallest index to label all the vertices in it. The *BC* is then represented by that vertex. Thus, all the vertices in the *BC* have a unique representative. Let D() be such a labeling function, i.e., D(u) = v for all $u \in BC$, where v is the vertex in *BC* with the smallest index.

Let T'' be a spanning tree of $\mathcal{G}(\mathcal{C})$. We now construct the tree $\mathcal{T}(\mathcal{C})$. Let $F_{\mathcal{T}(\mathcal{C})}(v)$ be the parent of v in $\mathcal{T}(\mathcal{C})$. Then, the vertex set of $\mathcal{T}(\mathcal{C})$ consists of all 2-edge connected components in $\mathcal{G}(\mathcal{C})$. The edges in $\mathcal{T}(\mathcal{C})$ are defined as follows: Initially, set $F_{\mathcal{T}(\mathcal{C})}(D(v)) := v$ for all $v \in BC$. Then, for every edge (x, y) in B, assign $F_{\mathcal{T}(\mathcal{C})}(D(x)) := D(y)$ if y is the parent of x in T''; $F_{\mathcal{T}(\mathcal{C})}(D(y)) := D(x)$ otherwise. This can be done in O(1) time with O(K) processors because B contains O(K) edges. \Box

Given an inverted tree $\mathcal{T}(\mathcal{C}_i)$ of \mathcal{C}_i , assigning every vertex in $\mathcal{T}(\mathcal{C}_i)$ a preorder numbering takes $O(\log K)$ time using O(K) processors on an EREW PRAM [31]. Then, all $E(\mathcal{C}_j, \mathcal{C}_i)$ s can be constructed in $O(\log K)$ time using O(K) processors by sorting the preorder numbering of vertices in increasing order because the degree of each vertex in $\mathcal{G}(\mathcal{C}_i)$ is no more than three, assuming that \mathcal{C}_j is a vertex cluster adjacent to \mathcal{C}_i .

The super graph $\hat{\mathcal{G}}$ can be constructed in O(1) time using $O((m/K)^2)$ processors because $\hat{\mathcal{G}}$ contains O(m/K) vertices. Thus, the equivalent graph $\overline{\mathcal{G}}$ of $\hat{\mathcal{G}}$ can be constructed in the same time and processor bounds because $\overline{\mathcal{G}}$ contains $O((m/K)^2)$ vertices and edges.

Consequently, the data structures pertinent to a vertex cluster C can be maintained in $O(\log K)$ time using O(K) processors and the data structures for $\hat{\mathcal{G}}$ and $\overline{\mathcal{G}}$ can be maintained in $O(\log(m/K) + \log K)$ time using $O((m/K)^2)$ processors on a CRCW PRAM, which will be proven in Section 3.4.

3.4 The Algorithm for Queries

Given a pair of vertices x and y, the query is about whether x and y are in the same 2-edge connected component in G. Instead of using G, an auxiliary graph $\mathcal{G}_{x,y}$ will be used to answer the query. Let \mathcal{C}_x and \mathcal{C}_y be the vertex clusters containing x and y, respectively. The simple graph $\mathcal{G}_{x,y}$ is obtained from $\overline{\mathcal{G}}$ through replacing vertices C_x , C_y and their adjacent artificial vertices and the edges incident to them by the full representations $\mathcal{F}(\mathcal{C}_x)$ and $\mathcal{F}(\mathcal{C}_y)$ of \mathcal{C}_x and \mathcal{C}_y . For example, Fig. 5b shows the resulting graph after replacing C_2 and its adjacent artificial vertices in $\overline{\mathcal{G}}$ in Fig. 5a by its full representation $\mathcal{F}(\mathcal{C}_2)$. Now, assume that x and y are not in the same 2edge connected component in G. Then, there must be at least one bridge $e \in E$ in G separating x from y and any such bridge can be one of the following two types: 1) If eis either internal to C_x , or C_y , or an edge between two different clusters; 2) if e is internal to a vertex cluster Cwith $C \neq C_x$ and $C \neq C_y$. To check these two types of bridges, we have the following theorem:

Theorem 2 [19]. Let x and y be any two vertices in G. Then, the following is true: 1) e is a Type 1 bridge in G if and only if the corresponding edge is a bridge in $G_{x,y}$; 2) if e is a Type 2 bridge

in G separating x from y and e is internal to a cluster C $(C \neq C_x \text{ and } C \neq C_y)$, then C is an articulation point in $\mathcal{G}_{x,y}$.

Following Theorem 2, the checking algorithm proceeds as follows: It checks Type 1 bridges in $\mathcal{G}_{x,y}$ first. If there is no Type 1 bridge, then it checks Type 2 bridges. If no Type 2 bridge exists, then, x and y are in the same 2-edge connected component in G; otherwise, x and y are in the different 2-edge connected components in G. The rest of this section, therefore, focuses on checking the two types of bridges. However, Type 1 bridge can be easily detected by the following lemma:

- **Lemma 7.** Given a simple connected graph with n_1 vertices and m_1 edges, to determine whether there exist bridges in any path between two given vertices takes $O(\log n_1)$ time using $O(n_1 + m_1)$ processors on a CRCW PRAM.
- Proof. First, find all biconnected components in the graph by applying Tarjan and Vishkin's algorithm and identify those biconnected components consisting of one edge only (the edges are the bridges of the graph). Let B be the set of such edges. Then, find a spanning tree T in the graph by applying any efficient spanning tree algorithm and identify the edges in path π in T between the two given vertices. Finally, if there is an edge in both *B* and π , then it is a bridge of the graph separating the two vertices. The time spent for finding biconnected components is $O(\log n_1)$ and the number of processors used is $O(m_1 + n_1)$. Constructing B takes O(1) time using $O(n_1)$ processors. T can be found in $O(\log n_1)$ time using $O(m_1 + n_1)$ processors. All the edges in π can be identified in $O(\log n_1)$ time using $O(n_1)$ processors. Therefore, all the bridges in the graph separating x from y can be found in $O(\log n_1)$ time using $O(m_1 + n_1)$ processors on a CRCW PRAM.

As a result, Type 1 bridges in *G* between *x* and *y* can be found in $O(\log K + \log(m/K))$ time using $O(K + (m/K)^2)$ processors on a CRCW PRAM because $\mathcal{G}_{x,y}$ contains $O(K + (m/K)^2)$ vertices and edges by Lemma 7.

3.4.1 Type 2 Bridge Checking

We now deal with Type 2 bridges. We first introduce the following notions and notations for later use:

Let a vertex cluster C be an articulation point in $\mathcal{G}_{x,y}$ separating x and y. Then, C is also an articulation point in both $\overline{\mathcal{G}}$ and $\hat{\mathcal{G}}$ separating \mathcal{C}_x from \mathcal{C}_y by the definition of equivalent graphs. Let $AC = \{Q_1, Q_2, \dots, Q_l\}$ be a set of vertex clusters in which each vertex cluster Q_j in $\hat{\mathcal{G}}$ is adjacent to $C, 1 \leq j \leq t$ and $t \geq 2$. Let $\mathcal{W}_1, \mathcal{W}_2, \dots, \mathcal{W}_s$ be the s connected components in the resulting graph $\hat{\mathcal{G}} - \{C\}$ after the removal of C and the edges incident to it from $\hat{\mathcal{G}}$, where each vertex in \mathcal{W}_i represents a vertex cluster, $1 \leq i \leq s$. Without loss of generality, let $\mathcal{W}_1 = \{\mathcal{X}_1, \mathcal{X}_2, \dots, \mathcal{X}_p\}$ be the connected component containing $\mathcal{C}_x, \mathcal{W}_2 = \{\mathcal{Y}_1, \mathcal{Y}_2, \dots, \mathcal{Y}_q\}$ be the connected component containing \mathcal{C}_y , and

$$\mathcal{W}_i = \{\mathcal{Z}_{i,1}, \mathcal{Z}_{i,2}, \dots, \mathcal{Z}_{i,r(i)}\}$$

be a connected component containing the vertex clusters in $AC - (W_1 \cup W_2)$, $3 \le i \le s$, where \mathcal{X}_{l_1} , \mathcal{Y}_{l_2} , and $\mathcal{Z}_{i,j}$ are the

vertex clusters, $1 \le l_1 \le p$, $1 \le l_2 \le q$, $1 \le j \le r(i)$, and $3 \le i \le s$. Clearly, $p + q + \sum_{i=3}^{s} r(i) = t$. Given the tree representation $\mathcal{T}(\mathcal{C})$ of \mathcal{C} and an edge e in it, let ${\mathcal T}_1({\mathcal C})$ and ${\mathcal T}_2({\mathcal C})$ be the resulting subtrees after the removal of e from $\mathcal{T}(\mathcal{C})$. The connected component \mathcal{W}_i $(= \{Z_{i,1}, Z_{i,2}, \dots, Z_{i,r(i)}\})$ defined is said to be *compatible* with *e* if all the edges between $\mathcal{Z}_{i,j}$ and \mathcal{C} , $1 \leq j \leq r(i)$, are incident to either $\mathcal{T}_1(\mathcal{C})$ or $\mathcal{T}_2(\mathcal{C})$, but not to both, $3 \le i \le s$. For example, the connected component W_3 , shown in Fig. 6a is compatible with edge (u, v), but the connected component \mathcal{W}_4 is incompatible with edge (c, d), which is explained as follows: There are two edges, (v_1, v_a) and (v_2, v_b) , in *G* such that both $v_1 \in Z_1$ and $v_2 \in Z_2$ while Z_i is a vertex cluster in W_4 , i = 1, 2. Let *a* and *b* be the 2-connected components in $\mathcal{G}(\mathcal{C})$ containing endpoints v_a and v_b of the two edges and $T_1(\mathcal{C})$ and $T_2(\mathcal{C})$ be the resulting trees including *c* and *d* after the removal of (c, d) from $\mathcal{T}(\mathcal{C})$. We then have $a \in \mathcal{T}_1(\mathcal{C})$, but $b \in \mathcal{T}_2(\mathcal{C})$ while both $v_1 \in Z_1$ in \mathcal{W}_4 and $v_2 \in Z_2$ in W_4 . Following the definition, W_4 is incompatible with edge (c, d).

Let $\mathcal{A}_1, \mathcal{A}_2, \ldots, \mathcal{A}_l$ be any l clusters in $\hat{\mathcal{G}}$ adjacent to cluster C, l > 0. Then, $E(A_i, C) \neq \emptyset$, for all i, and color each vertex in $E(\mathcal{A}_i, \mathcal{C})$ with the color *blue*, $1 \leq i \leq l$. Denote by $\lambda(\mathcal{A}_1, \mathcal{A}_2, \dots, \mathcal{A}_l)$ a vertex in $\mathcal{T}(\mathcal{C})$ which is the root of a subtree containing all the blue vertices but no other subtree rooted at a proper descendant of $\lambda(A_1, A_2, \dots, A_l)$ also contains all the blue vertices. In other words, $\lambda(\mathcal{A}_1, \mathcal{A}_2, \dots, \mathcal{A}_l)$ is such a vertex in $\mathcal{T}(\mathcal{C})$ that all the edges between A_i and C are incident below it, $1 \leq i \leq l$. Here, "below" means that a vertex is included in a subtree of $\mathcal{T}(\mathcal{C})$ rooted at $\lambda(\mathcal{A}_1, \mathcal{A}_2, \dots, \mathcal{A}_l)$. For example, given two clusters A_1 and A_2 adjacent to C, shown in Fig. 6b, vertex w is $\lambda(A_1, A_2)$, but vertex f is not because the subtree rooted at f does not include the blue vertices band e in $E(\mathcal{A}_2, \mathcal{C})$, where a, b, c, d, and e are blue vertices in $T(\mathcal{C})$.

Recall that C_x and C_y are the clusters including vertices x and y which are adjacent to C, $\mathcal{X}_i \in \mathcal{W}_1$, and $\mathcal{Y}_i \in \mathcal{W}_2$, $1 \leq i \leq p$ and $1 \leq j \leq q$. A vertex v in $\mathcal{T}(\mathcal{C})$ is colored *red* if there is an edge between \mathcal{X}_i and \mathcal{C} incident to $v, 1 \leq i \leq p$. Similarly, a vertex u in $\mathcal{T}(\mathcal{C})$ is colored *black* if there is an edge between \mathcal{Y}_j and \mathcal{C} incident to u, $1 \leq j \leq q$. Define the *top-red* vertex ρ as $\lambda(\mathcal{X}_1, \mathcal{X}_2, \dots, \mathcal{X}_p)$ and the *top-black* vertex β as $\lambda(\mathcal{Y}_1, \mathcal{Y}_2, \dots, \mathcal{Y}_q)$. We now check Type 2 bridges. Let \mathcal{C} be an articulation point in $\mathcal{G}_{x,y}$ separating *x* from *y*. Then, a Type 2 bridge internal to C separating x from y in G may exist. Our objective is to find the bridge if it exists. The intuition behind the algorithm is as follows: Assume that there is an edge e = (u, v) in $\mathcal{T}(\mathcal{C})$ and the corresponding edge of *e* in *G* is a bridge separating *x* from *y*. Let $\mathcal{T}_1(\mathcal{C})$ and $\mathcal{T}_2(\mathcal{C})$ be the two subtrees including u and v after the removal of *e* from $\mathcal{T}(\mathcal{C})$. Following the definition of \mathcal{C}_x and \mathcal{C}_y , every path in G between x and y must have the following structure. It starts with a path outside of C ending at a vertex of C that corresponds to a red vertex in T(C). It ends with a path outside C starting at a vertex of C that corresponds to a black vertex in $\mathcal{T}(\mathcal{C})$. Since the corresponding edge of e internal to C is the bridge of G separating xfrom y, then there is a path in G from x to y containing the corresponding edge of e. Thus, e satisfies the following two



Fig. 6. The notions on $\mathcal{T}(\mathcal{C})$. (a) W_3 is compatible with (u, v) while W_4 is incompatible with (c, d) in $\mathcal{T}(\mathcal{C})$. (b) An illustration of $\lambda(\mathcal{A}_1, \mathcal{A}_2)$.

conditions: 1) *e* must be in a path in $\mathcal{T}(\mathcal{C})$ between a red vertex and a black vertex and 2) each \mathcal{W}_i is *compatible* with *e*, $3 \leq i \leq s$. In other words, for a given \mathcal{W}_i , all the edges between \mathcal{W}_i and \mathcal{C} are incident to either $\mathcal{T}_1(\mathcal{C})$ or $\mathcal{T}_2(\mathcal{C})$, but not to both, $3 \leq i \leq s$. The query algorithm, thus, is to check whether there is such an edge *e* in $\mathcal{T}(\mathcal{C})$ in order to determine whether there is a bridge in *G* separating *x* from *y*. We have the following lemma:

Lemma 8 [19]. Let C be a vertex cluster that is an articulation point in $\mathcal{G}_{x,y}$ separating x from y. There is a bridge inside Cseparating x from y in G if and only if there is an edge e in T(C) that satisfies the following two conditions: 1) The removal of e separates red vertices from black vertices and 2) each connected component W_i is compatible with e, $3 \le i \le s$.

Given ρ and β in T(C), let $v = LCA(\beta, \rho)$ be the lowest common ancestor of ρ and β in T(C). Condition 1 of Lemma 8 can be further decomposed into four cases which are expressed in the following lemma, depending on the relative positions of the three vertices ρ , β , and v in T(C):

Lemma 9 [19]. Let ρ be the top-red, β the top-black vertices, and $v = LCA(\beta, \rho)$. Let $\pi_{\rho,\beta}$ be the path in T(C) between ρ and β . Then, 1) if $v = \rho = \beta$, no edge in T(C) can separate black and red vertices, 2) if $v \neq \rho$ and $v \neq \beta$, then all the edges in $\pi_{\rho,\beta}$ separate black and red vertices, 3) if $v = \beta \neq \rho$, an edge e separates black and red vertices if and only if e is in $\pi_{\rho,\beta}$ and there are no black vertices below child(e), where child(e) is an endpoint of e and another endpoint of e is the parent of *child*(*e*) *in* T(C), *and* 4) *if* $v = \rho \neq \beta$, *an edge e separates black from red vertices if and only if e is in* $\pi_{\rho,\beta}$ *and there are no red vertices below child*(*e*).

Lemmas 8 and 9 imply that we only need to check whether there is an edge e in $\pi_{\rho,\beta}$ of $\mathcal{T}(\mathcal{C})$ separating red vertices from black vertices and every W_i is compatible with $e, 3 \le i \le s$, in order to check whether there is a bridge inside C separating x from y in G. The algorithm proceeds as follows: It first considers all the edges between C and the clusters in \mathcal{W}_1 and \mathcal{W}_2 . If there is no edge e in $\pi_{\rho,\beta}$ separating red vertices from black vertices in $\mathcal{T}(\mathcal{C})$ (Case 1), then there is no bridge inside C separating x from y in G and the query is answered. Otherwise, it examines Cases 2-4 of Lemma 9. Here, we only deal with Case 3. Case 4 is analogous while Case 2 can be dealt by checking the edges in two paths $\pi_{\rho,v}$ and $\pi_{\beta,v}$ in $\mathcal{T}(\mathcal{C})$, respectively, using the very similar fashion. However, the algorithm for Case 3, due to Galil and Italiano, is highly sequential and, therefore, hard to parallelize. A parallel algorithm for Case 3 is proposed below.

Assume that *r* is the root of $\mathcal{T}(\mathcal{C})$ and each vertex *v* in $\mathcal{T}(\mathcal{C})$ has a preorder numbering *preorder*(*v*). Consider the connected components $\mathcal{W}_3, \mathcal{W}_4, \ldots, \mathcal{W}_s$ defined. Let $\gamma_i = \lambda(\mathcal{Z}_{i,1}, \mathcal{Z}_{i,2}, \ldots, \mathcal{Z}_{i,r(i)})$, where cluster $\mathcal{Z}_{i,j} \in \mathcal{W}_i$. If some γ_i is not in path $\pi_{\rho,r}$ of $\mathcal{T}(\mathcal{C})$ between ρ and *r*, then all the edges between $\mathcal{Z}_{i,j}$ and \mathcal{C} are incident to one of the two subtrees, $\mathcal{T}_1(\mathcal{C})$ and $\mathcal{T}_2(\mathcal{C})$, after the removal of any edge *e* in $\pi_{\rho,\beta}$ from $\mathcal{T}(\mathcal{C})$ for all *j*, $1 \leq j \leq r(i)$. This means that \mathcal{W}_i is always compatible with *e* in $\mathcal{T}(\mathcal{C})$ for any edge *e* in $\pi_{\rho,\beta}$ and all the edges between $\mathcal{Z}_{i,j}$ in \mathcal{W}_i and \mathcal{C} do not help



Fig. 7. Checking whether an edge in $\pi_{\rho,\beta}$ satisfies the two conditions of Lemma 8. (a) Condition 2 in the construction of G_1 . (b) An illustration of checking whether an edge in $\pi_{\gamma'_{lo},\beta}$ satisfies the two conditions of Lemma 8.

determine whether *e* separates red vertices from black vertices in $\mathcal{T}(\mathcal{C})$. So, discard all such γ_i s that are not in path $\pi_{\rho,r}$ for further consideration, $3 \le i \le s$. Sort the remaining γ_i s in $\mathcal{T}(\mathcal{C})$ by the preorder numbering in decreasing order and let $\gamma'_0, \gamma'_1, \gamma'_2, \ldots, \gamma'_{s'}$ be the sorted vertex sequence and $\mathcal{W}'_0, \mathcal{W}'_1, \ldots, \mathcal{W}'_{s'}$ be the corresponding connected components, then $preorder(\gamma'_l) > preorder(\gamma'_{l+1}), \ 0 \le l < s',$ $0 \le s' \le s, \ \text{and} \ \gamma'_0 = \rho$. Label each vertex cluster in \mathcal{W}'_i with γ'_i . Thus, each cluster $\mathcal{C}_{i'} \in \mathcal{W}'_i$ has a label $\gamma'_i, \ 0 \le i \le s'$.

The checking of Case 3 is implemented through constructing two auxiliary graphs G_i first, i = 1, 2, and then using the graphs to check whether there is an edge in $\pi_{\rho,\beta}$ in $\mathcal{T}(\mathcal{C})$ satisfying the two conditions in Lemma 8. The objective of constructing G_1 and G_2 is to find a maximal subpath $\pi_{\rho,\gamma'_{i_0}}$ of $\pi_{\rho,\beta}$ including ρ such that no edge in it can satisfy the two conditions in Lemma 8.

The auxiliary graph $G_1(V_1, E_1)$ is constructed as follows: $V_1 = \{\gamma'_0, \gamma'_1, \dots, \gamma'_{s'}\}$. There is an edge $(\gamma'_i, \gamma'_j) \in E_1$ if and only if 1) there is a cluster $C_{i'} \in W'_i$ and a cluster $C_{j'} \in W'_j$ and $E_{i',j'} \neq \emptyset$ or 2) there is a cluster $C_{j'} \in W'_j$ and there is an edge between $C_{j'}$ and C incident below γ'_i in T(C). Fig. 7a illustrates Condition (2) in the construction of G_1 , where urepresents a 2-edge connected component in $\mathcal{G}(C)$. There is an edge between $C_{j'}$ and C incident to u which is below $\gamma'_{i'}$, and $u \in C$. It is obvious that for a given pair of vertices γ'_i and γ'_j in G_1 , if either Condition (1) or Condition 2 holds, then there is an edge in G_1 and a corresponding path $\pi_{\gamma'_i, \gamma'_j}$ in T(C) in which no edge can satisfy the two conditions in Lemma 8 because there are some \mathcal{W}'_l s which are incompatible with e for every edge e in $\pi_{\gamma'_i, \gamma'_i}$.

Let *CC* be a connected component in $G_1(V_1, E_1)$ in which γ'_l and γ'_h are the two vertices with the smallest and the largest indexes l and h. In case *CC* contains only one vertex, then l = h. Clearly, $0 \le l, h \le s'$. We assign every vertex $v \in CC$ a pair of labels, l and h. Let [l, h] represent a closed interval with $l \le h$. Denote by $[l_i, h_i]$ the pair of labels assigned to γ'_i . Then, we have the following lemma:

- **Lemma 10.** Let CC be a connected component in $G_1(V_1, E_1)$ and every vertex in it is labeled with a pair of labels l_i and h_i as above. Then, there are $W'_j s$, which are incompatible with e for every edge e in path $\pi_{\gamma'_{l_i}, \gamma'_{l_i}}$ in $\mathcal{T}(C)$, $1 \le j \le s'$.
- **Proof.** Let CC be a connected component in G_1 in which every vertex has been labelled with l_i and h_i . Then, the vertices γ'_{l_i} and γ'_{h_i} are included in *CC* by the definition. Following the above discussion, we know that there is no edge in $\pi_{\gamma'_{c},\gamma'_{t}}$ in $\mathcal{T}(\mathcal{C})$ satisfying the two conditions in Lemma 8 for every edge $(\gamma'_a, \gamma'_b) \in E_1$ in G_1 by the construction of G_1 and $l_i \leq a, b \leq h_i$. Thus, every edge in G_1 corresponds to a subpath of $\pi_{\rho,r}$ in $\mathcal{T}(\mathcal{C})$ in which no edge satisfies the two conditions in Lemma 8. Now, consider two edges (γ'_a, γ'_b) and (γ'_b, γ'_c) sharing a common vertex γ'_b in a *CC*, then no edge in $\pi_{\gamma'_a,\gamma'_c}$ satisfies the two conditions in Lemma 8 because both $\pi_{\gamma'_a,\gamma'_b}$ and $\pi_{\gamma'_b,\gamma'_c}$ do not contain any edge satisfying the two conditions. Therefore, no edge in path $\pi_{\gamma'_{l_{\nu}},\gamma'_{h_{\nu}}}$ in $\mathcal{T}(\mathcal{C})$ satisfies the two conditions in Lemma 8 because $\pi_{\gamma'_{L},\gamma'_{h}}$ is the union of the subpaths generated by the corresponding edges of G_1 in the CC.

Given that each vertex $\gamma'_i \in V_1$ has a pair of labels l_i and h_i , $0 \leq i < s'$, another graph $G_2(V_2, E_2)$ is constructed as follows: $V_2 = V_1$ and $(\gamma'_i, \gamma'_j) \in E_2$ if and only if 1) either $(\gamma'_i, \gamma'_j) \in E_1$ or 2) the intersection of the two intervals is nonempty, i.e., $[l_i, h_i] \cap [l_j, h_j] \neq \emptyset$, which means that two paths $\pi_{\gamma'_l, \gamma'_{h_i}}$ and $\pi_{\gamma'_l, \gamma'_{h_i}}$ are overlapping. Since no edge in both $\pi_{\gamma'_l, \gamma'_{h_i}}$ and $\pi_{\gamma'_l, \gamma'_{h_i}}$ satisfies the two conditions of Lemma 8, no edge in the union of the two paths satisfies the two conditions in Lemma 8. We then have the following lemma:

Lemma 11. Let CC be a connected component in $G_2(V_2, E_2)$, including vertex $\rho = \gamma'_0$ such that i_0 is the maximum index of the vertices in it. Then, there is no edge e in $\pi_{\rho,\gamma'_{i_0}}$ in $\mathcal{T}(\mathcal{C})$ satisfying 1) the removal of e separates red vertices from black



Fig. 8. An illustration of the proof of Lemma 11.

vertices and 2) each connected component W'_i is compatible with $e, 1 \leq i \leq s'$.

Proof. Let CC' be a connected component in G_1 consisting of vertices $\gamma'_{i_1}, \gamma'_{i_2}, \ldots, \gamma'_{i_k}$. Assume that these vertices are sorted by their preorder numberings in $\mathcal{T}(\mathcal{C})$ in decreasing order, i.e., $preorder(\gamma'_{i_j}) > preorder(\gamma'_{i_{j+1}})$ and $1 \le j \le k - 1$. Then, there is no edge in path $\pi_{\gamma'_{i_1},\gamma'_{i_k}}$ in $\mathcal{T}(\mathcal{C})$ satisfying the two conditions in Lemma 8, which is derived by Lemma 10 directly.

Let CC_i and CC_j be the two connected components in G_1 and $[l_i, h_i]$ and $[l_j, h_j]$ the pairs of labels for them. If $[l_i, h_i] \cap [l_j, h_j] \neq \emptyset$, we can distinguish four cases:

- 1. $l_i > l_j$ but $h_i < h_j$, i.e., $[l_i, h_i] \subset [l_j, h_j]$. Path $\pi_{\gamma'_{l_i}, \gamma'_{h_i}}$ is a subpath in $\mathcal{T}(\mathcal{C})$ of $\pi_{\gamma'_{l_j}, \gamma'_{h_j}}$ while no edge in $\pi_{\gamma'_{l_j}, \gamma'_{h_j}}$ in $\mathcal{T}(\mathcal{C})$ satisfies the two conditions in Lemma 8.
- 2. $l_i < l_j$ and $h_i < h_j$. There is no edge in $\pi_{\gamma'_{l_i},\gamma'_{h_j}}$ in $\mathcal{T}(\mathcal{C})$ satisfying the two conditions in Lemma 8 because both $\pi_{\gamma'_{l_i},\gamma'_{h_i}}$ and $\pi_{\gamma'_{l_j},\gamma'_{h_j}}$ are the subpaths of $\pi_{\gamma'_{l_i},\gamma'_{h_i}}$.
- 3. $l_i < l_j$ but $h_i > h_j$. This case is similar to Case 1, omitted.
- 4. $l_i > l_j$ and $h_i > h_j$. This case is similar to Case 2 and omitted.

Thus, each edge (γ'_a, γ'_b) in a connected component, including ρ , corresponds to a path $\pi_{\gamma'_a,\gamma'_b}$ in $\mathcal{T}(\mathcal{C})$. So, we conclude that there is no edge in $\pi_{\rho,\gamma'_{i_0}}$ in $\mathcal{T}(\mathcal{C})$ satisfying the two conditions in Lemma 4 because $\pi_{\rho,\gamma'_{i_0}}$ is the union of the subpaths derived from the corresponding edges in the connected component including ρ in G_2 .

The example, shown in Fig. 8, illustrates the idea used in the proof of Lemma 11. Assume that $\gamma'_{j_1}, \gamma'_{j_2}, \gamma'_{j_3}$, and γ'_{j_4} are the vertices in a connected component in G_1 , $\gamma'_{i_1}, \gamma'_{i_2}, \gamma'_{i_3}$, and γ'_{i_4} are the vertices in another connected component in G_1 . We have shown that there is no edge in either $\pi_{\gamma'_{j_1},\gamma'_{j_4}}$ or $\pi_{\gamma'_{i_1},\gamma'_{i_4}}$ of $\mathcal{T}(\mathcal{C})$ satisfying the two conditions of Lemma 8, by Lemma 10. Suppose that $[j_1, j_4] \cap [i_1, i_4] \neq \emptyset$, then the vertices $\gamma'_{j_1}, \gamma'_{j_2}, \gamma'_{j_3}, \gamma'_{j_4}, \gamma'_{i_1}, \gamma'_{i_2}, \gamma'_{i_3}$, and γ'_{i_4} are in the same connected component in G_2 , by the definition of G_2 . Clearly, there is no edge in path $\pi_{\gamma'_{j_1},\gamma'_{i_4}}$ in $\mathcal{T}(\mathcal{C})$ satisfying the two conditions in Lemma 8 due to that $\pi_{\gamma'_{j_1},\gamma'_{i_4}}$ is the union of $\pi_{\gamma'_{j_1},\gamma'_{j_4}}$ and $\pi_{\gamma'_{i_1},\gamma'_{i_4}}$.

Having built G_1 and G_2 , to decide whether there is an edge in $\pi_{\rho,\beta}$ in $\mathcal{T}(\mathcal{C})$ satisfying the two conditions in Lemma 8 (Case 3), it proceeds as follows: If $preorder(\beta) > preorder(\gamma'_{i_0})$, then there is no edge in $\pi_{\rho,\beta}$ satisfying the two conditions in Lemma 8 because $\pi_{\rho,\beta}$ is a subpath of $\pi_{\rho,\gamma'_{i_0}}$ and there is no edge in $\pi_{\rho,\gamma'_{i_0}}$ satisfying the two conditions in Lemma 8 by Lemma 11, where vertices γ'_{i_0} and ρ are in the same connected component in G_2 and γ'_{i_0} has the maximum index i_0 in the connected component. Otherwise, although there is no edge in path $\pi_{\rho,\gamma'_{io}}$ satisfying the two conditions of Lemma 8, we cannot guarantee that this claim also holds in path $\pi_{\gamma'_{i_0},\beta}$. Path $\pi_{\rho,\beta}$ is the union of $\pi_{\rho,\gamma'_{i_0}}$ and $\pi_{\gamma'_{i_0},\beta}$. To check whether there is an edge in path $\pi_{\gamma'_{i_{\alpha}},\beta}$ satisfying the two conditions of Lemma 8, we proceed by checking whether there is a vertex cluster $\mathcal{Y}_j \in \mathcal{W}_2$ including C_y such that $u \in E(\mathcal{Y}_i, \mathcal{C})$ and

$$preorder(\gamma'_{i_0}) \le preorder(u) < preorder(\gamma'_{i_0}) + nd(\gamma'_{i_0}) - 1,$$
(1)

where nd(z) is the number of vertices in a subtree of $\mathcal{T}(\mathcal{C})$ rooted at z. In other words, we need to check whether there is an edge between \mathcal{Y}_j and \mathcal{C} incident to u and u is below γ'_{i_0} in $\mathcal{T}(\mathcal{C})$, $1 \leq j \leq q$. Fig. 7b illustrates this case. If (1) holds, then no edge in $\pi_{\gamma'_{i_0},\beta}$ satisfies the two conditions of Lemma 8, which means there is no bridge internal to \mathcal{C} separating x from y in G. We then examine the other articulation points \mathcal{C}' , if existing in $\mathcal{G}_{x,y}$, separating x from yto determine whether x and y are in the same 2-edge connected component in G. If none of the articulation points contains a bridge separating x from y in G, then x and y are in the same 2-edge connected component in G.

3.4.2 Implementing the Operations

Now, we discuss the implementation details of the proposed algorithm for query processing. Due to space limitations, some unimportant details are omitted. Assume that the data structures defined are available.

 $\mathcal{G}_{x,y}$ is constructed as follows: First, a graph \mathcal{G}' is generated from $\overline{\mathcal{G}}$ by removing the vertices \mathcal{C}_x and \mathcal{C}_y and their adjacent artificial vertices (i.e., the new vertices added in the generation of $\overline{\mathcal{G}}$). Then, $\mathcal{G}_{x,y}$ is generated by merging the adjacency lists of \mathcal{G}' , $\mathcal{F}(\mathcal{C}_x)$, and $\mathcal{F}(\mathcal{C}_y)$, which takes $O(\log K + \log(m/K))$ time using $O(K + (m/K)^2)$ processors on a CRCW PRAM because $\mathcal{G}_{x,y}$ contains $O(K + (m/K)^2)$ vertices and edges.

Given the graph $\mathcal{G}_{x,y}$, the next objective is to find all articulation points \mathcal{C} s in $\mathcal{G}_{x,y}$ separating x from y. Obviously, each such \mathcal{C} is also an articulation point in both $\overline{\mathcal{G}}$ and $\hat{\mathcal{G}}$, separating \mathcal{C}_x from \mathcal{C}_y , which can be found by the following lemma:

- **Lemma 12.** Finding all articulation points in \mathcal{G} separating C_x from C_y takes $O(\log(m/K) + \log K)$ time using $O((m/K)^3)$ processors on a CRCW PRAM.
- **Proof.** Since $\hat{\mathcal{G}}$ and $\overline{\mathcal{G}}$ are the equivalent graphs, an articulation point in $\hat{\mathcal{G}}$ is also an articulation point in $\overline{\mathcal{G}}$. We proceed with the following in parallel: For each vertex C in $\hat{\mathcal{G}}$, make a copy $\mathcal{G}_{\mathcal{C}}$ of $\overline{\mathcal{G}}$ for \mathcal{C} , delete C and its incident edges from $\mathcal{G}_{\mathcal{C}}$, and compute the connected components of the resulting graph $\mathcal{G}_{\mathcal{C}} - \{\mathcal{C}\}$. If vertices \mathcal{C}_x and C_y are not in the same connected component in $\mathcal{G}_{\mathcal{C}} - \{\mathcal{C}\}$, then \mathcal{C} is an articulation point in \mathcal{G} separating \mathcal{C}_x from C_y . Since the connected component computation in graph $\mathcal{G}_{\mathcal{C}} - \{\mathcal{C}\}$ for each \mathcal{C} requires $O(\log(m/K))$ time and $O((m/K)^2)$ processors on a CRCW PRAM [43], finding all articulation points in $\overline{\mathcal{G}}$ separating \mathcal{C}_x from \mathcal{C}_y requires $O(\log(m/K) + \log K)$ time and $O((m/K)^3)$ processors on a CRCW PRAM because $\overline{\mathcal{G}}$ contains O(m/K) vertex clusters (all the other vertices in $\overline{\mathcal{G}}$ are artificial vertices) and $O((m/K)^2)$ edges. П

Now, assume that every vertex v in $\mathcal{T}(\mathcal{C})$ has been assigned a preorder numbering preorder(v). Otherwise, such an assignment can be done in $O(\log K)$ time with O(K) processors using the Euler traversal technique [31] because $\mathcal{T}(\mathcal{C})$ is an inverted tree containing O(K) vertices and edges. Also, answering a LCA query in $\mathcal{T}(\mathcal{C})$ can be answered in O(1) time using one processor on an EREW PRAM [42].

- **Lemma 13.** Let $A_1, A_2, ..., A_l$ be l vertex clusters adjacent to a vertex cluster C. $\lambda(A_1, A_2, ..., A_l)$ can be found in $O(\log K + \log l)$ time with O(l) processors on a CRCW *PRAM*, where $1 \le l \le \lceil cm/K \rceil$ and c is constant.
- **Proof.** For every A_i , the two vertices in $E(A_i, C)$ with the smallest and the largest preorder numberings in T(C) can be found in $O(\log K)$ time using one processor due to the fact that $E(A_i, C)$ contains O(K) vertices and is maintained as a balanced binary search tree. Let u_i and v_i be the two vertices found from $E(A_i, C)$. Then, sort the sequence of 2l vertices in T(C) by their preorder numberings in increasing order, which can be done in $O(\log l)$ time with O(l) processors. $\lambda(A_1, A_2, \ldots, A_l)$ is the LCA of the two vertices with the smallest and the largest preorder numberings in the sorted sequence. So, the entire computation takes $O(\log l + \log K)$ time and uses O(l) processors on a CRCW PRAM.

- **Corollary 1.** ρ and β can be found in $O(\log(m/K) + \log K)$ time using O(m/K) processors on a CRCW PRAM.
- **Proof.** It can be derived from Lemma 13 with $l \le cm/K$ because $1 \le p, q \le cm/K$, *c* is constant.

Recall that $\gamma_i = \lambda(\mathcal{Z}_{i,1}, \mathcal{Z}_{i,2}, \dots, \mathcal{Z}_{i,r(i)})$, where $\mathcal{Z}_{i,j} \in \mathcal{W}_i$, $1 \leq j \leq r(i)$, and $3 \leq i \leq s$. Finding the sequence $\gamma_3, \gamma_4, \dots, \gamma_s$ takes $O(\log K + \log(m/K))$ time using

$$O\!\left(\sum_{i=3}^s r(i)\right) = O(m/K)$$

processors on a CRCW PRAM. Let $\gamma_1 = \rho$ and $\gamma_2 = \beta$, which can be found in the same time and processor bounds. Then, discard those γ_i s that are not in $\pi_{\rho,r}$. Let $\gamma'_1, \gamma'_2, \ldots, \gamma'_{s'}$ be the sorted resulting sequence by the preorder numbering of vertices in $\mathcal{T}(\mathcal{C})$ in decreasing order. This sequence can be obtained in

$$O(\log s) = O(\log(m/K))$$

time using O(s) = O(m/K) processors by checking if $\gamma_i = LCA(\rho, \gamma_i)$ which takes O(1) time using one processor, $1 \le i \le s \le cm/K$. Furthermore, let $\mathcal{W}'_0, \mathcal{W}'_1, \ldots, \mathcal{W}'_{s'}$ be the corresponding connected components of $\gamma'_0, \gamma'_1, \gamma'_2, \ldots, \gamma'_{s'}$, where $0 \le s' \le s$ and $\gamma'_0 = \rho$.

The complexity for the construction of G_1 is as follows: Checking Condition 1 takes O(1) time using $O((m/K)^2)$ processors because there are O(m/K) vertex clusters and checking Condition 2 takes $O(\log K)$ time using O(m/K) processors, which has been proven by the following lemma:

Lemma 14 [19]. Let C' be any cluster adjacent to C and let v be any vertex of T(C). Then, to check whether there are edges in G between C' and C incident below v in $O(\log K)$ time with one processor on a CRCW PRAM.

Notice that we use the same data structure as Galil and Italiano [21] did in this part. The proof of Lemma 14 is omitted.

The construction of G_1 thus requires $O(\log K)$ time and $O((m/K)^2)$ processors because $1 \le s' \le cm/K$. Finding connected components in G_1 requires $O(\log(m/K))$ time and $O((m/K)^2)$ processors on a CRCW PRAM because G_1 contains O(m/K) vertices and $O((m/K)^2)$ edges [43]. Broadcasting the smallest and the largest indices of the vertices in a connected component to all the vertices in it requires $O(\log(m/K))$ time and O(m/K) processors on an EREW PRAM. Similarly, the construction of G_2 and finding connected components in G_2 takes $O(\log(m/K))$ time using $O((m/K)^2)$ processors on a CRCW PRAM. As a result, the maximum index i_0 of a vertex in the connected component in G_2 containing ρ can be found in the above time and processor bounds. Thus, the path $\pi_{\rho,\gamma'_{\ell}}$ in $\mathcal{T}(\mathcal{C})$ can be found, which does not contain any edge satisfying the two conditions in Lemma 8. Checking whether there is any edge in $\pi_{\gamma'_{i,\beta}\beta}$ satisfying the two conditions of Lemma 8 can be done in $O(\log K)$ time using O(m/K) processors by Lemma 14, because there are O(m/K) vertices in $\mathcal{T}(\mathcal{C})$. Thus, we have the following theorem:

- **Theorem 3.** The query about whether x and y are in the same 2-edge connected component in a connected graph G(V, E) can be answered in $O(\log m)$ time using $O(m^{3/4})$ processors on a CRCW PRAM.
- **Proof.** The construction of $\mathcal{G}_{x,y}$ takes $O(\log K + \log(m/K))$ time using $O(K + (m/K)^2)$ processors. Following Theorem 2, Type 1 bridge checking takes

$$O(\log K + \log(m/K))$$

time using $O(K + (m/K)^2)$ processors. Type 2 bridge checking needs finding all articulation points in $\mathcal{G}_{x,y}$ separating x from y, which requires $O(\log K + \log(m/K))$ time using $O((m/K)^3)$ processors. Checking all the cases in Lemma 9 requires $O(\log(m/K) + \log K)$ time using $O((m/K)^2)$ processors. The theorem then follows, by setting $K = m^{3/4}$.

3.5 Inserting and Deleting Edges

In this section, we consider the InsertEdge(x, y) operation. Let C_x and C_y be the vertex clusters including x and y. When a new edge (x, y) is inserted, the degrees of vertices x and yin the original graph G increase by one. If none of them becomes four, only the data structures pertinent to C_x and C_y are required to be updated, i.e., $\mathcal{F}(\mathcal{C}_x)$, $\mathcal{T}(\mathcal{C}_x)$, $\mathcal{F}(\mathcal{C}_y)$, $\mathcal{T}(\mathcal{C}_y)$, $E(\mathcal{C}_x, \mathcal{C}_y), E(\mathcal{C}_y, \mathcal{C}_x)$, and \mathcal{G} needed be updated. Otherwise, if either the degree of x or the degree of y becomes four, then the transformation given in Section 2 will be applied. This leads to a constant number of extra vertices and edges in the resulting graph. Assume that the degree of x in G now is four. We apply the transformation to the cluster C_x . As a result, all data structures pertinent to C_x must be updated to reflect the change. However, to keep the vertex degree of a graph no greater than three, the insertion of extra vertices may lead to the size of C_x becoming greater than 3K/2. We deal with this latter case by splitting C_x into two vertex clusters C' and C''. Since the original size of C_x is between K/2 and 3K/2 and the degree of each vertex in it is no greater than three, the two new clusters C' and C'' as well as the data structures pertinent to them can be constructed in $O(\log K)$ time using $O(K + (m/K)^2)$ processors on a CRCW PRAM, which is explained as follows: The full representations $\mathcal{F}(\mathcal{C}')$ and $\mathcal{F}(\mathcal{C}'')$ and the tree representations $\mathcal{T}(\mathcal{C}')$ and $\mathcal{T}(\mathcal{C}'')$ of \mathcal{C}' and \mathcal{C}'' can be obtained in $O(\log K)$ time with O(K) processors by the discussion in Section 3.3. The balanced binary search tree for the other cluster C_i adjacent to C_x , $E(C', C_j)$, $E(C_j, C')$, $E(C'', C_j)$, and $E(C_j, C'')$ can be constructed in $O(\log K)$ time using O(K) processors because the degree of every vertex in C_x is at most three. The handling of C_y can be done similarly. The super graph $\hat{\mathcal{G}}$ and its equivalent graph $\overline{\mathcal{G}}$ can be updated in $O(\log(m/K) +$ $\log K$) time using $O((m/K)^2)$ processors because they contain $O((m/K)^2)$ vertices.

To perform an insertion of (x, y), we distinguish two cases depending on whether (x, y) is an edge internal to a cluster or an intercluster edge. 1) If $C_x = C_y = C$, then (x, y)is an internal edge. As a result, only the data structures related to C, $\mathcal{F}(C)$ and $\mathcal{T}(C)$, are updated, which takes $O(\log K)$ time using O(K) processors on a CRCW PRAM. 2) If $C_x \neq C_y$, then (x, y) is an intercluster edge. The data structures needed to be updated are $\mathcal{F}(C_x)$, $\mathcal{T}(C_x)$, $\mathcal{F}(\mathcal{C}_y)$, $\mathcal{T}(\mathcal{C}_y)$, $E(\mathcal{C}_x, \mathcal{C}_y)$, $E(\mathcal{C}_y, \mathcal{C}_x)$, and $\hat{\mathcal{G}}$. This takes $O(\log K + \log(m/K))$ time using $O((m/K)^2)$ processors on a CRCW PRAM. In summary, inserting an edge (x, y) causes at most two clusters to be split and at most constant clusters' data structures to be updated. Following the above discussion, each update requires $O(\log K + \log(m/K))$ time using $O(K + (m/K)^2)$ processors on a CRCW PRAM.

To keep $O(\log m)$ time complexity per update and per query in the worst case, we need to be cautious when dealing with the data structure maintenance. Let m_t be the number of edges in the graph at time t. We show an update at time t can be implemented in $O(\log m_t)$ time with $O(m_t^{3/4})$ processors on a CRCW PRAM. Let $K_t = \lceil m_t^{3/4} \rceil$. When the value of K changes due to an InsertEdge or DeleteEdge operation, there will be at least $[m_t^{3/4}/2]$ more updates before K becomes twice as large or one half as small as it was before. Following the idea of Galil and Italiano [19] that adjusts a constant number of clusters each time there is an update, this gives a total of $O(m_t^{3/4})$ cluster adjustments. Since there are no more than $O(m_t/K_t) =$ $O(m_t^{1/3})$ clusters that need to be adjusted and the adjustments can be accomplished before a new round of adjustment starts. Thus, whenever there is an insertion, the clusters can be scanned to find any cluster that is too small and a constant number of these clusters can be merged with a neighbor if needed.

The processing for the deletion case is similar to that for the insertion case, which is omitted here. Therefore, we have the following theorem:

- **Theorem 4.** Full dynamic maintenance of 2-edge connected components in a connected graph G(V, E) with m edges can be done in $O(\log m)$ time using $O(m^{3/4})$ processors on a CRCW PRAM.
- **Proof.** From the discussion above and Theorem 3, the theorem follows.

4 MAINTAINING 2-EDGE AND 3-EDGE CONNECTED COMPONENTS IN GENERAL GRAPHS

4.1 Maintaining 2-Edge Connected Components in Disconnected Graphs

We deal with the fully dynamic maintenance of 2-edge connected components in disconnected graphs by extending the results of connected graphs. Let G be a disconnected graph consisting of l connected components. We augment G by adding l-1 dummy edges such that the augmented graph is connected and no cycle in the augmented graph includes two dummy edges. Every edge in G is assigned weight 1 and every dummy edge is assigned weight 2. Such augmentation does not change the property of 2-edge connected components in G, i.e., two vertices x and y are in a 2-edge connected component in the augmented graph if and only if they are in the same 2-edge connected component in G. The key here is to maintain a vertex cluster partition of a minimum spanning tree T instead of maintaining a vertex cluster partition of an arbitrary spanning tree in the augmented graph.

When inserting an edge e = (x, y), we check whether there is already a dummy edge in *T* between them. If yes, we decrease the weight of *e* from 2 to 1; otherwise, we insert this edge with weight 1 to *T* and update *T*, which can be implemented in $O(\log m)$ time with $O(m^{2/3})$ processors on a CREW PRAM, using either Ferragina's algorithm [14] or Liang and McKay's algorithm [35]. If *e* becomes an edge in the new minimum spanning tree through deleting a dummy edge *e'*, then *e'* is deleted. When deleting an edge e = (x, y), we check whether *e* is in *T*. If yes, *T* is updated; otherwise, *e* is deleted. Therefore, we have the following theorem:

Theorem 5. Full dynamic maintenance of 2-edge-connected components in a disconnected graph G(V, E) can be done in $O(\log m)$ time using $O(m^{3/4})$ processors on a CRCW PRAM.

Proof. It is straightforward from Theorem 4.

The above proposed algorithm can be further improved if the sparsification technique of Eppstein et al. [10] is employed. In the following, we present an improved algorithm for the problem, using the sparsification technique.

If *G* is *k*-edge connected, we can use a sparse *k*-edge certificate $U_k \ (= \bigcup_{i=1}^k T_i)$ of *G* instead of *G* itself when dealing with the *k*-edge connectivity of *G*. Eppstein et al. [10] have shown that U_k is a strong *k*-edge certificate of *G*. Obviously, the sparse 2-edge certificate of *G* is a special case with k = 2. For any fixed *k*, U_k can be maintained in time $O(\log n)$ with $O(n^{2/3})$ processors per update, using the *k* copies of the algorithm for fully dynamic maintenance of maintain forest T_i , $i = 1, \ldots, k$. Also, U_k is stable since at most a single edge flips in one of the *k* minimum spanning forests when there is an update.

- **Theorem 6.** Full dynamic maintenance of 2-edge connected components in a graph G(V, E) can be done in $O(\log n \log \frac{m}{n})$ time using $O(n^{3/4})$ processors on a CRCW PRAM.
- Proof. By applying the sparsification technique, we first build a tree \mathcal{BT} and maintain the sparse 2-edge certificate at every node in \mathcal{BT} by two dynamic data structures, one for 1-edge connectivity which requires $O(\log n)$ time and $O(n^{2/3})$ processors on a CREW PRAM [14], [35] and another for 2-edge connectivity $U_2 = T_1 \cup T_2$, i.e., T_1 and T_2 . We then run the parallel algorithm for the maintenance of 2-edge connected components in U_2 , described in the beginning of this section, at the root node of \mathcal{BT} . So, for every edge insertion and deletion, the maintenance time spent for all the data structures involved is $O(\log n \log \frac{m}{n})$ and the number of processors used is $O(n^{3/4})$ because the height of \mathcal{BT} is $O(\log \frac{m}{n})$. Similarly, the query can be answered at the root of \mathcal{BT} in the same time and processor bounds as above.

4.2 Maintaining 3-Edge Connected Components

Theorem 7. Fully dynamic maintenance of 3-edge connected components in a graph G(V, E) requires

$$O(\log n \log \frac{m}{n} + \log n \log \log n / \alpha(3n, n))$$

time using $O(n\alpha(3n, n)/\log n)$ processors per update and O(1) time with a single processor per query on a CRCW PRAM. **Proof.** Let U_3 be a sparse 3-edge certificate of G. By applying the reduction technique of Galil et al. that reduces the *k*-edge connectivity of G into the *k*-vertex connectivity of another graph $\varphi_k(G)$, we first construct $\varphi_3(U_3)$ from U_3 , which takes $O(\log n)$ time using $O(n/\log n)$ processors on a CRCW PRAM because $\varphi_3(U_3)$ contains O(n) vertices and edges. We then maintain a tree \mathcal{BT} in which each node contains a sparse 3-vertex certificate of the subgraph induced by the edges at the leaf nodes in the subtree rooted at the node, using Eppstein et al.'s sparsification technique. The computational complexity of the fully dynamic maintenance of 3-vertex connected components is given by Theorem 9 in Section 5.3, the theorem then follows.

5 MAINTAINING BICONNECTED AND TRICONNECTED COMPONENTS IN GENERAL GRAPHS

5.1 Finding a Strong and Sparse *k*-Vertex Certificate of *G*

Eppstein et al.'s sparsification technique [10] is applicable only if the certificate \mathcal{P} is both sparse and strong. For *k*-vertex connectivity of a graph *G*, Eppstein et al. [10] have shown that $C_k (= \cup_{j=1}^k B_j)$ is a strong, sparse certificate of *G*, where B_j is a breadth-first search forest in graph $G - \bigcup_{i=1}^{j-1} B_i$. However, currently, the number of processors used in any NC algorithm for finding a breadth-first search forest is no better than the number of scalar operations for matrix multiplication, which is $O(n^{2.376})$ [7]. To reduce the number of processors used, we adopt an alternative approach to find a sparse *k*-vertex certificate and show the found certificate is a strong certificate.

To find a sparse k-vertex certificate of G(V, E), we use the *scan-first search* technique [4]. Let $C_k = G(V, \bigcup_{i=1}^k E_i)$, where E_i is the edge set of the spanning forest F_i obtained by applying the scan-first search in graph $G_i = G - \bigcup_{j=1}^{i-1} F_j$. It is obvious that C_k is sparse when k is fixed. In the following, we show that C_k is strong too:

Lemma 15. C_k is a strong k-vertex certificate of G(V, E).

Proof. Let *S* be a (k-1)-vertex cut in any graph $C_k \cup H$ which partitions the vertices in V-S into two subsets *A* and *B*. Obviously, *H* does not contain any edge between a vertex in *A* and a vertex in *B*. Let $s_i \in S - \{s_1, s_2, \ldots, s_{i-1}\}$ be the first vertex in the *i*th scan-first search of $G_i = G - F_1 - F_2 - \ldots - F_{i-1}$, where F_i is a scan-first search spanning forest of G_i .

Let *T* be a tree rooted at *r* containing s_1 and $T \in F_1$. We claim either $r \in A$, $r \in B$, or $r \in S$. If $r \in S$, then $r = s_1$ by the following argument: The index of *r* is smaller than the index of s_1 because *r* is visited before s_1 by the definition of scan-first search. This contradicts that s_1 has the smallest index in *S*. So, either $r \in A$ or $r \in B$. Without loss of generality, assume that $r \in A$. In the following, we show that all the edges between s_1 and the vertices in *B* are included in *T*, therefore, in F_1 because $T \in F_1$. We proceed as follows:

If $r = s_1$, then all such edges incident to s_1 are in T obviously. Now, consider $r \in A$. Assume that there is an edge $(s_1, v) \in E$ and $v \in B$, but $(s_1, v) \notin T$. We show this is impossible. Since s_1 and v are in the same connected

component in G, they must be in the same spanning tree in *G*. Therefore, both s_1 and v are included in *T*. Let p(v)be the parent of v in T. We then have either $p(v) \in A$, $p(v) \in \hat{B}$, or $p(v) \in S$ and only one of these holds. If $p(v) \in A$, then there is an edge between $p(v) \in A$ and v $(\in B)$ after the deletion of all the vertices in S and the edges incident to them. This contradicts that S is a (k-1)-vertex cut between A and B. Thus, either $p(v) \in B$ or $p(v) \in S$. Obviously, it is also impossible that $p(v) \in B$ because the index of s_1 is smaller than that of any other vertex in *B* by the definition of scan-first search. Thus, s_1 must be the parent of v in T because $(s_1, v) \in E$. Therefore, p(v) must be in S and $p(v) = s_1$, i.e., the edge (s_1, v) must be included in T. In other words, all the edges between s_1 and the vertices in B are included in F_1 instead of in $C_k - F_1$ and $S - \{s_1\}$ forms a (k-2)-vertex cut in $C_k - F_1$. We proceed in the same way on the resulting graph. As a result, each application of scan-first search will eliminate a vertex from G and in the end F_k must be disconnected. But this can only happen if *S* is also a vertex cut in *G* and, hence, in $G \cup H$. We have already shown that any vertex cut in $C_k \cup H$ is also a vertex cut in $G \cup H$. The converse follows immediately from the fact that C_k is a subgraph of G. Hence, C_k is a strong certificate of G. П

5.2 Maintaining Biconnected Components

Assume that *G* is connected. Otherwise, two types of data structures are maintained, one for 1-edge connectivity (connected components) and another for biconnectivity (2-vertex connected components). As a result, if there is a query to ask whether vertices *u* and *v* are in the same biconnected component in *G*, we first check the data structures for 1-edge connectivity to see whether *u* and *v* are connected. If yes, we then run the query algorithm on the data structures for biconnectivity. Otherwise, the response to the query is *false*. It is known that the fully dynamic maintenance of 1-edge connectivity can be done in $O(\log n \log \frac{m}{n})$ time using $O(n^{2/3})$ processors on a CREW PRAM [8], [35].

- **Theorem 8.** Fully dynamic maintenance of biconnected components of a graph G(V, E) requires $O(\log^2 n)$ time using $O(n\alpha(2n, n)/\log n)$ processors per update and O(1) time with one processor per query or requires $O(\log n \log \frac{m}{n})$ time using $O(n\alpha(2n, n)/\log n)$ processors per update and $O(\log n)$ time using $O(n\alpha(2n, n)/\log n)$ processors per query on a CRCW PRAM.
- **Proof.** When an edge is inserted or deleted, it results in only O(1) leaf node updates in \mathcal{BT} . Meanwhile, the sparse 2-vertex certificates of the involved leaf nodes and their ancestor nodes in the paths in \mathcal{BT} from the leaf nodes to the root node need to be updated. Finding the sparse 2-vertex certificate at a node requires $O(\log n)$ time and $O(n\alpha(2n,n)/\log n)$ processors by Cheriyan et al.'s algorithm [4]. At the root node, we run Tarjan and Vishkin's algorithm [44] to maintain biconnected components in the sparse 2-vertex certificate at the root node while finding biconnected components in a graph with *n* vertices and *m* edges requires $O(\log n)$ time and O(m + n) processors on a CRCW PRAM, which can also be implemented in $O(\log^2 n)$ with

 $O((m+n)/\log n)$ processors by Brent's theorem [3]. Therefore, updating all data structures requires

$$O(\log n \log \frac{m}{n} + \log^2 n) = O(\log^2 n)$$

time and $O(n\alpha(2n, n)/\log n)$ processors when there is an insertion or a deletion. For a query, we just check the root node of \mathcal{BT} to see whether the two vertices are in the same biconnected component, which takes O(1) time using one processor. If we want to reduce the update time further, there is another approach in which, for each update, we recompute the sparse 2-vertex certificates for all involved nodes in \mathcal{BT} and for each query about whether there exist k vertex disjoint paths, we run the algorithm by Khuller and Schieber [32] at the root of \mathcal{BT} . For the case where k = 2, their algorithm requires $O(\log n)$ time and $O(n\alpha(2n,n)/\log n)$ processors. So, it requires $O(\log n \log \frac{m}{n})$ time using $O(n\alpha(2n, n)/\log n)$ processors per update and $O(\log n)$ time using $O(n\alpha(2n,n)/\log n)$ processors per query on a CRCW PRAM.

Notice that there is a faster algorithm for testing k vertex disjoint paths between a pair of vertices [29], which requires $O(k \log \log k \log n)$ time and $O((n^2 + k)k(C(n, m) + kn))$ processors on an arbitrary CRCW PRAM. This algorithm is superior to the algorithm of Khuller and Schieber [32] only when k is not constant.

5.3 Maintaining Triconnected Components

Assume that *G* is biconnected. Otherwise, two types of data structures are maintained, one for biconnectivity and another for triconnectivity of *G*. That is, if there is a query to ask whether vertices *u* and *v* are in the same triconnected component, we first check the data structures for biconnectivity to see whether *u* and *v* are biconnected. If yes, we then run the query procedure on the data structures for triconnectivity. Otherwise, the response to the query is *false*. By Theorem 8, the fully dynamic maintenance of biconnectivity can be done in $O(\log n \log \frac{m}{n})$ time using $O(n\alpha(2n,n)/\log n)$ processors on a CRCW PRAM.

Theorem 9. Fully dynamic maintenance of triconnected components in a graph G(V, E) requires

$$O(\log n \log \frac{m}{n} + \log n \log \log n / \alpha(3n, n))$$

time using $O(n\alpha(3n, n)/\log n)$ processors per update and O(1) time with one processor per query. The algorithm runs in a CRCW PRAM.

Proof. The approach employed is the same as the one used for Theorem 8, except that we replace Tarjan and Vishkin's [44] algorithm for biconnectivity by the algorithm for triconnectivity due to Fussel et al. [17]. Their algorithm for the fully dynamic maintenance of triconnectivity of a graph with *n* vertices and *m* edges requires $O(\log n)$ time and $O((m + n) \log \log n / \log n)$ processors on a CRCW PRAM, which also can be implemented in $O(\log n \log \log n / \alpha(m, n))$ time using $O(m + n)\alpha(m, n) / \log n)$ processors by Brent's theorem.

CONCLUSIONS 6

In this paper, we have studied the problem of fully dynamic maintenance of k-connectivity of graphs and presented the first NC algorithms with O(n) work for it, where k = 2, 3. In particular, our NC algorithm for 2-edge connectivity uses o(n) processors only. For the biconnectivity problem, we have obtained an NC algorithm with O(n) processors. From this algorithm, we have also derived NC algorithms with O(n) processors for the fully dynamic maintenance of 3-edge connectivity and triconnectivity. However, for the addressed problems there is still a gap between the best sequential time complexity for them and the amount of work needed in parallel. It is interesting and challenging to find better NC algorithms for them to reduce or eliminate the gap.

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