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Faster Worst Case Deterministic Dynamic Connectivity

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Abstract

We present a deterministic dynamic connectivity data structure for undirected graphs with worst case update time $O\left(\frac{n(\log \log n)^2}{\log n}\right)$ and constant query time. This improves on the previous best deterministic worst case algorithm of Frederickson (SIAM J. Comput., 1985) and Eppstein Galil, Italiano, and Nissenzweig (J. ACM, 1997), which had update time $O(\sqrt{n})$. All other algorithms for dynamic connectivity are either randomized (Monte Carlo) or have only amortized performance guarantees.

1998 ACM Subject Classification F.2.2 Nonnumerical Algorithms and Problems, G.2.2 Graph Theory

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1 Introduction

Dynamic Connectivity is perhaps the single most fundamental unsolved problem in the area of dynamic graph algorithms. The problem is simply to maintain a dynamic undirected graph $G = (V,E)$ subject to edge updates and connectivity queries:

$\text{INSERT}(u, v) : \quad \text{Set } E \leftarrow E \cup \{(u, v)\}$.

$\text{DELETE}(u, v) : \quad \text{Set } E \leftarrow E \setminus \{(u, v)\}$.

$\text{CONN?}(u, v) : \quad \text{Determine whether } u \text{ and } v \text{ are in the same connected component in } G$.

Over thirty years ago Frederickson [10] introduced topology trees and 2-dimensional topology trees, which gave the first non-trivial solution to the problem. Each edge insertion/deletion is handled in $O(\sqrt{m})$ time and each query is handled in $O(1)$ time. Here $m$ is the current number of edges and $n$ the number of vertices. On sparse graphs (where

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\( m = O(n) \) Frederickson’s data structure has not been improved by any deterministic worst case algorithm. However, when the graph is dense Frederickson’s data structure can be improved using the general sparsification method of Eppstein, Galil, Italiano, and Nissenzweig [7]. Using simple sparsification [6] the update time becomes \( O(\sqrt{n} \log(m/n)) \) and using more sophisticated sparsification [7] the running time becomes \( O(\sqrt{n}) \). This last bound has not been improved in twenty years.

1.1 New Results

In this paper we return to the classical model of deterministic worst case complexity. We give a new dynamic connectivity structure with worst case update time on the order of

\[
\min \left\{ \sqrt{\frac{m (\log \log n)^2}{\log n}}, \sqrt{\frac{m \log^5 w}{w}} \right\},
\]

where \( w = \Omega(\log n) \) is the word size.\(^1\) These are the first improvements to Frederickson’s 2D-topology trees [10] in over 30 years. Using the sparsification reduction of Eppstein et al. [7] the running time expressions can be made to depend on ‘\( n \)’ rather than ‘\( m \)’, so we obtain \( O(\sqrt{\frac{m (\log \log n)^2}{\log n}}) \) bounds (or faster) for all graph densities.

1.2 Related Work

Most research on the dynamic connectivity problem has settled for amortized update time guarantees. Following [15, 16], Holm et al. [17] gave a very simple deterministic algorithm with amortized update time \( O(\log^2 n) \) and query time \( O(\log n / \log \log n) \).\(^2\) However, in the worst case Holm et al.’s [17] update takes \( \Omega(m) \) time, the same as computing a spanning tree from scratch! Recently Wulff-Nilsen [25] improved the update time of [17] to \( O(\log^2 n / \log \log n) \). Using Las Vegas randomization, Thorup [24] gave a dynamic connectivity data structure with an \( O(\log n (\log \log n)^3) \) amortized update time. In other words, the algorithm answers all connectivity queries correctly but the amortized update time holds with high probability.

In a major breakthrough Kapron, King, and Mountjoy [18] used Monte Carlo randomization to achieve a worst case update time of \( O(\log^4 n) \). However, this algorithm has three notable drawbacks. The first is that it is susceptible to undetected false negatives: \( \text{CONN}?(u, v) \) may report that \( u, v \) are disconnected when they are, in fact, connected. The second is that even when \( \text{CONN}?(u, v) \) (correctly) reports that \( u, v \) are connected, it is forbidden from exhibiting a connectivity witness, i.e., a spanning forest in which \( u, v \) are joined by a path. The Kapron et al. [18] \( \text{algorithm does maintain such a spanning forest internally, but if this witness were made public, a very simple attack could force the algorithm to answer connectivity queries incorrectly. Lastly, the algorithm uses } \Omega(n \log^2 n) \text{ space, which for sparse graphs is superlinear in } m. \) Very recently Gibb et al. [13] reduced the update time of [18] to \( O(\log^4 n) \).

On special graph classes, dynamic connectivity can often be handled more efficiently. For example, Sleator and Tarjan [22] maintain a dynamic set of trees in \( O(\log n) \) worst-case update time subject to \( O(\log n) \) time connectivity queries. (See also [1, 3, 15, 23].)

\(^1\) Our algorithms use the standard repertoire of \( \mathsf{AC}^0 \) operations: left and right shifts, bitwise operations on words, additions and comparisons. They do not assume unit-time multiplication.

\(^2\) Any connectivity structure that maintains (internally) a spanning forest can have query time \( O(\log_{t_u}/\log n) \) if the update time is \( t_u = \Omega(\log n) \).
Table 1 A survey of dynamic connectivity results. The lower bounds hold in the cell probe model with word size \( w = \Theta(\log n) \).

### Worst Case Data Structures

<table>
<thead>
<tr>
<th>Ref.</th>
<th>Update Time</th>
<th>Query Time</th>
<th>Notes</th>
</tr>
</thead>
<tbody>
<tr>
<td>[10]</td>
<td>( O(\sqrt{m}) )</td>
<td>( O(1) )</td>
<td></td>
</tr>
<tr>
<td>[7, 10]</td>
<td>( O(\sqrt{n}) )</td>
<td>( O(1) )</td>
<td>[10] + sparsification [7].</td>
</tr>
<tr>
<td>[18]</td>
<td>( O(c \log^3 n) )</td>
<td>( O\left(\frac{\log n}{\log \log n}\right) ) Randomized Monte Carlo; no connectivity witness; ( n^c ) ops. err with prob. ( n^{-c} ).</td>
<td></td>
</tr>
<tr>
<td>[13]</td>
<td>( O(c \log^4 n) )</td>
<td>( O\left(\frac{\log n}{\log \log n}\right) )</td>
<td></td>
</tr>
</tbody>
</table>

\( w = \Omega(\log n) \)

### Amortized Data Structures

<table>
<thead>
<tr>
<th>Ref.</th>
<th>Amort. Update</th>
<th>W.C. Query</th>
<th>Notes</th>
</tr>
</thead>
<tbody>
<tr>
<td>[15]</td>
<td>( O(\log^3 n) )</td>
<td>( O\left(\frac{\log n}{\log \log n}\right) ) Randomized Las Vegas.</td>
<td></td>
</tr>
<tr>
<td>[16]</td>
<td>( O(\log^2 n) )</td>
<td>( O\left(\frac{\log n}{\log \log n}\right) ) Randomized Las Vegas.</td>
<td></td>
</tr>
<tr>
<td>[17]</td>
<td>( O(\log^2 n) )</td>
<td>( O\left(\frac{\log n}{\log \log n}\right) )</td>
<td></td>
</tr>
<tr>
<td>[24]</td>
<td>( O(\log n(\log \log n)^3) )</td>
<td>( O\left(\frac{\log n}{\log \log \log n}\right) ) Randomized Las Vegas.</td>
<td></td>
</tr>
<tr>
<td>[25]</td>
<td>( O\left(\frac{\log^2 n}{\log \log n}\right) )</td>
<td>( O\left(\frac{\log n}{\log \log n}\right) )</td>
<td></td>
</tr>
</tbody>
</table>

### Amort./Worst Case Lower Bounds

<table>
<thead>
<tr>
<th>Ref.</th>
<th>Update Time ( t_u )</th>
<th>Query Time ( t_q )</th>
<th>Notes</th>
</tr>
</thead>
<tbody>
<tr>
<td>[11, 14, 19]</td>
<td>( t_q = \Omega\left(\frac{\log n}{\log(t_u \log n)}\right) )</td>
<td></td>
<td></td>
</tr>
<tr>
<td>[20]</td>
<td>( t_u = \Omega\left(\frac{\log n}{\log(t_q / t_u)}\right) ) ( t_q = \Omega\left(\frac{\log n}{\log(t_u / t_q)}\right) ) Implies ( \max{t_u, t_q} = \Omega(\log n) ).</td>
<td></td>
<td></td>
</tr>
<tr>
<td>[21]</td>
<td>( o(\log n) ) implies ( \Omega(n^{1-o(1)}) )</td>
<td></td>
<td></td>
</tr>
</tbody>
</table>
Connectivity in dynamic planar graphs can be reduced to the dynamic tree problem \cite{8, 9}, and therefore solved in $O(\log n)$ time per operation. The cell probe lower bounds of Pătraşcu and Demaine \cite{20} show that Sleator and Tarjan’s algorithm is optimal in the sense that some operation must take $\Omega(\log n)$ time. Superlogarithmic updates can be used to get modestly sublogarithmic queries, but Pătraşcu and Thorup \cite{21} prove the reverse is not possible. In particular, any dynamic connectivity algorithm with $o(\log n)$ update time has $n^{1-o(1)}$ query time. Refer to Table 1 for a history of upper and lower bounds for dynamic connectivity.

Compared to the ammortized algorithms \cite{17, 24, 25}, ours is better suited to online applications that demand a bound on the latency of every operation. 3 Compared to the Monte Carlo algorithms \cite{13, 18}, ours is attractive in applications that demand linear space, zero probability of error, and a public witness of connectivity.

## 2 The High Level Algorithm

The algorithm maintains a spanning tree of each connected component of the graph as a witness of connectivity. Each such witness tree $T$ is represented as an Euler tour $\text{Euler}(T)$\(^4\). Euler($T$) is the sequence of vertices encountered in some Euler tour around $T$, as if each undirected edge were replaced by two oriented edges. It has length precisely $2(|V(T)| - 1)$ if $|V(T)| \geq 2$ (the last vertex is excluded from the list, which is necessarily the same as the first) or length 1 if $|V(T)| = 1$. Vertices may appear in Euler($T$) several times. We designate one copy of each vertex the principal copy, which is responsible for all edges incident to the vertex. Each vertex in the graph maintains a pointer to its principal copy. Each $T$-edge $(u, v)$ maintains two pointers to the (possibly non-principal) copies of $u$ and $v$ that precede the oriented occurrences of $(u, v)$ and $(v, u)$ in Euler($T$), respectively. Note that cyclic rotations of Euler($T$) are also valid Euler tours; if Euler($T$) = $(u, \ldots, v)$ the last element of the list is associated with the tree edge $(v, u)$.

When an edge $(u, v)$ that connects distinct witness trees $T_0$ and $T_1$ is inserted, $(u, v)$ becomes a tree edge and we need to construct Euler($T_0 \cup \{(u, v)\} \cup T_1$) from Euler($T_0$) and Euler($T_1$). In the reverse situation, if a tree edge $(u, v)$ is deleted from $T = T_0 \cup \{(u, v)\} \cup T_1$ we first construct Euler($T_0$) and Euler($T_1$) from Euler($T$), then look for a replacement edge, $(\tilde{u}, \tilde{v})$ with $\tilde{u} \in V(T_0)$ and $\tilde{v} \in V(T_1)$. If a replacement is found we construct Euler($T_0 \cup \{(\tilde{u}, \tilde{v})\} \cup T_1$) from Euler($T_0$) and Euler($T_1$). Lemma 1 establishes the nearly obvious fact that the new Euler tours can be obtained from the old Euler tours using $O(1)$ of the following surgical operations: splitting and concatenating lists of vertices, and creating and destroying singleton lists containing non-principal copies of vertices.

\begin{lemma}
If $T = T_0 \cup \{(u, v)\} \cup T_1$ and $(u, v)$ is deleted, Euler($T_0$) and Euler($T_1$) can be constructed from Euler($T$) with $O(1)$ surgical operations. In the opposite direction, from Euler($T_0$) and Euler($T_1$) we can construct Euler($T_0 \cup \{(u, v)\} \cup T_1$) with $O(1)$ surgical operations. It takes $O(1)$ time to determine which surgical operations to perform.
\end{lemma}

\textbf{Proof.} Recall that cyclic shifts of Euler tours are valid Euler tours. Suppose without loss of generality that Euler($T$) = $(P_0, u, v, P_1, v, u, P_2)$ where $P_0$, $P_1$, and $P_2$ are sequences of

---

3 Amortized data structures are most useful when employed by offline algorithms that do not care about individual operation times. The canonical example is the use of amortized Fibonacci heaps \cite{12} to implement Dijkstra’s algorithm \cite{5}.

4 Henzinger and King \cite{15} were the first to use Euler tours to represent dynamic trees. G. Italiano (personal communication) observed that Euler tours could be used in lieu of Frederickson’s topology trees to obtain an $O(\sqrt{n})$-time dynamic connectivity structure.
vertices. (Note that Euler tours never contain immediate repetitions. If $P_1$ is empty then 
$\text{Euler}(T)$ would be just $(P_0, u, v, u, P_2)$; if both $P_0$ and $P_2$ are empty then $\text{Euler}(T) = (u, v, P_1, v).$ Then we obtain $\text{Euler}(T_0) = (P_0, u, P_2)$ and $\text{Euler}(T_1) = (v, P_1)$ with $O(1)$ surgical operations, which includes the destruction of non-principal copies of $u$ and $v$; at least one of the two copies must be non-principal. We could also set $\text{Euler}(T_1) = (P_1, v),$ which would be more economical if the $v$ following $P_1$ in $\text{Euler}(T)$ were the principal copy.

In the reverse direction, write $\text{Euler}(T_0) = (P_0, u, P_1)$ and $\text{Euler}(T_1) = (P_2, v, P_3),$ where the labeled occurrences are the principal copies of $u$ and $v.$ Then $\text{Euler}(T_0 \cup \{(u, v)\} \cup T_1) = (P_0, u, v, P_3, P_2, v, u, P_1),$ where the new copies of $u$ and $v$ are clearly non-principal copies. If $P_2$ and $P_3$ were empty (or $P_0$ and $P_1$ were empty) then we would not need to add a non-principal copy of $v$ (or a non-principal copy of $u.$)

Thus, we have reduced dynamic connectivity in graphs to implementing several simple operations on dynamic lists. Our algorithm maintains a pair $(\mathcal{L}, E),$ where $\mathcal{L}$ is a set of lists (containing principal and non-principal copies of vertices) and $E$ is the dynamic set of edges joining principal copies of vertices. In addition to the creation and destruction of single element lists we must support the following primitive operations.

\begin{itemize}
  \item **List** $(x):$ Return the list in $\mathcal{L}$ containing element $x.$
  \item **Join** $(L_0, L_1):$ Set $\mathcal{L} \leftarrow \mathcal{L} \setminus \{L_0, L_1\} \cup \{L_0L_1\},$ that is, replace $L_0$ and $L_1$ with their concatenation $L_0L_1.$
  \item **Split** $(x):$ Let $L = L_0L_1 \in \mathcal{L},$ where $x$ is the last element of $L_0.$ Set $\mathcal{L} \leftarrow \mathcal{L} \setminus \{L\} \cup \{L_0, L_1\}.$
  \item **ReplacementEdge** $(L_0, L_1):$ Return any edge joining elements in $L_0$ and $L_1.$
\end{itemize}

Our implementations of these operations will only be efficient if, after each Insert or Delete operation, there are no edges connecting distinct lists. That is, the ReplacementEdge operation is only employed by Delete when deleting a tree edge in order to restore Invariant 2.

\begin{itemize}
  \item **Invariant 2.** Each list $\mathcal{L}$ corresponds to the Euler tour of a spanning tree of some connected component.
\end{itemize}

The dynamic connectivity operations are implemented as follows. To answer a Conn?(u, v) query we simply check whether List(u) = List(v). To insert an edge (u, v) we do Insert(u, v), and if List(u) ≠ List(v) then make (u, v) a tree edge and perform suitable Splits and Joins to merge the Euler tours List(u) and List(v). To delete an edge (u, v) we do Delete(u, v), and if (u, v) is a tree edge in $T = T_0 \cup \{(u, v)\} \cup T_1,$ perform suitable Splits and Joins to create Euler($T_0$) and Euler($T_1$) from Euler($T$). At this point Invariant 2 may be violated as there could be an edge joining $T_0$ and $T_1.$ We call ReplacementEdge(Euler($T_0$), Euler($T_1$)) and if it finds an edge, say $(\hat{u}, \hat{v}),$ we perform more Splits and Joins to form Euler($T_0$ ∪ $(\hat{u}, \hat{v})$) ∪ $T_1$).

Henzinger and King [15] observed that most off-the-shelf balanced binary search trees can support Split, Join, and other operations in logarithmic time. However, they provide no direct support for the ReplacementEdge operation, which is critical for the dynamic connectivity application.

Section 3 gives a relatively simple instantiation of the high-level approach with update time $O(\sqrt{n} / w^{1/4}),$ $w = \Omega(\log n)$ being the word size. This is slightly slower than our claimed result. In Section 4 we describe the modifications needed to achieve the claimed bounds.
3 A New Data Structure for Dynamic Lists

3.1 Chunks and Superchunks

In order to simplify the maintenance of Invariant 3, stated below, we shall make two simplifying assumptions. We assume that we have a fixed upper bound $\hat{m}$ on the number of edges and that the maximum degree never exceeds $K$, where $K \approx \sqrt{\hat{m}}/\text{poly}(w)$. The first assumption is justified by the fact that the sparsification method of [7] creates instances in which $\hat{m}$ is known to be linear in the number of vertices. (It can also be removed by the standard technique of periodic rebuilding.) Refer to Section 5.1 for clean ways to remove the degree-bound assumption.

If $L'$ is a sublist of a list $L \in \mathcal{L}$, define $\text{mass}(L')$ to be the number of edges incident to elements of $L'$, counting an edge twice if both endpoints are in $L'$. The sum of list masses, $\sum_{L \in \mathcal{L}} \text{mass}(L)$, is clearly at most $2\hat{m}$, where $\hat{m}$ is the fixed upper bound on the number of edges. We maintain a partition of each list $L \in \mathcal{L}$ into chunks satisfying Invariant 3.

▶ Invariant 3. Let $L \in \mathcal{L}$ be an Euler tour. If $\text{mass}(L) < K$ then $L$ consists of a single chunk. Otherwise $L = C_0C_1\cdots C_{p-1}$ is partitioned into $\Theta(\text{mass}(L)/K)$ chunks such that $\text{mass}(C_l) \in [K, 3K]$ for all $l \in [p] \overset{\text{def}}{=} \{0, \ldots, p-1\}$.

The chunks are partitioned into contiguous sequences of $\Theta(h)$ superchunks according to Invariant 4. For the time being define $h = 2\lceil\sqrt{w}/2\rceil$, where $w$ is the word size.

▶ Invariant 4. A list in $\mathcal{L}$ having fewer than $h/2$ chunks forms a single superchunk with ID $\perp$. A list in $\mathcal{L}$ with at least $h/2$ chunks is partitioned into superchunks, each consisting of between $h/2$ and $h-1$ consecutive chunks. Each such superchunk has a unique ID in $[J] \overset{\text{def}}{=} \{0, \ldots, J-1\}$, where $J = 4\hat{m}/(Kh)$. (IDs are completely arbitrary. They do not encode any information about the order of superchunks within a list.)

Call an Euler tour list short if it consists of fewer than $h/2$ chunks. We shall assume that no lists are ever short, as this simplifies the description of the data structure and its analysis. In particular, all superchunks have proper IDs in $[J]$. Refer to Section 5.2 for a description of how to handle $\perp$ IDs and short lists.

3.2 Word Operations

When $h \leq \lceil\sqrt{w}\rceil$, Invariant 4 implies that we can store a matrix $A \in \{0, 1\}^{h \times h}$ in one word that represents the adjacency between the chunks within two superchunks $i$ and $j$. This matrix will always be represented in row-major order; rows and columns are indexed by $[h] = \{0, \ldots, h-1\}$. In this format it is straightforward to insert a new all-zero row above a specified row $k$ (and destroy row $h-1$) by shifting the old rows $k, \ldots, h-2$ down by one. It is also easy to copy an interval of rows from one matrix to another. Lemma 5 shows that the corresponding operations on columns can also be effected in $O(1)$ time with a fixed mask $\mu$ precomputable in $O(\log w)$ time.

▶ Lemma 5. Let $h = 2\lceil\sqrt{w}/2\rceil$ and let $\mu$ be the word $(1^h 0^{h/2})^h$. Given $\mu$ we can in $O(1)$ time copy/paste any interval of columns from/to a matrix $A \in \{0, 1\}^{h \times h}$, represented in row-major order.

---

5 Remember that edges are only incident to principal copies of vertices, so non-principal copies never contribute any mass.
Proof. Recall that the rows and columns are indexed by integers in $[h] = \{0, \ldots, h-1\}$. We first describe how to build a mask $\nu_k$ for columns $k, \ldots, h-1$ then illustrate how it is used to copy/paste intervals of columns. In C notation,\(^6\) the word $\nu'_k = (\mu \gg k) \& \mu$ is a mask for the intersection of the even rows and columns $k, \ldots, h-1$, so $\nu_k = \nu'_k \mid (\nu'_k \gg h)$ is a mask for columns $k$ through $h-1$.

To insert an all-zero column before column $k$ of $A$ (and delete column $h-1$) we first copy columns $k, \ldots, h-2$ to $A' = A \& (\nu_{k+1} \ll 1)$ then set $A = (A \& (\sim \nu_k)) \mid (A' \gg 1)$. Other operations can be effected in $O(1)$ time with copying/pasting intervals of columns, e.g., splitting an array into two about a designated column, or merging two arrays having at most $h$ columns together.

### 3.3 Adjacency Data Structures

In order to facilitate the efficient implementation of ReplacementEdge we maintain an $O(\hat{m}/K) \times O(\hat{m}/K)$ adjacency matrix between chunks, and a $J \times J$ adjacency matrix between superchunks. However, in order to allow for efficient dynamic updates it is important that these matrices be represented in a non-standard format described below. The data structure maintains the following information:

- Each list element maintains a pointer to the chunk containing it. Each chunk maintains a pointer to the superchunk containing it, as well as an index in $[h]$ indicating its position within the superchunk. Each superchunk maintains its ID in $[J] \cup \{\perp\}$ and a pointer to the list containing it.

- ChAdj is a $J \times J$ array of $h^2$-bit words ($h^2 \leq w$) indexed by superchunk IDs. The entry ChAdj$(i, j)$ is interpreted as an $h \times h$ 0-1 matrix that keeps the adjacency information between all pairs of chunks in superchunk $i$ and superchunk $j$. (It may be that $i = j$.) In particular, ChAdj$(i, j)(k, l) = 1$ iff there is an edge with endpoints in the $k$th chunk of superchunk $i$ and the $l$th chunk of superchunk $j$, so ChAdj$(i, j) = 0$ (i.e., the all-zero matrix) if no edge joins superchunks $i$ and $j$. The matrix ChAdj$(i, j)$ is stored in row-major order.

- Let $S$ be a superchunk with $ID(S) = \perp$. By Invariants 2 and 4, $S$ is not incident to any other superchunks and has fewer than $h/2$ chunks. We maintain a single word ChAdj$_S$ which stores the adjacency matrix of the chunks within $S$.

- For each superchunk with ID $i \in [J]$ we keep length-$J$ bit-vectors SupAdj$_i$ and Memb$_i$, where

$$\begin{align*}
\text{SupAdj}_i(j) &= 1 \text{ if } \text{ChAdj}(i, j) \neq 0 \text{ and } 0 \text{ otherwise, whereas} \\
\text{Memb}_i(j) &= 1 \text{ if } j = i \text{ and } 0 \text{ otherwise.}
\end{align*}$$

These vectors are packed into $[J/w]$ machine words, so scanning one takes $O([J/w])$ time.

- We maintain a list-sum data structure that allows us to take the bit-wise OR of the SupAdj$_i$ vectors or Memb$_i$ vectors, over all superchunks in an Euler tour. It is responsible for maintaining the \{SupAdj$_i$, Memb$_i$\} vectors described above and supports the following operations. At all times the superchunks are partitioned into a set $S$ of disjoint lists of superchunks. Each $S \in S$ (a list of superchunks) is associated with an $L \in \mathcal{L}$ (an Euler tour), though short lists in $\mathcal{L}$ have no need for a corresponding list in $S$.

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\(^6\) The operations $\&$, $\ll$, and $\sim$ are bit-wise AND, OR, and NOT; $\ll$ and $\gg$ are left and right shift.
SCInsert(i): Retrieve an unused ID, say i', and allocate a new superchunk with ID i' and all-zero vector SupAdj_i. Insert superchunk i' immediately after superchunk i in i's list in S. If no i is given, create a new list in S consisting of superchunk i'.

SCDelete(i): Delete superchunk i from its list and make ID i unused.

SCJoin(S_0, S_1): Replace superchunk lists S_0, S_1 ∈ S with their concatenation S_0S_1.

SCSplit(i): Let S = S_0S_1 ∈ S and i be the last superchunk in S_0. Replace S_0S_1 with two lists S_0, S_1.

UPDATEAdj(i, x ∈ \{0, 1\}^J): Set SupAdj_i ← x and update SupAdj_j(i) ← x(j) for all j ≠ i.

ADQUERY(S): Return the vector α ∈ \{0, 1\}^J where

\[ \alpha(j) = \bigvee_{i \in S} \text{SupAdj}_i(j) \]

The index i ranges over the IDs of all superchunks in S.

MEMBQUERY(S): Return the vector β ∈ \{0, 1\}^J, where

\[ \beta(j) = \bigvee_{i \in S} \text{Memb}_i(j) \]

We use the following implementation of the list-sum data structure. Each list of superchunks is maintained as any \(O(1)\)-degree search tree that supports logarithmic time inserts, deletes, splits, and joins. Each leaf is a superchunk that stores its two bit-vectors. Each internal node z keeps two bit-vectors, SupAdj^z and Memb^z, which are the bit-wise OR of their leaf descendants' respective bit-vectors. Because length-J bit-vectors can be updated in \(O([J/w])\) time, all “logarithmic time” operations on the tree actually take \(O(\log J \cdot J/w)\) time. The UPDATEAdj(i, x) operation takes \(O(\log J \cdot J/w)\) time to update superchunk i and its \(O(\log J)\) ancestors. We then need to update the ith bit of potentially every other node in the tree, in \(O(J)\) time. Since \(w = \Omega(\log n) = \Omega(\log J)\) the cost per UPDATEAdj is \(O(J)\). The answer to an ADQUERY(S) or MEMBQUERY(S) is stored at the root of the tree on S.

### 3.4 Creating and Destroying (Super)Chunks

There are essentially two causes for the creation and destruction of (super)chunks. The first is in response to a SPLIT operation that forces a (super)chunk to be broken up. (The SPLIT may itself be instigated by the insertion or deletion of an edge.) The second is to restore Invariants 3 and 4 after a JOIN or INSERT or DELETE operation. In this section we consider the problem of updating the adjacency data structures after four types of operations:

(i) splitting a chunk in two, keeping both chunks in the same superchunk, (ii) merging two adjacent chunks in the same superchunk, (iii) splitting a superchunk along a chunk boundary, and (iv) merging adjacent superchunks. Once we have bounds on (i)-(iv), implementing the higher-level operations in the stated bounds is relatively straightforward. Note that (i)-(iv) may temporarily violate Invariants 3 and 4.

#### 3.4.1 Splitting Chunks

Suppose we want to split the kth chunk of superchunk i into two pieces, both of which will (at least temporarily) stay within superchunk i. We first zero-out all bits of ChAdj(i, *)(k, *)

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7 Remember that ‘k’ refers to the actual position of the chunk within its superchunk whereas ‘i’ is an arbitrary ID that does not relate to its position within the list.
and ChAdj(\(i, k\)) in \(O(J)\) time. For each \(j\) we need to insert an all-zero row below row \(k\) in ChAdj(i, j) and an all-zero column after column \(k\) of ChAdj(j, i). This can be done in \(O(1)\) time for each \(j\), or \(O(J)\) in total; see Lemma 5.

In \(O(K)\) time we scan the edges incident to the new chunks \(k\) and \(k + 1\) and update the corresponding bits in ChAdj(i, *) for \(k'\) for \(k' \in \{k, k + 1\}\).

### 3.4.2 Merging Adjacent Chunks

In order to merge chunks \(k\) and \(k + 1\) of superchunk \(i\) we need to replace row \(k\) of ChAdj(i, j), for all \(j\), with the bit-wise OR of rows \(k\) and \(k + 1\) of ChAdj(i, j), zero out row \(k + 1\), then scoot rows \(k + 2, \ldots\) back one row. A similar transformation is performed on columns \(k\) and \(k + 1\) of ChAdj(j, i), which takes \(O(1)\) time per \(j\), by Lemma 5. In total the time is \(O(J)\), independent of \(K\).

### 3.4.3 Splitting Superchunks

Suppose we want to split superchunk \(i\) after its \(k\)th chunk. We first call SCINSERT(i), which allocates an empty superchunk with ID \(i'\) and inserts \(i'\) after \(i\) in its superchunk list in \(S\). In \(O(J)\) time we transfer rows \(k + 1, \ldots, h - 1\) from ChAdj(i, j) to ChAdj(i', j) and transfer columns \(k + 1, \ldots, h - 1\) from ChAdj(j, i) to ChAdj(j, i'). By Lemma 5 this takes \(O(1)\) time per \(j\).

At this point ChAdj is up-to-date but the list-sum data structure and \{SupAdj\} bit-vectors are not. We update SupAdj\(_j\), SupAdj\(_{i'}\) with calls to UPDATEADJ(i, x) and UPDATEADJ(i', x'). Using ChAdj, each bit of \(x\) and \(x'\) can be generated in constant time. This takes \(O(J)\) time.

### 3.4.4 Merging Superchunks

Let the two adjacent superchunks have IDs \(i\) and \(i'\). It is guaranteed that they will be merged only if they contain at most \(h\) chunks together. In \(O(J)\) time we transfer the non-zero rows of ChAdj(i', j) to ChAdj(i, j) and transfer the non-zero columns of ChAdj(j, i') to ChAdj(j, i). A call to SCDELETE(i') deletes superchunk \(i'\) from its list in \(S\) and retires ID \(i'\). We then call UPDATEADJ(i', x) with the new incidence vector \(x\). In this case we can generate \(x\) in \(O(J/w)\) time since it is merely the bit-wise OR of the old vectors SupAdj\(_j\) and SupAdj\(_{i'}\), with bit \(i'\) set to zero. Updating the list-sum data structure takes \(O(J)\) time.

### 3.5 Joining and Splitting Lists

Once we have routines for splitting and merging adjacent (super)chunks, implementing JOIN and SPLIT on lists in \(L\) is much easier. The goal is to restore Invariant 3 governing chunk masses and Invariant 4 on the number of chunks per superchunk.

#### 3.5.1 Performing JOIN(L₀, L₁)

Write \(L₀ = C₀, \ldots, C_{p-1}\) and \(L₁ = D₀, \ldots, D_{q-1}\) as a list of chunks. If both \(L₀\) and \(L₁\) are not short then they have corresponding superchunk lists \(S₀, S₁ \in S\). Call SCJOIN(S₀, S₁) to join \(S₀, S₁\) in \(S\), in \(O(J)\) time.
3.5.2 Performing Split($x$)

Suppose $x$ is contained in chunk $C_1$ of $L = C_0 \cdots C_{t-1} C_t C_{t+1} \cdots C_{p-1}$. We split $C_1$ into two chunks $C'_1 C''_1$, and split the superchunk containing $C_1$ along this line. Let $S$ be the superchunk list corresponding to $L$ and $i$ be the ID of the superchunk ending at $C'_1$. We split $S$ using a call to SC$\text{SPLIT}(i)$, which corresponds to splitting $L$ into $L_0 = C_0 \cdots C_{t-1} C'_1$ and $L_1 = C''_1 C_{t+1} \cdots C_{p-1}$. At this point $C'_1$ or $C''_1$ may violate Invariant 3 if mass($C''_1$) < $K$ or mass($C'_1$) < $K$. Furthermore, Invariant 4 may be violated if the number of chunks in the superchunks containing $C'_1$ and $C''_1$ is too small. We first correct Invariant 3 by possibly merging and resplitting $C_{t-1} C'_1$ and $C''_1 C_{t+1}$ along new boundaries. If the superchunk containing $C'_1$ has fewer than $h/2$ chunks, it and the superchunk to its left have strictly between $h/2$ and $3h/2$ chunks together, and so can be merged (and possibly resplit) into one or two superchunks satisfying Invariant 4. The same method can correct a violation of $C''_1$'s superchunk. This takes $O(K + J)$ time.

3.5.3 Performing ReplacementEdge($L_0, L_1$)

The list-sum data structure makes implementing the ReplacementEdge($L_0, L_1$) operation easy. Let $S_0$ and $S_1$ be the superchunk lists corresponding to Euler tours $L_0$ and $L_1$. We compute the vectors $\alpha \leftarrow \text{ADJQUERRY}(S_0)$ and $\beta \leftarrow \text{MEMBQUERY}(S_1)$ and their bit-wise AND $\alpha \land \beta$ with a linear scan of both vectors. If $\alpha \land \beta$ is the all-zero vector then there is no edge between $L_0$ and $L_1$. On the other hand, if $(\alpha \land \beta)(j) = 1$, then $j$ must be the ID of a superchunk in $S_1$ that is incident to some superchunk in $S_0$. To determine which superchunk in $S_0$ we walk down from the root of $S_0$’s list-sum tree to a leaf, say with ID $i$, in each step moving to a child $z$ of the current node for which SupAdj$^z(j) = 1$. Once $i$ and $j$ are known we retrieve any 1-bit in the matrix ChAdj($i, j$), say at position $(k, l)$, indicating that the $k$th chunk of superchunk $i$ and the $l$th chunk of superchunk $j$ are adjacent. We scan all its adjacent edges in $O(K)$ time and retrieve an edge joining $L_0$ and $L_1$. The total time is $O(J/w + \log J + K) = O(J/w + K)$.

3.5.4 Performing Insert($u, v$)

If List($u$) $\neq$ List($v$), first perform $O(1)$ Splits and Joins to restore the Euler tour Invariant 2. Now $u$ and $v$ are in the same list in $L$. Let $i, j$ be the IDs of the superchunks containing the principal copies of $u$ and $v$ and let $k, l$ be the positions of $u$ and $v$’s chunks within their respective superchunks. We set ChAdj$^{ij}(k, l) \leftarrow 1$. If ChAdj$^{ij}$ was formerly the all-zero matrix, we call UpdateAdj$^{ij}(i, x)$ to update superchunk $i$’s adjacency information with the correct vector $x$. Inserting one edge changes the mass of the chunks containing $u$ and $v$, which could violate Invariant 3. Invariants 3 and 4 are restored by splitting/merging $O(1)$ chunks and superchunks.

3.5.5 Performing Delete($u, v$)

Compute $i, j, k, l$ as defined above, in $O(1)$ time. After we delete ($u, v$) the correct value of the bit ChAdj$^{ij}(k, l)$ is uncertain. We scan chunk $k$ of superchunk $i$ in $O(K)$ time, looking for an edge connected to chunk $l$ of superchunk $j$. If we do not find such an edge we set ChAdj$^{ij}(k, l) \leftarrow 0$, and if that makes ChAdj$^{ij}(i, j) = 0$ (the all-zero matrix), we call

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8 Since $x$ only differs from the former SupAdj, at position SupAdj$^{ij}(j)$, this update to the list-sum tree takes just $O(\log J)$ time since it only affects ancestors of leaves $i$ and $j$. 

$$1 + 2$$

which takes time

3.6 Running Time Analysis

The time for a search can actually be improved to 9. Given ChAdj($i, j$), we perform Splits and Joins to replace Euler($T$) with Euler($T_0$), Euler($T_1$), which may violate Invariant 2 if there is a replacement edge between $T_0$ and $T_1$. We call ReplacementEdge(Euler($T_0$), Euler($T_1$)) to find a replacement edge. If one is found, say ($u, \hat{v}$), we form Euler($T_0$) with a constant number of Splits and Joins.

4 Speeding Up the Algorithm

Observe that there are $\Theta((\hat{m}/Kh)^2)$ matrices (ChAdj($i, j$)) but only $\hat{m}$ edges, so for $K = \sqrt{\frac{m}{h}}$, the average $h \times h$ matrix has $O(h)$ 1s. Thus, storing each such matrix verbatim, using $h^2$ bits, is information theoretically inefficient on average. By storing only the locations of the 1s in each matrix we can represent each matrix in $O(h \log h)$ bits on average and thereby hope to solve dynamic connectivity faster with a larger ‘h’ parameter.

4.1 The Encoding

In this encoding we index rows and columns by indices in $\{1, \ldots, h\}$ rather than $[h]$. Let $m_{i,j} = m_{j,i}$ be the number of 1s in ChAdj($i, j$). We encode ChAdj($i, j$) by listing its 1 positions in $O(m_{i,j} \log h/w)$ lightly packed words. Each word is partitioned into fields of $1 + 2\lceil \log(h + 1) \rceil$ bits: each field consists of a control bit (normally 0), a row index, and a column index. Each word is between half-full and full, the fields in use being packed contiguously in the word. This invariant allows us to insert a new field after a given field in $O(1)$ time. We list the 1s of either ChAdj($i, j$) or ChAdj($j, i$) = ChAdj($i, j$)\top in row-major order, with a bit indicating which of the two representations is used.

4.2 Fast Operations

Given ChAdj($i, j$) in row-major order, we can determine if ChAdj($i, j$)($k, l$) = 1 in $O(\log h)$ time by doing a binary search to find the correct word in the list, then a binary search within the word to find an entry ($k, l$), if any.\footnote{The time for a search can actually be improved to $O(\log((m_{i,j} \log h)/w))$, which is faster when ChAdj($i, j$) has average density ($m_{i,j}$ is close to $h$) but is still $O(\log h)$ in the worst case. We still do a binary search over the first field in each word to determine which word (if any) has a field containing ($k, l$): the binary encoding of ($k, l$). This takes $O(\log((m_{i,j} \log h)/w))$ time since there are $\Theta(m_{i,j} \log h/w)$ lightly packed words. If we add $2^{2\lceil \log(h + 1) \rceil} = (k, l)$ to each field in the word, the control bits for all fields that are equal to or greater than ($k, l$) will be flipped to 1. Similarly, if we set all control bits to 1}
In the same time bound we can also identify the positions of the first and last 1s in row $k$. Thus, we can perform the following operations on $\text{ChAdj}(i,j)$ in $O((m_{i,j} \log h)/w)$ time: setting a row to zero, incrementing/decrementing the row-index of some interval of rows, or copying an interval of rows.

The operations sketched above are only efficient if $\text{ChAdj}(i,j)^\top$ is in row-major order. If we have $\text{ChAdj}(i,j)^\top$ in row-major order we can effect a transpose by (1) swapping the row and column indices in each field using masks and shifts, and (2) sorting the fields. In general, sorting $x$ words of $O(w/\log h)$ fields takes $O(x(\log^2(w/\log h) + \log x \log \log(w/\log h)))$ time using Albers and Hagerup’s implementation [2] of Batcher’s bitonic mergesort [4].

We sort each word in $O(\log(\log^2(w/\log h)))$ time, resulting in $x$ sorted lists, then iteratively merge the two shortest lists until one list remains. Merging two lists containing $y$ words takes $O(y \log(w/\log h))$ time: we can merge the next $w/\log h$ fields of each list in $O(\log(w/\log h))$ time [2] and output at least $w/\log h$ items to the merged list.

Alternatively, if $w = \log n$ we can sort and merge lists of $\epsilon \log n/\log h$ fields in unit time using table lookup to precomputed tables of size $O(n \epsilon)$. In this case sorting $x$ packed words takes $O(x \log x)$ time.

### 4.3 Splitting and Joining

The cost of splitting and joining (super)chunks is now slightly more expensive. When handling superchunk $i$ (or any chunk within it) we first put each $\text{ChAdj}(i,j)$ in row-major order, in

$$\sum_{j=1}^{J} O\left(\frac{m_{i,j} \log h}{w} \log^2 h\right) = O(J \log^2 h + (Kh/w) \log^3 h)$$

since, by Invariants 3 and 4, $\sum m_{i,j} = O(Kh)$. Once the relevant superchunks are in the correct format, splitting or joining $O(1)$ (super)chunks takes $O(K \log h + J + (Kh/w) \log h)$ time. Since $J = O(\hat{m}/(Kh))$, the overall update time is

$$O\left(K \log h + \frac{\hat{m} \log^2 h}{K h} + \frac{Kh \log^3 h}{w}\right).$$

Setting $h = w$ and $K = \sqrt{\frac{\hat{m} \log^3 w}{w}}$, the overall time is $O(\sqrt{\frac{\hat{m} \log^3 w}{w}})$. When $w = O(\log n)$ the cost of taking the transpose is cheaper since sorting and merging a packed word takes unit time via table lookup. Setting $h = \log n$, the total time is

$$O\left(K \log \log n + \frac{\hat{m}}{K \log n} + K(\log \log n)^2\right),$$

which is $O(\sqrt{\frac{\hat{m}(\log \log n)^2}{\log n}})$ when $K = \sqrt{\frac{\hat{m}}{\log n(\log \log n)^2}}$.

and subtract $(k, l) + 1$ from each field, the control bits of fields that are equal to or less than $(k, l)$ will be flipped to 0. Thus, we can single out the control bit for an occurrence of $(k, l)$ (if any) with $O(1)$ bit-wise operations. If $(k, l)$ is not present, the control bits reveal the field in the word after which it could be inserted, if we need to set $\text{ChAdj}(i,j)(k, l) \leftarrow 1$.

Albers and Hagerup also require that the fields to be sorted begin with control bits.
5 Loose Ends

5.1 Removing the Bounded Degree Assumption

Invariants 3 and 4 imply that there are \( J = \Theta(\hat{m}/(Kh)) \) superchunks with non-\( \perp \) IDs. However, Invariant 3 cannot be satisfied (as stated) unless the maximum degree is bounded by \( O(K) \). One way to guarantee this is to physically split up high degree vertices, replacing each \( v \) with a cycle on new vertices \( v_1, \ldots, v_{\deg(v)/O(K)} \), each of which is responsible for \( \Theta(K) \) of \( v \)'s edges. This is the method used by Frederickson [10], who actually demanded that the maximum degree be 3 at all times!

This vertex-splitting can be effectively simulated in our algorithm as follows. If \( \deg(v) \geq K/2 \), replace the principal copy of \( v \) in its Euler tour with an interval of artificial principal vertices \( v_1, \ldots, v_{\deg(v)/(K/2)} \), each of which is responsible for between \( K/2 \) and \( K \) of \( v \)'s edges. Invariant 3 is therefore maintained w.r.t. this modified tour. To keep the mass of artificial vertices between \( K/2 \) and \( K \), each edge insertion/deletion may require splitting an artificial vertex or merging two consecutive artificial vertices. When the Euler tour changes we always preserve the invariant that \( v \)'s artificial vertices form a contiguous interval in the tour.

5.2 Dealing with Short Lists

Until now we have assumed for simplicity that all superchunks have proper IDs in \( [J] \). It is important that we not give out IDs to short lists (consisting of less than \( h/2 \) chunks) because the running time of the algorithm is linear in the maximum ID \( J \). The modifications needed to deal with short lists are tedious but minor.

Consider an \( \text{INSERT}(u, v) \) operation where \( u \) and \( v \) are in lists \( L_0 \), \( L_1 \) and \( L_1 \) is a short list consisting of one superchunk \( S \) with \( \text{ID}(S) = \perp \). If \( L_0 \) is not short (or if it is short but the combined list \( L_0L_1 \) will not be short) then we retrieve an unused ID, say \( i \), set \( \text{ID}(S) \leftarrow i \), set \( \text{ChAdj}(i, i) \leftarrow \text{ChAdj}_S \), and destroy \( \text{ChAdj}_S \). By Invariant 2, \( S \) was not incident to any other superchunk, so \( \text{ChAdj}(i, j) = 0 \) (the all-zero matrix) for all \( j \neq i \). At this point \( S \) violates Invariant 4 (it is too small), so we need to merge it with the last superchunk in \( L_0 \) and resplit it along a different chunk boundary, in \( O(1) \) time.

The modifications to \( \text{DELETE}(u, v) \) are analogous. If we delete a tree edge \( (u, v) \), splitting its component into \( T_0 \) and \( T_1 \) having associated Euler tours \( L_0 \) and \( L_1 \), and \( \text{REPLACE}(u, v) \) fails to find an edge joining \( L_0 \) and \( L_1 \), we need to check whether \( L_0 \) (and \( L_1 \)) are short. If so let \( S \) be the superchunk in \( L_0 \). We allocate and set \( \text{ChAdj}_S \leftarrow \text{ChAdj}([\text{ID}(S)/\text{ID}(S)]) \), then set \( \text{ChAdj}([\text{ID}(S)/\text{ID}(S)]) \leftarrow 0 \) and finally retire \( \text{ID}(S) \).

The implementation of \( \text{REPLACE}(L_0, L_1) \) is different if \( L_0 \) and \( L_1 \) were originally in a short list \( L = \text{Euler}(T) \) before a tree edge in \( T \) was deleted. Suppose \( L \) originally had one superchunk \( S \), whose chunk adjacency was stored in \( \text{ChAdj}_S \). After \( O(1) \) splits and joins, both \( L_0 \)'s chunks and \( L_1 \)'s chunks occupy \( O(1) \) intervals of the rows and columns of \( \text{ChAdj}_S \). Of course \( \text{ChAdj}_S \) is represented as a list of its \( 1 \) positions in row-major order, so we can isolate the correct intervals of rows and columns in \( O(h^2 \log^3 h/w) \) time. If there is any \( 1 \) there, say at location \( \text{ChAdj}_S(k, l) \), then we know that there is an edge between \( L_0 \) and \( L_1 \), and can find it in \( O(K) \) time by examining chunks \( k \) and \( l \). The permutation of rows/columns in \( \text{ChAdj}_S \) must be updated to reflect any splits and joins that take place, and if no replacement edge is discovered, \( \text{ChAdj}_S \) must be split into two lists representing matrices \( \text{ChAdj}_{S_0} \) and \( \text{ChAdj}_{S_1} \), to be identified with the single superchunks \( S_0 \) and \( S_1 \) in \( L_0 \) and \( L_1 \), respectively.
References


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