An $O(m\log n)$-Time Algorithm for the Maximal Planar Subgraph Problem

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ABSTRACT

Based on a new version of Hopcroft and Tarjan's planarity testing algorithm, we develop an $O(m\log n)$-time algorithm to find a maximal planar subgraph.

Key words. algorithm, complexity, depth-first-search, embedding, planar graph, selection tree

AMS(MOS) subject classifications. 68R10, 68Q35, 94C15

1. Introduction

In [15], Wu defined the problem of planar graphs in terms of the following four subproblems:

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P1. Decide whether a connected graph G is planar.

P2. Find a minimal set of edges the removal of which will render the remaining part of G planar.

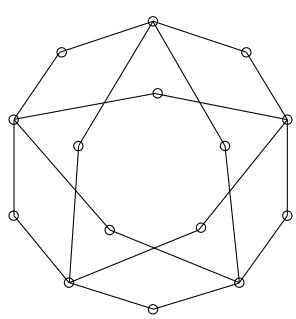
P3. Give a method of embedding G in the plane in case G is planar.

P4. Give a description of the totality of possible planar embeddings of G in the plane in case G is planar.

Linear-time algorithms for P1, P3, and P4 have been known for a long time. The first linear-time solution (which we call the H-T algorithm) for problem P1 (the planarity-testing problem) was given by Hopcroft and Tarjan [7] in 1974 using depth-first search (DFS) trees. A P-Q tree solution for P1 based on an earlier algorithm given by Lempel, Even, and Cederbaum [11] was proved to have a linear-time implementation in 1976 partly by Even and Tarjan [4] and partly by Booth and Lueker [1]. The P-Q tree approach is conceptually simpler, but its implementation is more complicated than that of the H-T algorithm. Linear-time solutions for P3 and P4, also based on P-Q trees, were given by Chiba *et al.* [2] in 1985.

Wu [15] gave an algebraic solution for all four problems. He proved that a graph is planar if and only if a certain system of linear equations is solvable. In case the graph is planar, an actual embedding can be obtained by considering another system of quadratic equations. His solution is elegant, but his algorithm takes $O(m^2)$ time on an m-edge graph.

Recently, Jayakumar *et al.* [9] studied problem P2 (the maximal planar subgraph problem). For the special case in which a biconnected spanning planar subgraph is given, their algorithm runs in $O(n^2)$ time and $O(mn)$ space on a graph with n vertices and m edges. For more general situations, their algorithm runs in $O(mn)$ time. Their algorithm is also based on P-Q trees. Note that not every biconnected graph has a biconnected spanning planar subgraph (See Fig. 1.)



A biconnected graph that does not have a
biconnected spanning planar subgraph

Fig. 1

In this paper we give an $O(m\log n)$-time and $O(m)$-space solution to P2. For sparse graphs (*i.e.*, graphs with $m = O(n^{1+\varepsilon})$, where $\varepsilon < 1$), it beats the algorithm of Jayakumar *et al.* even in the special case when a biconnected spanning planar subgraph is given. Independent of our work, Di Battista and Tamassia [3] have claimed an $O(logn)$-time-per-operation solution to the problem of maintaining a planar graph under edge additions. Their algorithm also solves the minimal planar subgraph problem in $O(m\log n)$ time. Their method is much more complicated than ours, however, as it is designed to solve a more general problem. Recently, Kobayashi *et al.* [10] have shown that if a Hamiltonian tour of the graph is given, then P2 can be solved in liner lime. We show that this result can be easily derived from our algorithm as a special case.

The maximal planar subgraph problem is closely related to the planarity-testing problem. In fact, a graph is planar iff it is the maximal planar subgraph of itself. Our solution to the maximal planar subgraph problem is based on the H-T algorithm. But for our purpose, we need to modify the algorithm. The main difference is that our version of the algorithm admits a more general ordering than the original H-T algorithm does in processing the successors of each tree edge. Also, the H-T algorithm processes one path at a time, while our algorithm processes one edge at a time. In this sense, our algorithm is a more recursive version of the H-T algorithm.

For the above reason, many of our lemmas and theorems are similar, but not identical, to those in [7]. Instead of referring the readers to [7] for the proofs, we find it more convenient and accurate to supply all main proofs in this paper.

The rest of this paper is organized as follows. Section 2 gives preliminary definitions. Section 3 is a new version of the H-T planarity testing algorithm, which leads to our maximal planar subgraph algorithm in Section 4. Section 5 is a summary.

2. Preliminaries

Consider an undirected graph $G_0 = (V_0, E_0)$ with edge set E_0 and vertex set V_0. Let $n = |V_0|$ and $m = |E_0|$. We can draw a picture G_0' of G_0 in the plane as follows: for each vertex $v \in V_0$, we draw a distinct point v'; for each edge $(u, v) \in E_0$, we draw an simple arc connecting the two points u' and v'. We call this arc an *embedding* of the edge (u, v). For brevity, we will sometimes identify graphs with their pictures thus drawn on the plane. If no arcs of G_0' cross each other, then we call G_0' a *planar embedding*, or simply an *embedding*, of G_0. If G_0 has an embedding, then we say that G_0 is *planar*.

The following facts are important to our discussion:

Observation 1. Let C be a simple closed curve in the plane as in Fig. 2; let a be a point inside C and b be a point outside C. Then any curve that joins a and b crosses C.

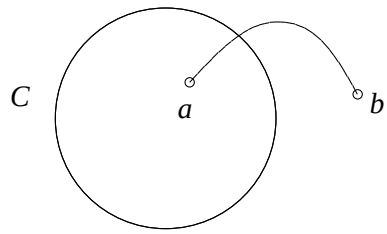


Fig. 2

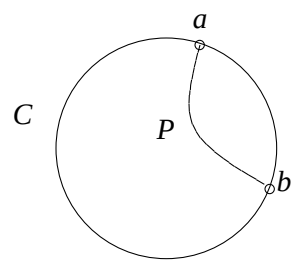


Fig. 3

Observation 2. Let G_1 be the undirected graph represented by Fig. 3, in which P is a path joining the two vertices a and b on cycle C. Then in any embedding of G_1, all the edges of path P are on the same side of the cycle C (either inside or outside).

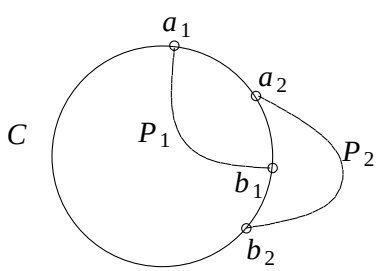


Fig. 4

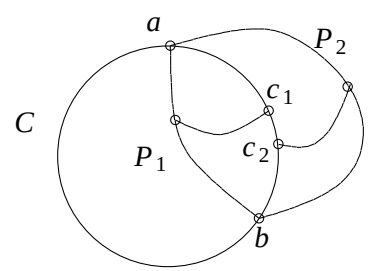


Fig. 5

Observation 3. Let G_2 be the undirected graph represented by Fig. 4, in which a_1, a_2, b_1 and b_2 are four distinct vertices that appear in order on C. Then in any embedding of G_2, the two paths P_1 and P_2 are on opposite sides of the cycle C.

Observation 4. Let G_3 be the undirected graph represented by Fig. 5, in which a, c_1, c_2 and b are vertices that appear in order on C, and c_1 and c_2 may be the same. Then in any embedding of G_3, the two subgraphs P_1 and P_2 are on opposite sides of the cycle C.

All four observations above are intuitively obvious and follow from the Jordan Curve Theorem [6, 14].

A depth-first-search (*abbr*. DFS) [7] will convert the undirected graph $G_0 = (V_0, E_0)$ into a directed graph $G = (V, T, B)$, where V is the set of DFS numbers of vertices in V_0, T is the set of tree edges, and B is the set of back edges. Each edge of G_0 is converted into either a tree edge or a back edge. All the tree edges form a DFS forest. If $[a, b]$ is a tree edge, then $a < b$. If $[a, b]$ is a back edge, then $b < a$, and there is a tree path in T from b to a. In either case, a is called the *tail* of $[a, b]$, and b is called the *head* of $[a, b]$. The union of T and B will be denoted by E.

For notational convenience, we will frequently identify undirected graphs with their DFS representations. Since we are interested only in graphs with no isolated vertices, we will represent

graphs with their edge sets.

We define *successor*s for both vertices and edges. If $[a, b]$ is a tree edge, then b is a *successor* of a. If $[a, b]$ is a tree edge and $[b, c]$ is any edge, then $[b, c]$ is a *successor* of $[a, b]$. Back edges have no successors. We also define descendants and ancestors for both vertices and edges. A *descendant* of vertex (resp. edge) x is defined recursively as either x itself or a successor of a descendant of x. If y is a descendant of x, then x is an *ancestor* of y.

Let $e = [a, b] \in E$. Let Y be the set of vertices y such that for some x, $[x, y]$ is a back edge and also a descendant of e. If Y is not empty, we define $low_1(e)$ to be the smallest integer in Y, and $low_2(e)$ to be the second smallest integer in $Y \cup \{n+1\}$. Otherwise, we define $low_1(e) = low_2(e) = n+1$. The two mappings low_1 and low_2 can be computed in $O(m)$ time during the depth-first-search on G_0 [7]. If a is not the root of a DFS tree, and $low_1(e) \geq a$, then a is an articulation point of G [12].

If $e = [a, b]$ is any edge in E, then we define the function ϕ on E as follows.

$$\phi(e) = \begin{cases} 2\, low_1(e) & \text{if } low_2(e) \geq a \\ 2\, low_1(e) + 1 & \text{otherwise} \end{cases}$$

We arrange the successors of each tree edge in increasing order on their ϕ values. This ordering can be computed in $O(m)$ time using a bucket sort [7]. If $e_1, ..., e_k$ are the successors of e ordered this way, we will call e_i the ith successor of e for $i = 1 .. k$.

As in [7], for $e = [a, b]$, we define $S(e)$, the *segment* of e, to be the subgraph of G that consists of all the descendants of e. We use $ATT(e)$ to denote the set of back edges $[c, d]$ in $S(e)$ such that d is an ancestor of a, including a itself. Each back edge in $ATT(e)$ is called an *attachment* of e.

For any edge $e = [a, b]$, we define $cycle(e)$ as follows: if e is a back edge, then $cycle(e) = \{e\} \cup \{e'$: e' belongs to the tree path from b to $a\}$; if e is a tree edge and $low_1(e) > a$, then $cycle(e) = \{\}$; otherwise, $cycle(e) = cycle(e_1)$, where e_1 is the first successor of e. We use $sub(e)$ to denote the subgraph $S(e) \cup cycle(e)$. It is easy to see that if $cycle(e)$ is not empty, then the vertex $low_1(e)$ is always on $cycle(e)$. Also, if $low_1(e) \geq a$, then $sub(e) = S(e)$; if $low_1(e) < a$, then $sub(e) - S(e) = \{e'$: e' belongs to the tree path from $low_1(e)$ to $a\}$.

Fig. 6 illustrates some of these definitions, where $low_1(e) = 1$; $low_2(e) = 2$; $cycle(e) = \{$ [1, 2], [2, 3], [3, 4], [4, 5], [5, 6], [6, 7], [7, 8], [8, 1] $\}$; $S(e)$ contains all the edges in the graph except [1, 2], [2, 3], [3, 4]; $sub(e)$ is the whole graph; $ATT(e) = \{$[8, 1], [9, 3], [12, 1], [14, 2], [13, 4].

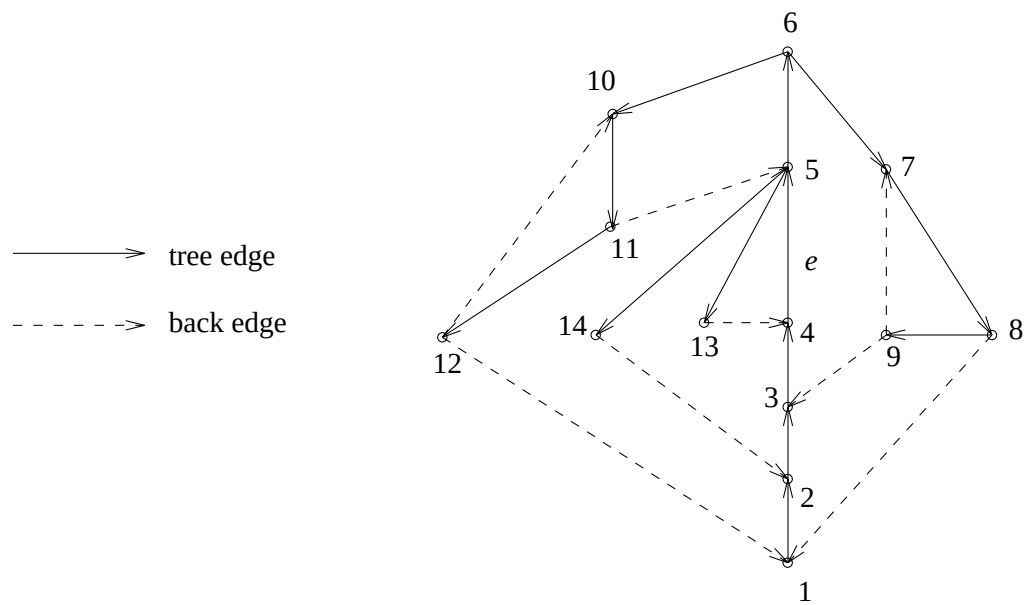


Fig. 6

3. Planarity testing

As explained in [7], a graph is planar if and only if each of its biconnected components is planar. Also, a graph of one edge is always planar. Thus, we need only consider how to test the planarity of biconnected graphs with more than one edge. Let $G = (V, T, B)$ be a DFS representation of such a graph. Then T forms a single tree with only one tree edge leaving the root. Call this tree edge e_0. Since $sub(e_0)$ is the whole graph, we can determine the planarity of G with a procedure that can determine the planarity of $sub(e)$ for all e in G.

We say that an edge e is planar if $sub(e)$ is planar. To determine the planarity of an edge e, we consider two cases. If e is a back edge, then $sub(e) = cycle(e)$, which is always planar. Otherwise, e is a tree edge having at least one successor. In this case we first determine the planarity of each of its successors. If all these successors are planar, then we determine the planarity of e based on the structure of its attachments. The details follow.

3.1. Structure of attachments

The planarity of an edge $e = [a, b]$ directly depends on the structure of its attachments. Since we assume that G is a biconnected graph with more than one edge, then $low_1(e) \le a$, and both $ATT(e)$ and $cycle(e)$ are not empty. If e is planar, then we can partition the edges of $ATT(e)$ into *block*s as follows. We put two back edges of $ATT(e)$ in the same block if they are on the same side of $cycle(e)$ in every embedding of $sub(e)$. Two blocks B_1 and B_2 of $ATT(e)$ *interlace* if they are on opposite sides of $cycle(e)$ in every embedding of $sub(e)$. Each block B_i of $ATT(e)$ can interlace at most one other block, since two attachments of e that cannot be

embedded on the same side of $cycle(e)$ as B_i must be in the same block.

The back edge on $cycle(e)$ is the only attachment of e that will not be embedded on either side of $cycle(e)$. By convention, this back edge forms a block by itself, called the *singular block* of e, which does not interlace other blocks of $ATT(e)$.

In Fig. 6, $ATT(e)$ consists of four blocks: B_1 = {[8, 1]}, B_2 = {[12, 1], [14, 2]}, B_3 = {[9, 3]}, and B_4 = {[13, 4]}. B_1 is singular. B_2 and B_3 are interlacing.

If $e' = [u, v]$ is an attachment of e, then $low_1(e) \le v \le a$. If $low_1(e) < v < a$, then we say that e' is *normal*. Otherwise we say that e' is *special*. A block of attachments of e is *normal* if it contains some normal attachment of e. Otherwise we say that it is *special*. In Fig. 6, B_1 and B_4 are special, and other blocks are all normal. We say that $sub(e)$ is *strongly planar* w.r.t. e if e is planar and if all the normal blocks of $ATT(e)$ can be embedded on the same side of $cycle(e)$. If $sub(e)$ is strongly planar (w.r.t. e), then we say that e is strongly planar. We have

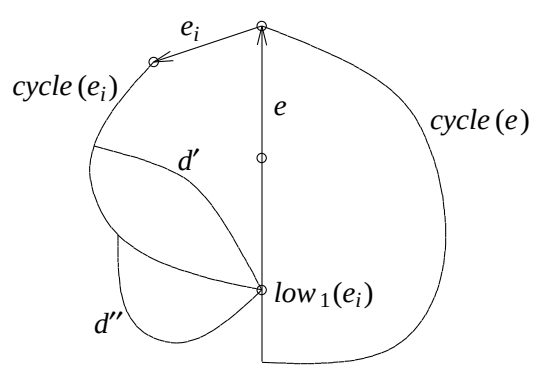
LEMMA 1. *Let e_i be the ith successor of e, where $i > 1$. Then e_i is strongly planar iff the subgraph $S(e_i) \cup cycle(e)$ is planar.*

Proof

$\Rightarrow$ If e_i is strongly planar, then there is an embedding of $sub(e_i)$ such that all its normal blocks are on the same side of $cycle(e_i)$. Thus we can add $cycle(e)$ to the other side of $cycle(e_i)$ to get an embedding of $S(e_i) \cup cycle(e)$.

$\Leftarrow$ If $S(e_i) \cup cycle(e)$ is planar, then in any embedding of $S(e_i) \cup cycle(e)$, all the normal blocks of $ATT(e_i)$ must be on the same side of $cycle(e_i)$. ■

Note that in an embedding of $S(e_i) \cup cycle(e)$, the special blocks of e_i do not have to be on the same side of $cycle(e_i)$. (See Fig. 7.)



The two special attachments d' and d'' of e_i can be on different sides of $cycle(e_i)$, although they are on the same side of $cycle(e)$.

Fig. 7

We will represent a block of back edges $H = \{[b_1, a_1], [b_2, a_2], ..., [b_t, a_t]\}$ by a list $K = [a_1, a_2, ..., a_t]$, where $a_1 \le a_2 \le ... \le a_t$. Frequently, we will identify blocks with their list representations. Define $first(H) = first(K) = a_1$, and $last(H) = last(K) = a_t$. If K is empty, we define $first(H) = first(K) = n+1$, and $last(H) = last(K) = 0$. We can further organize the blocks of $ATT(e)$ as follows: if two blocks X and Y interlace, we put them into a pair $[X, Y]$, assuming $last(X) \ge last(Y)$; if a nonempty block X does not interlace any other block, we form a pair $[X, [\ ]]$. Let $[X_1, Y_1]$ and $[X_2, Y_2]$ be two pairs of interlacing blocks. We say $[X_1, Y_1] \le [X_2, Y_2]$ iff $last(X_1) \le min(first(X_2), first(Y_2))$.

We say a list of interlacing pairs $[q_1,...,q_s]$ is *well-ordered* if $q_1 \le \cdots \le q_s$. Empty lists or lists of one pair are well-ordered by convention. We will see that all the interlacing pairs of $ATT(e)$ can be organized into a well-ordered list $[p_1, ..., p_t]$. We call this list $att(e)$.

In Fig. 6, $att(e) = [p_1, p_2, p_3]$, where $p_1 = [[1], [\ ]]$, $p_2 = [[3], [1, 2]]$, and $p_3 = [[4], [\ ]]$.

3.2. Computing *att* (*e*)

Now we are ready to compute $att(e)$. The planarity of e will be decided at the same time.

Consider any edge $e = [a, b]$. If e is a back edge, then its only attachment is e itself. Therefore $att(e) = [[[b], [\ ]]]$. Otherwise, let $e_1, ..., e_k$ be the successors of e in increasing order by their ϕ values. We first recursively compute $att(e_i)$ for each successor e_i of e. Then we compute $att(e)$ in four steps:

Algorithm 1

Step 1 For $i = 1 .. k$, delete all occurrences of b appearing in blocks within $att(e_i)$. Because these occurrences appear together at the end of the blocks that are contained in the last pairs of $att(e_i)$ only, a simple list traversal suffices to delete all these occurrences in time $O(k + \textit{number of deletions})$. After this, initialize $att(e)$ to be $att(e_1)$.

Step 2. For $i = 2 .. k$, merge all the blocks of $att(e_i)$ into one intermediate block B_i. See Fig. 8.

According to Lemma 1, this step can only be done for a given value of i if the normal blocks of $att(e_i)$ do not interlace. (If a pair of normal blocks of $att(e_i)$ interlace, the graph is not planar, and the computation fails.) To merge the blocks for a given value of i, we traverse the list of pairs $att(e_i)$, concatenating blocks to form B_i. Initially, B_i is empty. To process pair $[X, Y]$, if X and Y are both normal, the computation fails, since the graph is not planar. Otherwise, we concatenate X and Y onto the end of B_i in order and continue. In case that Y is a special block, we know that Y has only attachments ending in $low_1(e_i)$, since b was deleted from all blocks of $att(e_i)$ in Step 1. Thus the correct ordering of attachments is maintained by this process. This

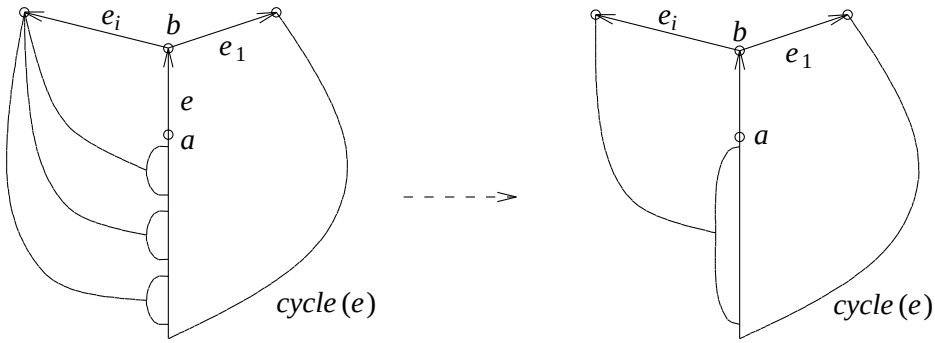


Fig. 8

step takes $O(1 + \textit{number of blocks in att}(e_i))$, resulting in one block for each i.

Step 3. Merge blocks in $att(e)$. See Fig. 9.

By Observation 3, all blocks D in $att(e)$ with $last(D) > low_1(e_2)$ must be merged into one block B_1. (If any two of these blocks interlace, the graph is not planar, and the computation fails.) This is achieved by merging from the high end of $att(e)$. The time is $O(1 + \textit{reduction in number of blocks})$. This step turns $att(e)$ into a list of pairs $p_1 \leq \cdots \leq p_h$ with only p_h possibly having a block D with $last(D) > low_1(e_2)$. Note that $low_1(e_2)$ is the lowest among the vertices $low_1(e_2)$, ..., $low_1(e_k)$.

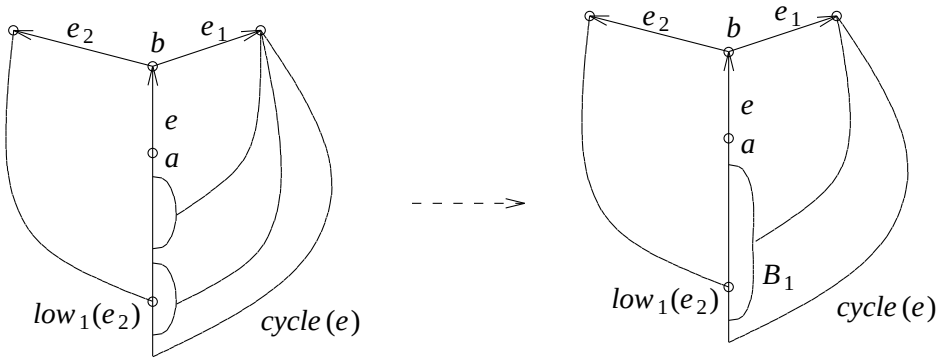


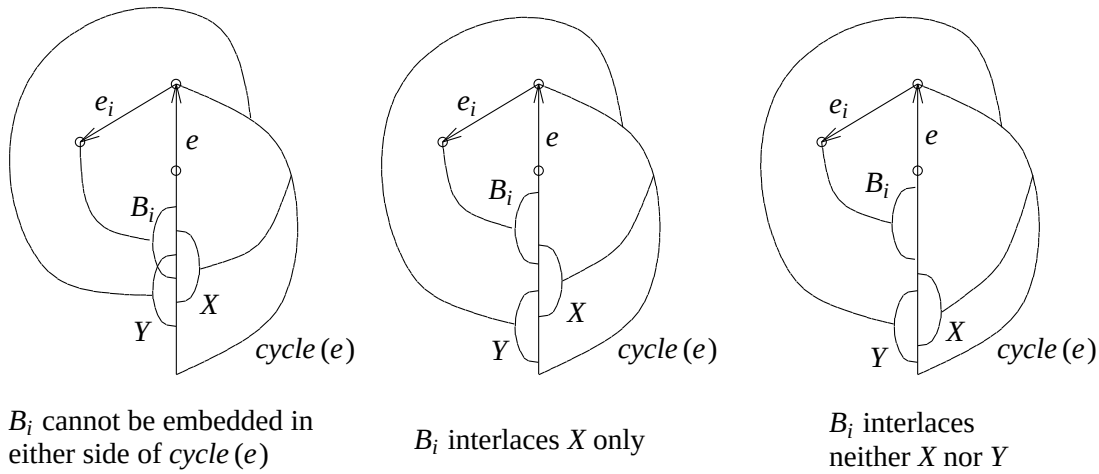
Fig. 9

Step 4. For $i = 2, ..., k$, add blocks B_i into $att(e)$.

To process B_i, consider the highest pair $P : [\ X,\ Y\ ]$ of $att(e)$. Consider three subcases:

i. If B_i cannot be embedded on either side of $cycle(e)$, then the computation of $att(e)$ fails.

ii. If B_i interlaces X only, then merge B_i into Y by concatenating their ordered list representations. Next, switch X and Y if $last(X) < last(Y)$.



B_i cannot be embedded in either side of $cycle(e)$

B_i interlaces X only

B_i interlaces neither X nor Y

Fig. 10

iii. If B_i interlaces neither X nor Y, then add $[B_i, [\ ]]$ to the high end of $att(e)$; $P := [B_i, [\ ]]$.

By the following lemma, testing whether B_i interlaces X or Y takes $O(1)$ time. Also by that lemma, it is not possible that B_i interlaces Y only, since $last(X) \geq last(Y)$ (see Fig. 10). ■

LEMMA 2. *In Step 4, B_i and D can be embedded on the same side of $cycle(e)$ iff $low_1(e_i) \geq last(D)$, where $D = X$ or $D = Y$.*

Proof

$\Rightarrow$ Assume $low_1(e_i) < last(D)$. Then there must be a path P_1 in $S(e_i)$ from b to $low_1(e_i)$ containing some back edge in B_i, and another path P_2 in $S(e_1) \cup ... \cup S(e_{i-1})$ from a vertex w on $cycle(e)$ to $last(D)$ containing some back edge in D but no edge on $cycle(e)$. We consider two cases (see Fig. 11). If $w > b$, then by Observation 3, P_1 and P_2 cannot be embedded on the same side of $cycle(e)$. If $w = b$, then the first edge on P_2 is e_j for some $1 < j < i$, which implies $i > 2$ and $\phi(e_j) \leq \phi(e_i)$. Consequently, $low_1(e_j) \leq low_1(e_i) < last(D) < b$, which implies that $low_2(e_j) < b$. If $low_1(e_j) < low_1(e_i)$, then there must be an undirected simple path P_3 between $last(D)$ and $low_1(e_j)$ containing back edges in D but no edges on $cycle(e)$. By observation 3 again, P_1 and P_3 cannot be embedded on the same side of $cycle(e)$. If $low_1(e_j) = low_1(e_i)$, then $low_2(e_i) < b$ (recall $low_1(e_j) \leq low_1(e_i)$). By observation 4, $S(e_i)$ and $S(e_j)$ cannot be embedded on the same side of $cycle(e)$. All of the above cases imply that B_i and D cannot be embedded on the same side of $cycle(e)$.

$\Leftarrow$ See the proof of Lemma 4 in the next section. ■

THEOREM 1.

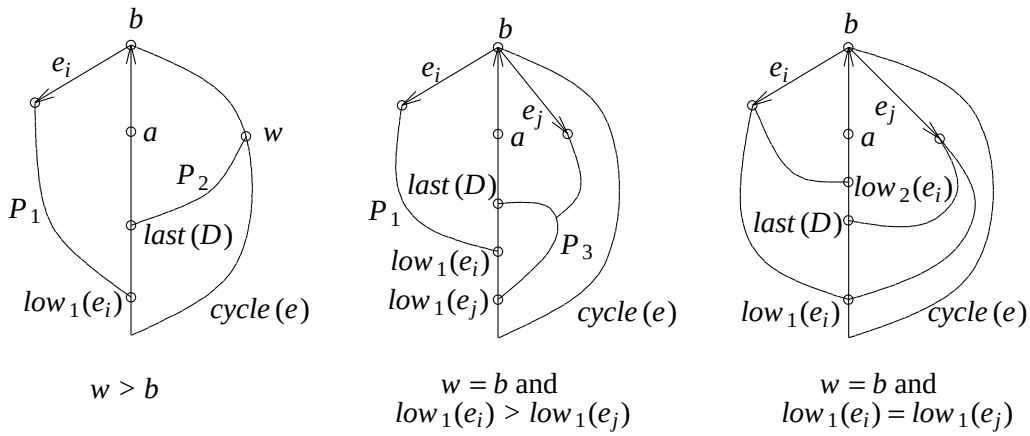


Fig. 11

1. Algorithm 1 computes $att(e)$ successfully iff e is planar.

2. If e is planar, then Algorithm 1 computes $att(e)$ correctly.

Proof See the next section. ■

3.3. Correctness

In the following proofs, unless stated otherwise, we will use $att(e)$ to mean the list $att(e)$ computed by Algorithm 1. But we will prove that this $att(e)$ correctly implements the $att(e)$ defined in Section 3.1.

During the presentation of Algorithm 1, we explained that two nonempty blocks form a pair within $att(e)$ only if they cannot be embedded on the same side of $cycle(e)$, and the computation of $att(e)$ fails only when e is not planar. Also we can see that the singular block of e is not merged with any other block. To prove Theorem 1, we still have to show that the following assertions are true:

(1) if computation of $att(e)$ succeeds, then e is planar;

(2) if any two nonempty nonsingular blocks of $att(e)$ do not form a pair, then these blocks can be embedded on the same side as well as on different sides of $cycle(e)$;

(3) $att(e)$ is well-ordered.

We first prove (3), *i.e.*,

LEMMA 3. *The list of pairs att(e) computed by Algorithm 1 is well-ordered.*

Proof We prove this lemma by induction on the number of descendants of e. If e has no successor, then e is a back edge, and the lemma is trivially true. Now assume that e is a tree edge with successors e_1, ..., e_k in increasing order by ϕ value, and that $att(e_1)$, ..., $att(e_k)$ are all well-ordered. After Steps 1 and 2 are executed, $att(e_1)$, ..., $att(e_k)$ are still well-ordered. Thus, $att(e)$

is well-ordered when it is initialized to $att(e_1)$. In Step 3, only blocks in the highest pairs of $att(e)$ are merged, and therefore $att(e)$ is still well-ordered after the merge. Then consider the moment in Step 4 just before B_i is added to $att(e)$. Assume $att(e)$ is well-ordered at this moment. Let $P: [X, Y]$ be the last pair of $att(e)$. We need only consider the two cases in which the computation does not fail.

1. $last(Y) \le low_1(B_i) < last(X)$. Then B_i is merged with Y. If P is the only pair in $att(e)$, then $att(e)$ is well-ordered by definition. Otherwise, let $Q: [X_1, Y_1]$ be the pair next to P. Then we have $Q \le P$ before merge. We need only to show that this is still true after the merge. If $i = 2$, then Step 3 guarantees that $first(B_2) = low_1(e_2) \ge \max(last(X_1), last(Y_1))$. If $i > 2$, then B_{i-1} is contained in either X or Y. Since $first(B_i) = low_1(e_i) \ge low_1(e_{i-1}) = first(B_{i-1}) \ge \min(first(X), first(Y))$, then merging B_i into Y does not change the value of $\min(first(X), first(Y))$. Thus, after merging Y and B_i, we still have $Q \le P$.

2. $low_1(B_i) \ge last(X)$. Then $[B_i, [\ ]]$ becomes the last pair of $att(e)$. Since $last(X) \le low_1(e_i) = first(B_i)$ in this case, we have $P \le [B_i, [\ ]]$.

Thus, $att(e)$ is still well-ordered after each B_i is added, $i = 2 .. k$. Therefore $att(e)$ is well-ordered after Step 4. ■

Next we prove the *if* part of Lemma 2: In Step 4 of Algorithm 1, if $low_1(e_i) \ge last(D)$, then B_i and D can be embedded on the same side of $cycle(e)$, where $D = X$ or $D = Y$.

Proof Consider an embedding of $sub(e)$ before B_i is added. Assume without loss of generality that X is embedded on the left hand side of $cycle(e)$ in this embedding. Let $h = last(X)$. Then $h = \max\{last(Z)$: Z is a block in $att_i\}$, and there is a face F on the left of $cycle(e)$ in the current embedding of $sub(e)$ such that the tree path from h to b is on the boundary of F. See Fig. 12.

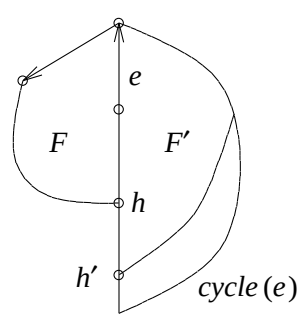


Fig. 12

Thus, if $low_1(B_i) \ge last(X)$, then B_i can be embedded in F. Similarly, let $h' = \max\{last(U)$: U is block in att_i embedded on the right hand side of $cycle(e)\}$. Then there is a face F' on the right of $cycle(e)$ in the current embedding of $sub(e)$ such that the tree path from h' to b is on the

boundary of F'. According to the proof of Lemma 3, if $low_1(B_i) \geq last(Y)$, then $low_1(B_i) \geq h'$ also. Therefore B_i can be embedded in F' in this case. ■

Then we prove the following lemma that implies the assertions (1) and (2). We say that a set W of blocks of $att(e)$ is *consistent* w.r.t. e if for all $X, Y \in W$, neither $[X, Y]$ nor $[Y, X]$ is in $att(e)$.

LEMMA 4.. *If Algorithm 1 does not fail, and D_1 and D_2 are two disjoint consistent sets of nonsingular blocks from $att(e)$, then there is an embedding of $sub(e)$ such that blocks of D_1 are on one side of $cycle(e)$ and blocks of D_2 are on the other side of $cycle(e)$.*

Proof The lemma is trivially true if e has no successors. If e has successors, let $e_1, ..., e_k$ be the list of successors of e in increasing order by their ϕ values. Assume that the lemma holds for each of these successors. We want to construct an embedding of $sub(e)$ such that D_1 and D_2 are embedded on different sides of $cycle(e)$.

If W is a set of blocks, then a *W-attachment* is an attachment contained in some block of W. For $j = 1, 2$, let $D_j' = \{ X : [X, Y]$ or $[Y, X]$ is a pair in $att(e)$ and $Y \in D_j\}$. Let $H_1 = \{ X$: ($[X, Y]$ is a pair in $att(e)$) and (X and Y are not in $D_1 \cup D_2$) }, and let $H_2 = \{ Y$: ($[X, Y]$ is a pair in $att(e)$) and (X and Y are not in $D_1 \cup D_2$) }. Let $C_1 = D_1 \cup D_2' \cup H_1$, and $C_2 = D_2 \cup D_1' \cup H_2$. For $j = 1, 2$, let $K_j = \{X : X$ is a block in $sub(e_1)$ containing some C_j-attachment$\}$. Then K_1 and K_2 are two disjoint consistent subsets of blocks of $sub(e_1)$.

Initially, we construct an embedding of $sub(e_1)$ such that K_1 and K_2 are on different sides of $cycle(e_1)$. As a result, those C_1-attachments and C_2-attachments contained in $sub(e_1)$ are on different sides of $cycle(e)$ (which is $cycle(e_1)$). This embedding exists by the induction hypothesis. Take this embedding to be the initial embedding of $att(e)$. Then for $i = 2, ... , k$, we add $sub(e_i)$ to this embedding one by one as follows.

Since the normal blocks of $att(e_i)$ do not interlace, we can, by induction, find an embedding of $sub(e_i)$ such that all of its normal blocks are embedded on the same side of $cycle(e_i)$. We call this embedding E_i, and its mirror image E_i'. Let B_i, P and $[X, Y]$ be the same as in Step 4 of algorithm 1. Assume without loss of generality that X is embedded on the left hand side of $cycle(e)$. Consider two cases:

Case 1: $last(Y) \leq low_1(e_i) < last(X)$. Then B_i is merged with Y. According to Lemma 2, one of E_i or E_i' can be embedded on the right of $cycle(e)$. If B_i contains any C_1-attachment, then X contains some C_2-attachment; if B_i contains any C_2-attachment, then X contains some C_1-attachment. In any case, C_1-attachments and C_2-attachments are still on different sides of $cycle(e)$ after B_i is embedded.

Case 2: $low_1(e_i) \geq last(X)$. Again by to Lemma 2, one of E_i or E_i', say E_i, can be embedded on the left of $cycle(e)$, and the other, E_i', can be embedded on the right of $cycle(e)$. One of

these two choices will result in an embedding such that C_1-attachments are on one side of $cycle(e)$, and C_2-attachments are on the other side.

When all the B_i's are added, we get an embedding of $sub(e)$ such that D_1 is on one side of $cycle(e)$, and D_2 is on the other side of $cycle(e)$. This is true because every D_1-attachment is a C_1-attachment, and every D_2-attachment is a C_2-attachment. ■

This completes the proof of Theorem 1, and establishes that the list $att(e)$ computed by Algorithm 1 has the properties discussed at the end of Section 3.1.

Let e, e_i, B_i, att_i, X, Y, and h' be the same as in the above proofs. Let $L = \{Z: [Z, U]$ is a pair in $att_i\}$, $R = \{U: [Z, U]$ is a pair in $att_i\}$, and $h_2 = \{last(U): U$ is a block in $R\}$. Then it is easy to see that $h' \geq h_2$. According to Lemma 4, there exists a embedding of $sub(e)$ (before adding B_i) such that L is embedded in one side of $cycle(e)$ and R is embedded on the other side. According to the proof of Lemma 2, $low_1(e_i) \geq last(Y)$ iff $low_1(e_i) \geq h_2$. Therefore we also have

COROLLARY 1.1. *B_i cannot be embedded in either side of $cycle(e)$ iff $low_1(e_i) < h_2$.* ■

Corollary 1.1 gives a test whether B_i can be added to att_i without referring to the top pair of att_i. This is useful in our maximal planar subgraph algorithm, where we need to test whether B_2 can be added to $att(e)$ before Step 3 is performed.

3.4. Data structure and running time

As suggested in [7], we can implement blocks as linked lists. An interlacing pair of blocks can be represented as a record containing two pointers to the two linked lists representing these two blocks. Then $att(e)$ can be represented as a linked list of such records. In this way, the time cost for Step 1 is $O(k + number\ of\ deletions)$. The cost for Step 2, 3 and 4 is $O(k + reduction\ in\ number\ of\ blocks)$. The cumulate expense of executing these steps over the whole graph is $O(m)$. The initial DFS in which low_1 values are computed takes time $O(m)$. Arranging the successors in increasing order by ϕ value for all tree edges takes $O(m)$ time using a bucket sort. Thus the whole algorithm runs in $O(m)$ time. It is well known that any $O(m)$-time algorithm for planarity testing can be implemented in $O(n)$ time since $m = O(n)$ for a planar graph [7].

3.5. A modification to Algorithm 1

Consider Step 4 of Algorithm 1. Lemma 2 requires that the successors of each tree edge be ordered by ϕ values. Maintaining this ordering causes difficulties in solving the maximal planar subgraph problem. Fortunately, we can modify Algorithm 1 so that it requires only the low_1 ordering of the successors of each tree edge.

Let $e = [a, b]$ be a tree edge, and $e_1, ..., e_k$ be the list of its successors in increasing order by low_1 values. Still define $cycle(e) = cycle(e_1)$. Then Step 1, 2, and 3 can be performed w.r.t. this ordering without any modification.

Next we want to merge $B_2, ..., B_k$ into $att(e)$ in that order. In general, successors ordered by low_1 values may not be ordered by ϕ values. Consequently, there may be some $1 < i \le k$ such that $\phi(e_{i-1}) > \phi(e_i)$. But if this happens, we know that $low_1(e_{i-1}) = low_1(e_i)$ and $low_2(e_i) \ge b$. If $i = 2$, Lemma 2 still applies, and we can merge B_2 into $att(e)$ as before. Otherwise, the following lemma says that we do not have to merge B_i into $att(e)$ at all:

LEMMA 5. *If for some* $2 < i \le k$, $low_1(e_{i-1}) = low_1(e_i)$, $low_2(e_i) \ge b$, *and* e_i *is planar, then* G *is planar iff* $G - S(e_i)$ *is planar.*

Proof The *only if* part is trivial, so we just prove the *if* part. Consider an embedding E_i of $G_i = G - S(e_i)$. Under the condition of the lemma, e_i has no normal attachments. Since e_i is planar, then e_i is strongly planar. Also, b and $low_1(e_i)$ are the only two vertices shared by $S(e_i)$ and G_i. Therefore $S(e_i)$ can be embedded in any face of E_i whose boundary contains the two vertices b and $low_1(e_i)$.

Let P be the tree path $cycle(e_{i-1}) \cap cycle(e)$ and let C be the closed curve $cycle(e_{i-1}) \cup cycle(e) - P$. Then C contains edges from both $S(e_{i-1})$ and $G_i - S(e_{i-1})$. By Observation 2, P is on one side of C. Call this side of C S_1, and the other side S_2. Let U be the set of faces in S_2 whose boundaries contain edges from $S(e_{i-1})$ only, and let W be the set of faces in S_2 whose boundaries contain edges from $G_i - S(e_{i-1})$ only. Then faces in U and faces in W do not share common boundaries. Thus, within S_2 there must be some face F whose boundary contains edges from both $S(e_{i-1})$ and $G_i - S(e_{i-1})$, and therefore contains at least two vertices common to $S(e_{i-1})$ and $G_i - S(e_{i-1})$. But all the vertices common to $S(e_{i-1})$ and $G_i - S(e_{i-1})$ are on P, and among them only b and $low_1(e_{i-1})$ are on the boundary of S_2. Therefore these two vertices must be on the boundary of F. Thus we can embed $S(e_i)$ in F to get an embedding of G. ■

Therefore, under the conditions of Lemma 5, in deciding the planarity of G, we can ignore its subgraph $S(e_i)$. Since the condition $low_1(e_{i-1}) = low_1(e_i)$ and $low_2(e_i) \ge b$ is implied by $low_1(e_{i-1}) \le low_1(e_i)$ and $\phi(e_{i-1}) > \phi(e_i)$, we can modify Step 4 as follows:

Step 4'. Add blocks $B_2, ..., B_k$ into $att(e)$ in that order, assuming $low_1(e_1) \le low_1(e_2) \le \cdots \le low_1(e_k)$. Initially, let $j = 1$ and $i = 2$. To process B_i, we consider two cases. If $j = 1$ or $\phi(e_j) \le \phi(e_i)$, we do the same thing as in Step 4, and then let $j = i$; otherwise, we do nothing.

The list $att(e)$ computed by the modified algorithm may not contain all the attachments of e. Some attachments may be omitted by Step 4', because their existence does not affect the planarity of the whole graph G.

4. The maximal planar subgraph problem

Now we consider the maximal planar subgraph problem: find a minimal set of edges whose deletion results in a planar graph. The resulting graph is called a *maximal planar subgraph* of G. We can always find a maximal planar subgraph of G by deleting back edges only, since all the tree edges form a forest, which is planar.

We will not assume that the input graph is biconnected, since deletion of back edges may turn a biconnected graph into a graph with articulation points. But without loss of generality we can assume that the input graph is connected. Thus the tree edges of G form a single tree with root r. Let $t_1, ..., t_s$ be the tree edges leaving the root. If $s = 1$, then $sub(t_1)$ is the whole graph G. If $s > 1$, then r is the only vertex common to $sub(t_1), ..., sub(t_s)$. Thus, to find a maximal planar subgraph of G, we can just find a maximal planar subgraph for each of the subgraphs $sub(t_1), ..., sub(t_s)$, and then simply put these subgraphs together. Therefore, what we need is a procedure that can find a maximal planar subgraph of $sub(e)$ for any given edge e of G.

4.1. Maximal *l–planar* subgraphs

We cannot build a maximal planar subgraph of $sub(e)$ by simply putting together the maximal planar subgraphs of $sub(e_1), ..., sub(e_k)$, and deleting those back edges causing failure in Algorithm 1. The reason is that after these edges are deleted, it may turn out that some other edges, which we deleted for making $sub(e_1), ..., sub(e_k)$ planar, would not have had to be deleted at all. We avoid this difficult situation by constructing such maximal subgraphs $S_1, ..., S_k$ of $sub(e_1), ..., sub(e_k)$ that they can be used to construct a planar subgraph S of $sub(e)$ without further deletion of edges. Two measures are taken for this purpose. Firstly, those back edges in $sub(e_i)$ that can cause failure in Step 3 or Step 4 of Algorithm 1 are deleted before a maximal subgraph of S_i is recursively computed. Secondly, the information where blocks of $sub(e_i)$ are allowed to interlace is passed to the recursive call that computes S_i, so that when the returned S_i is merged to $sub(e)$, Step 2 of Algorithm 1 can also be performed successfully without deletion. Since the planar subgraph S of $sub(e)$ computed by our algorithm may be used to build a larger planar subgraph of G in the same way as we use $S_1, S_2, ..., S_k$ to build S, we also need to know where in S blocks are allowed to interlace. This approach leads naturally to the concept of *l–planar subgraph*s, which is a generalization of the concept of strongly planar subgraphs.

Consider an edge $e = [a, b]$ and a vertex l on the tree path from $low_1(e)$ to a. An attachment $[u, v]$ of e is *l–normal* if $low_1(e) < v < l$. A block of attachments is *l–normal* if it contains some l-normal attachment. Let D be the list representation of a nonempty block of attachments. Define $second(D)$ to be the second smallest element in the set $\{x: x \in D\} \cup \{n+1\}$, and define $second([]) = n+1$. Then D is l-normal iff $low_1(e) < first(D) < l$ or $second(D) < l$. The two mappings *first* and *second* can be maintained during the computation of $att(e)$ in $O(1)$ time

for each modification to $att(e)$.

We say that the subgraph $sub(e)$ is l*–planar* if e is planar and the l-normal blocks of $att(e)$ do not interlace. (*c.f.* Fig. 13, where (a) is l-planar, but (b) is not.) Edge e is l*–planar* if $sub(e)$ is l-planar.

If H is a subgraph obtained from $sub(e)$ by deleting back edges only, then we can define the l-planarity for H (w.r.t. e) in the same way as we did for $sub(e)$. We will talk about l-planar subgraphs of $sub(e)$ in this sense. An l-planar subgraph of $sub(e)$ is *maximal* if it can be obtained from $sub(e)$ by deleting a minimal set of back edges.

Consider edge $e = [a, b]$. According to our definition, e is planar iff e is $low_1(e)$-planar, and e is strongly planar iff e is a-planar. Therefore, if we can find a maximal l-planar subgraph of $sub(e)$ for any l with $low_1(e) \le l \le a$, then we can compute a maximal planar subgraph of $sub(e)$.

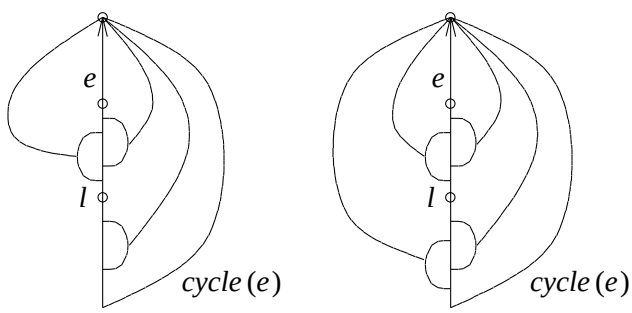


Fig. 13

The following is an outline of our maximal l-planar subgraph algorithm, where l is a given integer with $low_1(e) \le l \le a$ and remains fixed during the processing of an edge. Let $e = [a, b]$, and consider three cases:

Case 1: e is a back edge. Assign [[[b], [ ]]] to $att(e)$, and return.

Case 2: e is a tree edge with no successors. Assign [ ] to $att(e)$, and return.

Case 3: e is a tree edge with successors $e_1, ..., e_k$, among which e_1 has the smallest low_1 value. We construct a sequence $G_1, ..., G_k$ of l-planar subgraphs of $sub(e)$ such that G_1 is a maximal l-planar subgraph of $sub(e_1)$ and $low_1(e_1)$ remains unchanged; G_k is a maximal l-planar subgraph of $sub(e)$; and each G_i, $1 < i \le k$, is obtained from G_{i-1} by adding to it a strongly planar subgraph S_i of $sub(e_i)$, where e_i is some successor of e not contained in G_{i-1}. During the construction, we compute $att(e)$ using the modified version of Algorithm 1. We describe below in rough terms how we compute S_i:

1 select an edge e_i with the smallest low_1 value from successors of e not contained in G_{i-1};

2 **while** there exists a maximal strongly planar subgraph of $sub(e_i)$ whose addition to G_{i-1} destroys its l-planarity **do**

3 delete some attachments from $sub(e_i)$;

4 **if** the deletion changes the low_1 value of e_i **then**

5 select a possibly new edge e_i with the smallest low_1 value from successors of e not contained in G_{i-1};

6 **end if**;

7 **od**;

8 recursively construct a maximal strongly planar subgraph of $sub(e_i)$ without changing $low_1(e_i)$ further. We take this subgraph as S_i.

In the procedure sketched above, lines 1, 4, 5, 6, and 8 guarantee that subgraphs S_i are generated in increasing order by new low_1 values of the corresponding successors. For each $1 < i \leq k$, once S_i is computed, no edges will be deleted further from it. There are still two questions remaining to be answered: how the testing in line 2 can be done without constructing a maximal strongly planar subgraph of $sub(e_i)$, and how the attachments are chosen so that the deletion in line 3 makes the set of deleted edges minimal. These two questions are closely related and will be explained together in the next section.

Remark In Algorithm 1, we do not need the concept of l-planarity, since our purpose is to check the planarity of G. If some interlacing blocks of $sub(e)$ are found not being able to fit in the whole graph after returning from several levels of recursive calls, we simply declare that the graph is not planar. But if we want to construct a maximal planar subgraph of G, then it is too late to delete edges efficiently by that time. Therefore we use the parameter l to pass the information where blocks of $sub(e)$ are allowed to interlace, to the recursive calls, so that the correct edges are already deleted during the processing of $sub(e)$.

The need to generalize to l-planarity arises in the following way in the algorithm sketched above. To compute a maximal planar subgraph of the input graph, the recursive calls that construct S_i for $i = 2, ..., k$ must construct maximal strongly planar graphs. Within one of these recursive calls, the initial second-level recursive call (to construct a maximal b-planar subgraph of $sub(e_{i1})$, where e_{i1} is the first successor of edge e_i) and more deeply nested recursive calls of the same kind construct maximal l-planar subgraphs for general values of l. ■

4.2. Algorithm for deleting back edges

Let $e = [a, b]$, and consider the **while**-loop in the procedure sketched above. If $low_1(e_i) \geq b$, then b is the only vertex common to $sub(e_i)$ and $G - sub(e_i)$. In this case, we can apply the maximal planar subgraph to $sub(e_i)$ separately, and do not have to consider the effect on the

whole graph. Next, we consider the case when $low_1(e_i) < b$. Assume that $sub(e_i)$ is made strongly planar by deleting some back edges. Suppose that the low_1 value and the low_2 value of e_i are not changed by these deletions. We want to see whether the union of $sub(e_i)$ and G_{i-1} is l-planar.

As in planarity testing, let B_i be the block of attachments obtained by merging $att(e_i)$; let att_i be the current value of $att(e)$; let B_j be the last block merged into att_i by Step 4'; let h_1 = max{$last(Z)$: $[Z, U]$ is a pair in att_i} and h_2 = max{$last(U)$: $[Z, U]$ is a pair in att_i}. (Initially, we set j =1 and $att(e) = att(e_1)$ after removing all the occurrences of b from $att(e_1)$.) Finally, let h_3 = max{$last(Z)$: Z is an l-normal block of att_i}.

The two variables h_1, h_2 can be maintained in $O(1)$ time per modification to $att(e)$ by maintaining two lists L and R (as suggested in [7]), where L is the ordered list of nonempty blocks X such that $[X, Y]$ is a pair in $att(e)$, and R is the ordered list of nonempty blocks Y such that $[X, Y]$ is a pair in $att(e)$. If B_L and B_R are the highest blocks of L and R respectively, then $h_1 = last(B_L)$ and $h_2 = last(B_R)$. Lists L and R also let h_3 be maintained easily. If e is a back edge, $h_3 = 0$ from its definition. If e is a tree edge, we get the initial value of h_3 from the computation of $att(e_1)$, and modify it in $O(1)$ time for each modification of $att(e)$. The details will not be discussed here.

By Lemma 5, in case that e_i is strongly planar, $sub(e_i)$ can affect the planarity of G only if any of the following conditions holds:

a. $j = 1$, *i.e.*, no block B_j has been merged into $att(e)$ yet,

b. $low_1(e_j) < low_1(e_i)$, or

c. $low_2(e_i) < b$.

If any of these conditions is true, we consider two additional cases:

1. The union of $sub(e_i)$ and G_{i-1} is not planar. By Corollary 1.1, this happens iff $low_1(e_i) < h_2$.

2. The union of $sub(e_i)$ and G_{i-1} is planar, but not l-planar. Then B_i is l-normal, and it interlaces an l-normal block of att_i. We know that B_i is l-normal iff $low_1(e_1) < low_1(e_i) < l$ or $low_2(e_i) < l$. Also, it is easy to see that B_i interlaces an l-normal block of att_i iff $low_1(e_i) < h_3$.

This means, under the conditions a, b, or c, that the union of $sub(e_i)$ and G_{i-1} is not l-planar iff any of the following conditions holds

i. $low_1(e_i) < h_2$

ii. $low_1(e_1) < low_1(e_i) < \min\{h_3, l\}$

iii. $low_2(e_i) < l$ and $low_1(e_i) < h_3$

Therefore, if any of the condition i, ii, and iii holds, some back edge has to be deleted to change

either the low_1 value or the low_2 value of e_i. Such testing and deletion can be done even before making $sub(e_i)$ strongly planar. For this purpose, we combine the above conditions (a or b or c) and (i or ii or iii) into two groups according to whether they involve $low_2(e_i)$ or not:

Condition AA.

($j = 1$ and $low_1(e_i) < h_2$) or

($low_1(e_j) < low_1(e_i) < h_2$) or

($low_1(e_j) < low_1(e_i) < \min\{h_3, l\}$)

Condition BB.

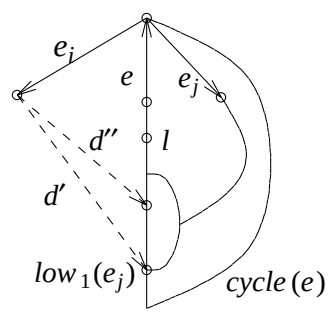
($low_2(e_i) < b$ and $low_1(e_i) < h_2$) or

($low_2(e_i) < b$ and $low_1(e_1) < low_1(e_i) < \min\{h_3, l\}$) or

($low_2(e_i) < l$ and $low_1(e_i) < h_3$)

It can be checked that the condition ((a or b or c) and (i or ii or iii)) is equivalent to the condition (AA or BB).

If Condition AA is true, we can make it false only by changing the value $low_1(e_i)$. In this case, we delete all the back edges of $sub(e_i)$ entering the vertex $low_1(e_i)$. After the deletion, we choose a possibly new e_i with the smallest low_1 value.



The edge e_i satisfies condition BB. If we choose to delete d', then d'' will also be deleted later because of Condition AA, and the resulting graph will not be maximal.

Fig. 14

If Condition AA is false, then we test Condition BB. If the result is true, we know that $low_1(e_i) = low_1(e_j)$. This is because $low_1(e_i) < low_2(e_i)$, which means that BB implies that $low_1(e_i) < h_2$ or $low_1(e_i) < \min\{h_3, l\}$, from which it follows that $low_1(e_j) = low_1(e_i)$; otherwise AA would be true. (We have $low_1(e_j) \le low_1(e_i)$ by the ordering of the successors of e.) To make Condition BB false, we can change the value of either $low_1(e_i)$ or $low_2(e_i)$. If we choose to change $low_2(e_i)$ consistently, then at least one of the back edges $[u, v]$ of $sub(e_i)$ with $v = low_1(e_i)$ will survive. But if we choose to change $low_1(e_i)$, it may happen that all the attachments in $ATT(e_i) \cap ATT(e)$ are deleted eventually and that the resulting graph is not maximal.

Therefore, in this case we choose to delete all the back edges $[u, v]$ of $sub(e_i)$ with $v = low_2(e_i)$ (see Fig. 14).

We test and delete repeatedly as described above until we find an edge e_i that does not satisfy AA or BB. Then we can construct S_i recursively from $sub(e_i)$ and merge it into G_{i-1}. Since no edge is added to $sub(e_i)$ during the construction of S_i, conditions AA and BB remain false after the construction. Thus, the resulting graph G_i will be planar, and no l-normal blocks will interlace.

To see that the deleted set of back edges is minimal, let $[u, v]$ be an edge deleted by the above algorithm, and add it back to G_i. If $[u, v]$ was deleted because of Condition AA, then $low_1(e_i) = v$ now, and Condition AA is true again. If $[u, v]$ was deleted because of Condition BB, then $low_2(e_i) = v$ now, and Condition BB is true again. Notice that, in the latter case, the low_1 value of e_i has remained unchanged since the deletion of $[u, v]$. In either case, G_i will not be l-planar.

4.3. Data structures and running time

In the algorithm described above, we need to repeatedly select an unprocessed successor of e with the smallest low_1 value, and the low_1 values of tree edges are constantly changing. Therefore we maintain a heap [13] based on low_1 values of the unprocessed successors of the tree edge e currently being processed. Since the algorithm is recursive, we actually maintain simultaneously a heap of unprocessed successors for each tree edge along the path to the currently active tree edge. The total size of all such heaps is $O(m)$. The initialization of all these heaps takes a total of $O(m)$ time. When the low_1 value of some element in a heap increases, we modify the heap accordingly. It is important to note that any two edges in active heaps are unrelated; thus deletion of a single attachment can modify the low_1 value of only a single such edge. It follows that the total number of modifications to and deletions from heaps is $O(m)$. The time for the heap operations is $O(\log n)$ time per operation, for a total of $O(m \log n)$ time. (Since $m < n^2$, $\log m = O(\log n)$.)

We also need a data structure for the back edges of $sub(e)$ so that the following operations can be done efficiently:

1. delete an attachment $[u, v]$ of e with $v = low_1(e)$ or $v = low_2(e)$;

2. maintain the low_1 and low_2 values of e;

3. split the data structure into several pieces, one for each successor of e.

One easy solution that meets these requirements is the *selection tree* [8]. To represent a set of edges E_0 as a selection tree T_0, we store edges of E_0 inside the leaves of T_0 from left to right in increasing order (by DFS number) of their tails. Edges with the same tail are ordered

arbitrarily. Each internal node w of T_0 has two children *w.lchild* and *w.rchild*. Let S_w be the set of edges stored in the leaves of the subtree rooted at w; let $lb = \min\{x : [x, y] \in S_w\}$ and $rb = \max\{x : [x, y] \in S_w\}$; let $low_1 = \min\{y : [x, y] \in S_w\}$, and $low_2 = \min(\{y : [x, y] \in S_w \mid y \neq low_1\} \cup \{n+1\})$. Then the four values lb, rb, low_1 and low_2 are stored in the four fields $w.lb$, $w.rb$, $w.low_1$, and $w.low_2$ of w respectively. If w is a leaf storing the edge $[x, y]$, then $w.lb = x$, $w.rb = x$, $w.low_1 = y$, and $w.low_2 = n+1$. The values in each internal node can be computed from the values in the children (in constant time).

In the following discussion, we will refer to a tree by its root. Let r_1 and r_2 be two selection trees representing the two disjoint sets of edges E_1 and E_2. If $u_1 \le u_2$ for all $[u_1, v_1] \in E_1$ and $[u_2, v_2] \in E_2$, then we can merge r_1 and r_2 to get the selection tree for $E_1 \cup E_2$ in $O(1)$ time:

```
procedure merge(r1, r2);
begin  if r1 = null then
               return r2;
       end if;
       if r2 = null then
               return r1;
       end if;
       r := newnode();
       r.lchild := r1;
       r.rchild := r2;
       r.lb := r1.lb;
       r.rb := r2.rb;
       r.low1 := min{r1.low1, r2.low1};
       r.low2 := min({r1.low1, r2.low1, r1.low2, r2.low2} - {r.low1});
       return r;
end;
```

Let r be a selection tree representing a set of edges E_0. To split E_0 into two sets $E_1 = \{[u, v] \in E_0 \mid u \le u_x\}$ and $E_2 = \{[u, v] \in E_0 \mid u > u_x\}$, we split r with respect to u_x as follows:

```
procedure split(r , ux);
begin  if ux < r.lb then
               return [null, r];
       elseif  ux ≥ r.rb then
               return [r, null];
       else    [rl, rr] := [r.lchild, r.rchild];
               if ux < rl.rb then
                       [rl1, rl2] := split(rl, ux);
                       return [rl1, merge(rl2, rr)];
               else    [rr1, rr2] := split(rr, ux);
                       return [merge(rl, rr1), rr2];
               end if;
       end if;
end;
```

The height of any tree that results from splitting a tree r can be no greater than the height of r. To select and delete an edge $[x, v]$ from a tree r, where $v \in \{r.low_1, r.low_2\}$, we do the following:

```
procedure delete(r, v);
begin  if r is a leaf then
               mark the back edge stored in r as 'deleted';
               return null;
       else    [r_l, r_r] := [r.lchild, r.rchild];
               if v = r_l.low_1 or v = r_l.low_2 then
                      return merge(delete(r_l, v), r_r);
               else   return merge(r_l, delete(r_r, v));
               end if;
       end if;
end;
```

Assuming the input graph G is connected, we know that all the tree edges form a tree. Let the root be 1. For technical reasons, we add a dummy edge $e_0 = [0, 1]$ to the tree edges. To get a maximal planar subgraph of G, we just construct a 0-planar subgraph of $sub(e_0)$, and then delete e_0 from it. Initially, we construct a balanced selection tree $tree(e_0)$ to store all the back edges of G. The height of this tree is $O(\log n)$. The time and space needed to initialize $tree(e_0)$ are both $O(m)$.

When we begin to construct a maximal l-planar subgraph for a tree edge e, we first split $tree(e)$ into several pieces $tree(e_1)$, ..., $tree(e_k)$, where e_1, ..., e_k are the successors of e not marked as 'deleted'. For each such successor e_i, $tree(e_i)$ is a selection tree representing the set of back edges in $sub(e_i)$, and can be obtained as follows. If e_i is a back edge, then $tree(e_i)$ consists of a single edge, and can be constructed in $O(1)$ time. If e_i is a tree edge, let $e_i = [b, c_i]$, and let n_i be the number of descendants of c_i. It is well known that a back edge $[u, v]$ is a descendant of e_i iff $c_i \le u < c_i + n_i$. Then we can use the procedure *split* to get $tree(e_i)$ from $tree(e)$ in $O(logn)$ time. Since there are at most n tree edges in G, then the splitting takes takes $O(n\log n)$ time for the whole algorithm. After each split, the total size of the trees is still $O(m)$.

To select and delete an attachment $[x, v]$ of e_i, where $v \in \{low_1(e_i), low_2(e_i)\}$, we execute $delete(tree(e_i), v)$, which takes $O(log(n))$ time. There can be at most $O(m)$ such invocations of *delete*, so the total cost for executing *delete* is $O(m\log n)$. Given the selection tree $tree(e_i)$, the values $\phi(e_i)$, $low_1(e_i)$, and $low_2(e_i)$ can be computed from $tree(e_i)$ in $O(1)$ time: if $tree(e_i)$ is null, we just set these values to $n+1$; otherwise, they can be computed from $tree(e_i).low_1$ and $tree(e_i).low_2$. Thus, the total cost of selection tree operations is $O(m\log n)$.

We have mentioned that the total cost of heap operations is also $O(m\log n)$. The other costs of the algorithm are the same as in planarity-testing. Thus the total cost of our maximal planar subgraph algorithm is $O(m\log n)$.

4.4. The complete algorithm

Now we summarize our maximal planar subgraph algorithm. We take a connected undirected graph as input, and convert it into a DFS representation $G = (V, T, B)$. At the same time, we compute the two mappings *succ* and N, where, for each $e \in T$, $succ(e)$ gives the successor edges of e in increasing order of their tails, and for each $v \in V$, $N(v)$ gives the number of descendants of v. We assume that there is a dummy edge $e_0 = [0, 1]$ such that $succ(e_0)$ gives the list of tree edges leaving the root. The whole preprocessing takes $O(m)$ time.

We summarize the maximal l-planar subgraph algorithm below.

```
procedure lplanar(e, l);
begin  let e = [a, b];
       if e ∈ B then
              return [[[b], []]];
       end if;
       if e has no successors then
              return [];
       end if;
       let e_1, ..., e_k be the successors of e not marked as 'deleted';
       split tree(e) into tree(e_1), ..., tree(e_k);
       organize e_1, ..., e_k into a heap based on their low_1 values, with the smallest one on the top;
       let e_1 be the edge on the top of the heap;
       delete e_1 from the heap;
1      att(e) := lplanar(e_1, l);
       delete all the occurrences of b from the top blocks of att(e);
       j := 1;
       i := 2;
       while heap is not empty do
              let e_i be the edge on the top of the heap;
              if low_1(e_i) ≥ b then
                     delete e_i from the heap;
2                    dummy := lplanar(e_i, b);
              elseif Condition AA is true then
                     v := tree(e_i). low_1;
                     while v = tree(e_i). low_1 do
3                           tree(e_i) := delete(tree(e_i), v);
                     end while;
                     if e_i is a back edge then
                            delete e_i from heap;
                     else modify heap;
                     end if;
              elseif Condition BB is true then
                     v := tree(e_i). low_2;
                     while v = tree(e_i). low_2 do
4                           tree(e_i) := delete(tree(e_i), v);
                     end while;
              else   delete e_i from the heap;
5                    att(e_i) := lplanar(e_i, b);
```

```
6                    merge blocks of att(e_i) into one block B_i;
                     delete all the occurrences of b from the top blocks of att(e_i);
                     if i = 2 then
7                          perform Step 3 of Alg.1;
                     end if;
8                    merge B_i into att(e) as described in Step 4';
                     i := i + 1;
                end if
            end while;
            return att(e);
        end;
```

The procedure *lplanar* (e, l) implicitly constructs a maximal l-planar subgraph of $sub(e)$ by deleting a minimal set of back edges. The parameter l specifies where the blocks of $att(e)$ are not allowed to interlace in the resulting subgraph, so that this subgraph can be used to build a larger planar subgraph without further deletion when we process the predecessor of e. For the initial call where $e = e_0$, we have $l = 0$, meaning that we need to construct a maximal planar subgraph of $sub(e_0)$. In the recursive calls for the successors of e, the l values are determined as follows. Since no l-normal blocks are allowed to interlace in $sub(e)$, then no l-normal blocks are allowed to interlace in $sub(e_1)$ either. Thus the recursive call of *lplanar* for e_1 (line 1) has the same parameter l as for the edge e. The remaining calls for e_2, ..., e_k (line 5) just construct maximal strongly planar subgraphs, therefore have b as their l values. Thus when we merge blocks at line 6, no normal blocks of $att(e_i)$ interlace. At line 7, we merge all normal blocks of $att(e)$ above $low_1(e_2)$ in to one block; at line 8, we merge B_i into $sub(e)$. Because of the deletions at lines 3 and 4, these steps can be performed successfully (without any further deletion). At line 2, $sub(e_i)$ is detected to be a biconnected component and is processed separately.

To compute a maximal planar subgraph, we simply do the following:

1. Organize B into a selection tree $tree(e_0)$;

2. Execute $lplanar(e_0, 0)$;

Then $T \cup B - B'$ gives a maximal planar subgraph of G, where B' is the set of back edges deleted by the procedure *delete* in the preceding algorithm.

Remark The procedure *lplanar* can be greatly simplified if we know that all the tree edges of G are on a same cycle. In this case, the low_1 values need not be dynamically maintained: if $e = [a, b]$ is a back edge, then $low_1(e) = b$; otherwise $low_1(e) = 1$. The low_2 values, which are used only for testing condition BB when $i > 1$, need not be maintained either, since for $i > 1$, e_i is a back edge. As a result, the selection trees are no longer useful, and the heaps storing the successors of tree edges can be replaced by lists precomputed as in Algorithm 1. With these simplifications, our algorithm gives the following result which is first reported in [10] by *Kobayashi et at.*: A maximal planar subgraph can be constructed in linear time provided that a

Hamiltonian tour of the graph is given. ■

5. Summary

The problem of drawing graphs in the plane arises naturally in circuit layout. Since finding a maximum planar subgraph is *NP*-complete [5], a maximal planar subgraph seems to be a reasonable approximation. Because planarity-testing can be done in linear time, it is easy to solve the maximal planar subgraph problem in $O(mn)$ time: start with a graph H with no edge; for each edge of the input graph G, add it to H if the resulting graph is planar, and reject it otherwise. The resulting graph H will be a maximal planar subgraph of G. However, a better solution seemed to be hard to find for a long time. Jayakumar *et al.* [9] even made the conjecture that "no maximal planarization algorithm of complexity better than $O(mn)$ will be possible." Our $O(m\log n)$ solution disproves this conjecture, as does the method of Di Battista and Tamassia [3].

We have assumed that the input graph to our algorithm is connected. For a more general graph, we can find a maximal planar subgraph by applying our algorithm to each of its connected components.

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