# **Fast Minor Testing in Planar Graphs**

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**Abstract** Minor Containment is a fundamental problem in Algorithmic Graph Theory used as a subroutine in numerous graph algorithms. A model of a graph H in a graph G is a set of disjoint connected subgraphs of G indexed by the vertices of H, such that if  $\{u, v\}$  is an edge of H, then there is an edge of G between components  $C_u$  and  $C_v$ . A graph H is a minor of G if G contains a model of H as a subgraph. We give an algorithm that, given a planar n-vertex graph G and an G-vertex graph G-verte

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#### 1 Introduction

For two input graphs G and H, the MINOR CONTAINMENT problem is to decide whether H is a minor of G. This is a classical NP-complete problem [18], and remains NP-complete even when both graphs G and H are planar, as it is a generalization of the HAMILTONIAN CYCLE problem. When H is fixed, by Robertson and Seymour's celebrated result [30], there is an algorithm to decide if H is a minor of an input graph G that runs in time  $f(h) \cdot n^3$ , where n is the number of vertices of G, h is the number of vertices in H, and f is some recursive function. One of the significant algorithmic implications of this result is that, combined with Robertson and Seymour's Graph Minor Theorem [32], it shows the polynomial-time solvability of many graph problems, some of which were previously not even known to be decidable [17]. However, these algorithmic results are highly *non-practical*. This triggered an ongoing quest in the Theory of Algorithms since then—related to trying to simplify the 20-papers proof of the Graph Minors Theorem—for making Graph Minors constructive and for making its algorithmic proofs practical for a wide range of applications (e.g., [10, 23]).

Unfortunately, in Robertson and Seymour's minor testing algorithm [30], the function f(h) has an *immense* (super-)exponential growth, which makes the algorithm absolutely impractical even for very simple patterns (see [24] for recent theoretical improvements of this function). There were several attempts to improve the running time of Robertson and Seymour's algorithm. One direction of such improvements is decreasing the degree of the polynomial in n. For example, Reed and Li gave a linear time algorithm solving  $K_5$ -minor containment [29]. The second direction of improvements is towards reducing the exponential dependency in the function f(h), which is a natural direction of study for Parameterized Complexity [16]. A significant step in this direction was done by Hicks [21], who provided in graphs of branchwidth k and m edges an  $\mathcal{O}(3^{k^2} \cdot (h+k-1)! \cdot m)$  time algorithm, following the algorithm sketched by Robertson and Seymour [30]. Recently, this was improved to  $\mathcal{O}(2^{(2k+1)\log k} \cdot h^{2k} \cdot 2^{2h^2} \cdot m)$  on general graphs, and in planar, and more generally, in graphs of bounded genus, to  $2^{\mathcal{O}(k)} \cdot h^{2k} \cdot 2^{\mathcal{O}(h)} \cdot n$  [1].

In this paper we focus on the case where the input graph G is planar.

PLANAR H-MINOR CONTAINMENT

**Input:** A planar graph G.

**Objective:** Either find a model M of H in G, or conclude that

G does not contain such a model.

Over the last four decades, many different algorithmic techniques in planar graphs were developed for different type of problems and algorithms, including approximation [5, 9], exact [14, 26], and parameterized algorithms [3, 8, 15]. However, it seems that none of these approaches can be used to speed up the algorithm for PLANAR *H*-MINOR CONTAINMENT.



## 1.1 Our Results and Key Ideas

By arguments inspired by Bidimensionality Theory [7], it can be easily shown that the  $2^{\mathcal{O}(k)} \cdot h^{2k} \cdot 2^{\mathcal{O}(h)} \cdot n$  time algorithm from [1], combined with the grid minor Theorem of Robertson, Thomas, and Seymour [31], can be used to solve Planar H-Minor Containment in time  $\mathcal{O}(2^{\mathcal{O}(h\log h)} \cdot n + n^2 \cdot \log n)$ . This directly sets up the challenge of designing a single-exponential (on the size h of the pattern H) algorithm for this problem.

Our main result is the following theorem.

**Theorem 1** Given a planar graph G on n vertices and a graph H on h vertices, PLANAR H-MINOR CONTAINMENT is solvable in time  $\mathcal{O}(2^{\mathcal{O}(h)} \cdot n + n^2 \cdot \log n)$ .

That is, we prove that when G is planar the behavior of the function f(h) can be made *single-exponential*, improving over all previous results for this problem [1, 21, 30]. In addition, we can *enumerate* and *count* the number of models within the same time bounds. Let us remark that by Theorem 1, PLANAR H-MINOR CONTAINMENT is solvable in polynomial time when the size of the pattern graph H is  $\mathcal{O}(\log n)$ , therefore substantially improving the existing algorithms for small patterns [12].

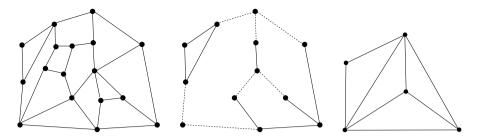
In order to prove Theorem 1, we introduce a novel approach of dynamic programming in planar graphs of bounded branchwidth, namely partially embedded dynamic programming. This approach is extremely helpful in computing graph minors, and we believe that this technique can be used in many related problems including PLANAR DISJOINT PATHS. Our technique is inspired by the technique of embedded dynamic programming introduced in [13] for solving PLANAR SUBGRAPH ISOMORPHISM for a pattern of size h and an input graph of size n in time  $2^{\mathcal{O}(h)} \cdot n$ . There, one controls the partial solutions by the ways the separators of G can be routed through the pattern. The difference (and difficulty) concerning PLANAR H-MINOR CONTAINMENT is that we look for a model M of size  $\mathcal{O}(n)$  out of  $2^{\mathcal{O}(n)}$  possible non-isomorphic models of H in G. In partially embedded dynamic programming, we look for potential models of H in G with a "magnifying glass" only at a given separator S of G. That is, we consider a collection A of graphs A arising from 'decontracting' a part of H, namely the part interacting with S. In this way, each graph A behaves like a subgraph of G inside the intersection with S, and outside that intersection A behaves like a minor of G; this is why we call our dynamic programming technique "partially embedded".

After giving some preliminaries in Sect. 2, we first show in Sect. 3 how PLANAR H-MINOR CONTAINMENT can be solved in polynomial time for input graphs of large branchwidth (in comparison to the pattern size). If the branchwidth is small, we compute the collection  $\mathcal{A}$  in Sect. 3.2 and give the partially embedded dynamic programming approach in Sect. 3.3.

## 2 Preliminaries

We use standard graph terminology, see for instance [11].





**Fig. 1** On the left hand side we illustrate a graph G. In the middle we have a subgraph M of G. Contracting the edges of M indicated by the dashed lines, we obtain minor H illustrated on the right hand side. M is a model of H in G. The dashed lines illustrate its c-edges, the solid lines its m-edges

## 2.1 Graphs and Graph Minors

All graphs considered in this article are simple and undirected. Given a graph G, we denote by V(G) and E(G) the vertex set and the edge set of G, respectively. A graph H is a *subgraph* of a graph G,  $H \subseteq G$ , if  $V(H) \subseteq V(G)$  and  $E(H) \subseteq E(G)$ . We define graph the operation of *contracting* an edge  $e = \{x, y\} \in G$  by removing e, x and y from G, and adding a new vertex  $v_e$  adjacent to both the former neighbors of x and the former neighbors of y (excluding x and y).

A graph H is a *minor* of a graph G (denoted by  $H \leq G$ ), if H can be obtained from a subgraph of G by a (possibly empty) sequence of edge contractions. In this case we also say that G is a *major* of H. A graph H is a *contraction minor* of a graph G (denoted by  $H \leq_c G$ ), if H can be obtained from G by a (possibly empty) sequence of edge contractions.

A model M of a minor H in G is a subgraph of G, where the edge set E(M) is partitioned into c-edges (contraction edges) and m-edges (minor edges) such that the graph resulting from contracting all c-edges is isomorphic to H.

For an illustration, see Fig. 1.

#### 2.2 Branchwidth

A branch decomposition  $(T, \mu)$  of a graph G consists of an unrooted ternary tree T (i.e., all internal vertices have degree three) and a bijection  $\mu: L \to E(G)$  from the set L of leaves of T to the edge set of G. We define for every edge e of T the middle set  $\operatorname{mid}(e) \subseteq V(G)$  as follows: Let  $T_1$  and  $T_2$  be the two connected components of  $T \setminus \{e\}$ . Then let  $G_i$  be the graph induced by the edge set  $\{\mu(f): f \in L \cap V(T_i)\}$  for  $i \in \{1, 2\}$ . The middle set is the intersection of the vertex sets of  $G_1$  and  $G_2$ , i.e.,  $\operatorname{mid}(e) = V(G_1) \cap V(G_2)$ . The width of  $(T, \mu)$  is the maximum order of the middle sets over all edges of T, i.e.,  $\operatorname{width}(T, \mu) := \max\{|\operatorname{mid}(e)|: e \in E(T)\}$ . The branchwidth of G is defined as  $\operatorname{bw}(G) := \min\{\operatorname{width}(T, \mu) \mid (T, \mu) \operatorname{branch} \operatorname{decomposition} \operatorname{of} G\}$ . Note that for each  $e \in E(T)$ ,  $\operatorname{mid}(e)$  is a separator of G, unless  $\operatorname{mid}(e) = \emptyset$ .

Remark 1 For every two edges  $e, f \in E(T)$  with  $e \cap f \neq \emptyset$ , we have  $|\text{mid}(e) \cup \text{mid}(f)| \leq 1.5 \cdot width(T, \mu)$ .



Intuitively, a graph G has small branchwidth if G is close to being a tree. The fundamental Grid Minor Theorem says that, roughly, a graph has either small branchwidth, or it contains a large grid as a minor. We use the variant for planar graphs.

**Proposition 1** [6, 20, 31] Given a planar graph G on n vertices with  $bw(G) \ge k$ , a model of the  $(\lfloor k/3 \rfloor \times \lfloor k/3 \rfloor)$ -grid in G can be found in time  $\mathcal{O}(n^2 \cdot \log n)$ .

On the other hand, every planar graph is a minor of a large enough grid.

**Proposition 2** [31] *If H is a planar graph with*  $|V(H)| + 2|E(H)| \le \ell$ , then H is isomorphic to a minor of the  $(2\ell \times 2\ell)$ -grid.

# 2.3 Planar Graphs and Equivalent Drawings

Let  $\Sigma$  be the unit sphere. A *planar drawing*  $\Phi$ , or simply *drawing*  $\Phi$ , of a graph G with vertex set V(G) and edge set E(G) maps vertices to points in the sphere, and edges to simple curves between their end-vertices, such that edges do not cross, except in common end-vertices. A *plane graph*  $\langle G, \Phi \rangle$  is a graph G together with a planar drawing G. A *planar graph* is a graph that admits a planar drawing. The set of *faces* G of a plane graph G is defined as the collection of the connected regions of G of a subgraph of a plane graph G, induced by the vertices and edges incident to a face G of a plane graph G of a planar graph G. A *homeomorphism* of G onto G is a homeomorphism of G onto itself which maps vertices, edges, and faces of G onto vertices, edges, and faces of G onto the other.

**Proposition 3** (e.g. [28]) Every planar n-vertex graph has  $2^{\mathcal{O}(n)}$  non-equivalent drawings. A set of all such drawings can be computed in time  $2^{\mathcal{O}(n)}$ .

**Proposition 4** [37] *The number of non-isomorphic edge-maximal planar graphs on* n *vertices is*  $2^{\mathcal{O}(n)}$ .

#### 2.4 Nooses and Combinatorial Nooses

A *noose* of a plane graph G is a simple closed curve in  $\Sigma$  that meets G only in vertices. From the Jordan Curve Theorem (e.g. [27]), it follows that nooses separate  $\Sigma$  into two regions.

Let  $V(N) = N \cap V(G)$  be the vertices and F(N) be the faces intersected by a noose N. The *length* of N is |V(N)|, the number of vertices in V(N). The clockwise order in which N meets the vertices of V(N) is a cyclic permutation  $\pi$  on the set V(N).

A combinatorial noose  $N_C = [v_0, f_0, v_1, f_1, \dots, f_{\ell-1}, v_\ell]$  in a plane graph G is an alternating sequence of vertices and faces of G, such that

<sup>&</sup>lt;sup>1</sup>If the drawing  $\Phi$  is clear from the context, we may also simply write *plane graph G*.



- $f_i$  is a face incident to both  $v_i$ ,  $v_{i+1}$  for all  $i < \ell$ ;
- $v_0 = v_\ell$  and the vertices  $v_1, \ldots, v_\ell$  are mutually distinct; and
- if  $f_i = f_j$  for any  $i \neq j$  and  $i, j = 0, ..., \ell 1$ , then the vertices  $v_i, v_{i+1}, v_j$ , and  $v_{j+1}$  do not appear in the order  $(v_i, v_j, v_{i+1}, v_{j+1})$  on the bound of face  $f_i = f_j$ .

The *length* of a combinatorial noose  $[v_0, f_0, v_1, f_1, \dots, f_{\ell-1}, v_\ell]$  is  $\ell$ .

Remark 2 The order in which a noose N intersects the faces F(N) and the vertices V(N) of a plane graph G gives a unique alternating face-vertex sequence of  $F(N) \cup V(N)$  which is a combinatorial noose  $N_C$ . Conversely, for every combinatorial noose  $N_C$  there exists a noose N with face-vertex sequence  $N_C$ .

We will refer to combinatorial nooses simply as nooses if it is clear from the context.

**Proposition 5** [13] Every plane n-vertex graph has  $2^{\mathcal{O}(n)}$  combinatorial nooses.

## 2.5 Sphere Cut Decompositions

For a plane graph G, we define a *sphere cut decomposition* (or *sc-decomposition*)  $\langle T, \mu, \pi \rangle$  as a branch decomposition, which for every edge e of T has a noose  $N_e$  that divides  $\Sigma$  into two regions  $\Delta_1$  and  $\Delta_2$  such that  $G_i \subseteq \Delta_i \cup N_e$ , where  $G_i$  is the graph induced by the edge set  $\{\mu(f): f \in L \cap V(T_i)\}$  for  $i \in \{1, 2\}$  and  $T_1 \cup T_2 = T \setminus \{e\}$ . Thus  $N_e$  meets G only in  $V(N_e) = \text{mid}(e)$  and its length is |mid(e)|. The vertices of  $\text{mid}(e) = V(G_1) \cap V(G_2)$  are ordered according to a cyclic permutation  $\pi = (\pi_e)_{e \in E(T)}$  on mid(e).

**Theorem 2** [15, 19, 36] Let G be a plane graph satisfying  $bw(G) \le k$  without vertices of degree one. Then there exists a sc-decomposition of G of width at most k. In addition, such a sc-decomposition can be computed in time  $O(n^2 \log n)$ .

## 3 Minor Testing in Planar Graphs

For solving PLANAR H-MINOR CONTAINMENT in single-exponential time  $\mathcal{O}(2^{\mathcal{O}(h)} \cdot n + n^2 \cdot \log n)$ , we now introduce the method of *partially embedded dynamic programming*. We present Algorithm 3.1 as a roadmap on how we proceed in proving our main Theorem 1.

We divide Algorithm 3.1 into three parts, presented in the following three subsections.

Namely, in Sect. 3.1 we show how to find in polynomial time a model of H in G in the case of G having large branchwidth. Otherwise, we may assume that  $\mathrm{bw}(G) \leq c \cdot h$  for some constant c. In this case, PRE-PROC(H) (basically) computes a list of all plane majors A of H up to a fixed size linear in h. This "preprocessing step" is presented in Sect. 3.2. In the sequel, we are interested in graphs A of our list, if A is a minor of G obtained from H by "decontracting" some part, such that on a given subset  $S \subseteq V(G)$ , our graph A looks like a subgraph of G. That is, we consider such



## **Algorithm 3.1**: The main routine for PLANAR *H*-MINOR CONTAINMENT

**Input**: A planar graph G.

**Output**: A model M of H in G, if it exists.

Compute sc-decomposition  $\langle T, \mu, \pi \rangle$  of G of width bw(G).

**if** bw(G) > 42 · h **then** Compute M (Proposition 6)

else Run PRE-PROC(H) to produce a collection  $\mathcal A$  of plane majors of H (Sect. 3.2, Algorithm 3.2); Run partially embedded dynamic programming on  $\langle T, \mu, \pi \rangle$  to find a model M of H in G by using  $\mathcal A$  (Sect. 3.3).

A that have a model  $M_A$  in G such that every edge of  $M_A[S]$  is an m-edge and for every pair of vertices  $u, v \in V(M_A[S])$  there is no path of c-edges in  $M_A$  connecting u and v. Finally, in Sect. 3.3, we proceed by partially embedded dynamic programming bottom-up along a sphere cut decomposition of G. Here we make use of the fact that every middle set S yields a separating noose in an embedding of G. If H has a model  $M \subseteq G$  that intersects S, then the noose comes from a noose in M, which in turn is present in some major A of H of our list. We use this fact to restrict the number of candidates A we need to consider in every single dynamic programming step.

## 3.1 Case of Large Branchwidth

The following proposition allows us to find a model of H in G in the case of G having large branchwidth.

**Proposition 6** Let G and H be planar graphs with |V(G)| = n and |V(H)| = h. There exists a constant  $c \le 42$  such that if  $bw(G) > c \cdot h$ , then G contains a model of H, which can be found in time  $\mathcal{O}(n^2 \cdot \log n + h^4)$ .

*Proof* Since any planar graph H with at least three vertices satisfies  $|E(H)| \le 3|V(H)| - 6$ , by Proposition 2 any planar graph on h vertices is isomorphic to a minor of the  $(14h \times 14h)$ -grid.

With the algorithm of [36], we can find bw(G) in time  $\mathcal{O}(n^2)$ . If bw(G) > 42h, then by Proposition 1, we can find in time  $\mathcal{O}(n^2\log n)$  a model of a  $(14h\times 14h)$ -grid in G. From this grid we find a model of H in G using Proposition 2. To conclude, let us discuss how the proof of Proposition 2 provided in [31] can be easily made constructive. For the sake of presentation, we omit many details that can be found in [31]. Indeed, the proof of [31, Fact (1.5)] consists in subdividing H to obtain an auxiliary planar graph  $H_1$  with  $|V(H_1)| \le 7h$ , which is isomorphic to a minor of a planar Hamiltonian graph  $H_2$  with  $|V(H_2)| \le 14h$ . It is then proved that such a graph  $H_2$  is isomorphic to a minor of a  $(|V(H_2)| \times |V(H_2)|)$ -grid. Then the proof uses simple operations on vertices, edges, and separating triangles. At one point it uses a Hamiltonian cycle, which existence is guaranteed by Whitney's Theorem<sup>2</sup> [38]. There exist a constructive version of Whitney's Theorem, i.e. an algorithm finding a Hamiltonian cycle in triangulated planar graph in linear time [34]. One can then check

<sup>&</sup>lt;sup>2</sup>Whitney's Theorem states that each triangulated planar graph without separating triangles is Hamiltonian.



that the overall running time is dominated by the inductive proof of [31, Fact (1.4)], in which one must find  $\mathcal{O}(h)$  times a separating triangle in a graph on 7h vertices. This procedure can be naïvely done in time  $\mathcal{O}(h^4)$ . Therefore, a rough upper bound for the algorithm that follows from Proposition 2 is  $\mathcal{O}(h^4)$ .

## 3.2 Preprocessing

If the branchwidth of G is at most  $c \cdot h$ , then we compute a sphere cut decomposition of width  $\mathcal{O}(h)$  in time  $\mathcal{O}(n^2)$  by using the algorithm of [19], and we continue with dynamic programming.

In the first step we do preprocessing. Namely, we compute for H a list of auxiliary graphs A with  $H \leq A$ , such that A is a candidate for a model M in G. We will need this collection in the dynamic programming algorithm described in Sect. 3.3. To be precise, we compute a collection  $\mathcal{A}$  of edge-colored plane graphs, each consisting of

- a planar graph  $A_{m,c}$  with  $h \le |V(A_{m,c})| \le h + 1.5 \cdot \text{bw}(G)$ , and  $|E(A_{m,c})| \ge |E(H)|$ , such that H is a contraction minor of  $A_{m,c}$ ,
- a bipartition of the edge set  $E(A_{m,c})$  into m-edges and c-edges such that contracting the c-edges of  $A_{m,c}$  creates a graph isomorphic to H; and
- a drawing  $\Phi$  of  $A_{m,c}$ .

## **Algorithm 3.2**: The preprocessing algorithm: PRE-PROC

```
Input: Planar graph H on h vertices and \ell edges.

Output: Collection \mathcal{A} of edge-colored plane graphs.

Compute all non-isomorphic planar graphs A satisfying h \leq |V(A)| \leq 1.5 \cdot \text{bw}(G) + h and |E(A)| \geq \ell.

for every such graph A do

for every subset X of E(A) of size \ell do

mark the edges in X as m-edges and the edges in E(A) \setminus X as C-edges, resulting in a edge-colored graph A_{m,c}.

for every graph A_{m,c} do

if the graph resulting from contracting the C-edges of A_{m,c} is isomorphic to H then compute its non-equivalent drawings \Phi_1, \ldots, \Phi_q, and add the plane graphs A_{m,c}, \Phi_i for all 1 \leq i \leq q to A.
```

We describe the preprocessing in Algorithm 3.2: the routine PRE-PROC takes as input H and outputs the collection  $\mathcal{A}$  of edge-colored plane graphs  $\langle A_{m,c}, \Phi \rangle$ .

When doing dynamic programming in Sect. 3.3, we compute in each dynamic programming step a subset of collection A consisting of minors of M which represent both H and M.

**Lemma 1** For every planar graph H on h vertices and every constant d, the cardinality of the collection A of non-isomorphic edge-colored plane graphs on  $d \cdot h$  vertices containing a minor isomorphic to H is  $2^{\mathcal{O}(h)}$ . Furthermore, we can compute A in time  $2^{\mathcal{O}(h)}$ .



*Proof* By Proposition 4, the number of non-isomorphic planar graphs A on  $d \cdot h$  vertices (for a constant d) is  $2^{\mathcal{O}(h)}$ . We compute this set in time  $2^{\mathcal{O}(h)}$  using the algorithm of [25]. We partition the edge set of each A into three subsets: the edges that we need to delete, the c-edges, and the m-edges. There are again  $2^{\mathcal{O}(h)}$  possible such partitions, which can be computed in time  $2^{\mathcal{O}(h)}$ . We use the linear time algorithm for planar graph isomorphism [22] to check if after applying the graph operations the resulting graph A' is isomorphic to A. If so, we generate all non-equivalent drawings of A and add them to A by using Proposition 3 and Algorithm 4.1 in [13], again in time  $2^{\mathcal{O}(h)}$ .

Using Lemma 1, we get the following corollary.

**Corollary 1** Algorithm 3.2 is correct and runs in time  $2^{\mathcal{O}(bw(G)+h)}$ .

## 3.3 Partially Embedded Dynamic Programming

From now on, we will refer to an edge-colored plane graph  $\langle A_{m,c}, \Phi \rangle \in \mathcal{A}$  simply as A. In this section, we present the technique of partially embedded dynamic programming. Before proceeding to a formal description, we provide the basic intuition behind our algorithm. Towards this, let us consider graphs  $A \in \mathcal{A}$  satisfying  $H \leq_c A$  and  $A \leq G$ .

We define subgraphs PAST, PRESENT, and FUTURE of A with

- $V(A) = V(PAST) \cup V(PRESENT) \cup V(FUTURE);$
- $E(A) = E(PAST) \dot{\cup} E(PRESENT) \dot{\cup} E(FUTURE);$
- PRESENT  $\subseteq G$ , (i.e., we can obtain A as a minor of G with PRESENT being subgraph of G); and
- E(PAST) ⊆ E(H) (i.e., we can obtain H as a contraction minor of A without contracting edges in PAST).

Here, we slightly abuse notation by assuming that vertex and edge sets in different graphs are actually the same, instead of introducing bijective mappings. Note that we make no assumption about the edges in FUTURE. Intuitively speaking, in partially embedded dynamic programming, we look for potential models M of H with a magnifying glass only in the separators of the sc-decomposition of G. By decontracting H at the separators, we obtain the part PRESENT, which yields a subgraph of G for which we are enabled to apply embedded dynamic programming. For memorizing the rest of the potential model M, we contract all necessary edges to PAST in the processed graph and (almost) all edges to FUTURE in the graph remainder. The picture will be concretized in what follows.

Given a sc-decomposition of G, we proceed with dynamic programming: Every edge e of the sc-decomposition defines a separator  $\operatorname{mid}(e) \subseteq V(G)$  and an associated noose  $N_e$ , which separates the graph  $G_{\operatorname{sub}} \subseteq G$  processed so far from  $G \setminus G_{\operatorname{sub}}$ . At every edge e of the sc-decomposition, we check for every graph A of A all the ways in which the graph  $G_{\operatorname{sub}}$  can be obtained as a major of  $A_{\operatorname{sub}} \subseteq A$  with  $A_{\operatorname{sub}} = (V(\operatorname{PAST}) \cup V(\operatorname{PRESENT}), E(\operatorname{PAST}) \cup E(\operatorname{PRESENT}))$ , where  $\operatorname{mid}(e)$  determines  $V(\operatorname{PRESENT})$ . The noose  $N_e$  comes from a noose in A, and this is controlled



by the ways in which  $N_e$  can be routed through the vertices of A. The number of solutions we get—the *valid partial solutions*—is bounded by the number of combinatorial nooses in A onto which we can map  $N_e$ . When updating the valid partial solutions at two incident edges of the sc-decomposition, we unite PRESENT and PAST of two solutions and set the graph remainder to FUTURE. In a post-processing step, we contract part of PRESENT, namely those edges with at most one endpoint in the newly obtained separator of the sc-decomposition; this part becomes PAST. We then decontract some edges of FUTURE for the next updating step. This concludes the informal description of the algorithm.

In the remaining part of this section, we will precisely describe and analyze the dynamic programming routine with which we achieve the following result:

**Lemma 2** For a plane graph G with a given sc-decomposition  $\langle T, \mu, \pi \rangle$  of G of width bw(G) and a planar graph H on h vertices, we can decide in time  $2^{\mathcal{O}(bw(G)+h)}$ . n whether G contains a model M of H.

## 3.4 Dynamic Programming

We root the sc-decomposition  $\langle T, \mu, \pi \rangle$  at some node  $r \in V(T)$ . For each edge  $e \in T$ , let  $L_e$  be the set of leaves of the subtree rooted at e. The subgraph  $G_e$  of G is induced by the edge set  $\{\mu(v) \mid v \in L_e\}$ . The vertices of mid(e) form a combinatorial noose N that separates  $G_e$  from the residual graph.

Let A be a given plane graph in A. If A is a minor of G, then there exists a plane model M of A in G. Furthermore, for above noose N the intersection  $M \cap N$  forms a noose, both in model M and in candidate A. One basic point of partially embedded dynamic programming is to check how the vertices of the combinatorial noose N are mapped to faces and vertices of A. For a combinatorial noose  $N_A$  in A, we can map N to  $N_A$  bounding (clockwise) a unique subgraph  $A_{\text{sub}}$  of A.

In each step of the algorithm, we compute the solutions for a sub-problem in  $G_e$ , where each solution consists of three parts, namely

- a plane edge-colored graph  $A \in \mathcal{A}$ ;
- a combinatorial noose N<sub>A</sub> in A; and
- a mapping  $\gamma$  from combinatorial noose N to  $N_A$  (defined below).

 $N_A$  has the properties that (a) it bounds (clockwise) a subgraph  $A_{\text{sub}} \subseteq A$  and (b) no vertex in  $V(A_{\text{sub}}) \setminus V(N_A)$  is incident to a c-edge. The subgraph  $A_{\text{sub}}$  is representing the part of model M already computed, whereas the residual graph of A represents the part of M which still has to be verified. For every middle set, we store this information in an array of triples  $\langle A, N_A, \gamma \rangle$ .

We define now valid mappings between combinatorial nooses and describe how partial solutions are stored in the dynamic programming. Then, we give the different DP-steps and finally verify the approach.

## 3.5 Valid Partial Solutions

For a middle set mid(e) of the rooted sc-decomposition  $\langle T, \mu, \pi \rangle$  of plane graph  $G, N = N_e$  is the associated combinatorial noose in G with face-vertex sequence



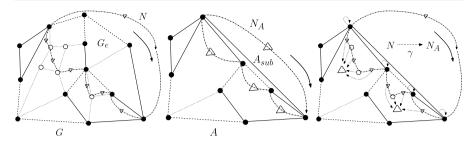


Fig. 2 The figure on the left hand side illustrates the graph G of Fig. 1 with an oriented noose N (dashed) enclosing subgraph  $G_e$ . The figure in the middle shows a graph A of collection A (that is, A is a major of the graph H of Fig. 1) with a noose  $N_A$  enclosing  $A_{\text{sub}}$ . The figure on the right hand side shows a mapping  $\gamma$  from the faces and vertices of N to  $N_A$ 

of  $F(N) \cup V(N)$  separating  $G_e$  from the residual graph. Let  $\mathfrak{N}$  denote the set of all combinatorial nooses of A whose length is at most the length of N and which bound (clockwise) a subgraph  $A_{\text{sub}} \subseteq A$  such that no vertex in  $V(A_{\text{sub}}) \setminus V(N_A)$  is an end-vertex of a c-edge. We now map N to nooses  $N_A \in \mathfrak{N}$ , preserving the order. More precisely, we map vertices of N to both vertices and faces of N. Therefore, we consider partitions of  $V(N) = V_1(N) \cup V_2(N)$  where vertices in  $V_1(N)$  are mapped to vertices of V(N) and vertices in  $V_2(N)$  to faces of V(N).

We define a mapping  $\gamma: V(N) \cup F(N) \to V(A) \cup F(A)$  relating N to the combinatorial nooses in  $\mathfrak{N}$ . For every  $N_A \in \mathfrak{N}$  on faces and vertices of set  $F(N_A) \cup V(N_A)$  and for every partition  $V_1(N) \cup V_2(N)$  of V(N), mapping  $\gamma$  is valid if

- (a)  $\gamma$  restricted to  $V_1(N)$  is a bijection to  $V(N_A)$ ;
- (b) every  $v \in V_2(N)$  and  $f \in F(N)$  satisfy  $\gamma(v) \in F(N_A)$  and  $\gamma(f) \in F(N_A)$ ;
- (c) for every  $v_i \in V(N)$  and subsequence  $[f_{i-1}, v_i, f_i]$  of N: if  $v_i \in V_2(N)$ , then face  $\gamma(v_i)$  is equal to both  $\gamma(f_{i-1})$  and  $\gamma(f_i)$ , and if  $v_i \in V_1(N)$ , then vertex  $\gamma(v_i)$  is incident to both  $\gamma(f_{i-1})$  and  $\gamma(f_i)$ ; and
- (d)  $A_{\text{sub}}$  is a minor of  $G_e$  with respect to (a)–(c).

Items (a) and (b) say where to map the faces and vertices of N to. Item (c) (with (a)) makes sure that if two vertices  $v_h, v_j$  in sequence  $N = [\dots, v_h, \dots, v_j, \dots]$  are mapped to two vertices  $w_i, w_{i+1}$  that appear in sequence  $N_A$  as  $[\dots, w_i, f_i, w_{i+1}, \dots]$ , then every face and vertex between  $v_h, v_j$  in sequence N (here underlined) is mapped to face  $f_i$ . Item (d) simply makes clear that the boundary of  $G_e$  is mapped to the boundary  $A_{\text{sub}}$  such that contracting some part of  $G_e \setminus N$  leads to  $A_{\text{sub}}$ . For an illustration, see Fig. 2.

We assign an array  $\Psi_e$  to each mid(e) consisting of triples, where each triple  $\langle A, N_A, \gamma \rangle$  represents a minor candidate A together with a valid mapping  $\gamma$  from a combinatorial noose N corresponding to mid(e) to a combinatorial noose  $N_A \in \mathfrak{N}$ . The vertices and faces of N are oriented clockwise around the drawing of  $G_e$ . Without loss of generality, we assume for every  $\langle A, N_A, \gamma \rangle$  the orientation of  $N_A$  to be clockwise around the drawing of subgraph  $A_{\text{sub}}$  of A.

Step 0: Initializing the leaves. For every parent edge e of a leaf v of T, we initialize for every  $A \in \mathcal{A}$  the valid mappings from the combinatorial noose bounding the



edge  $\mu(v)$  of G to every combinatorial noose of length at most two in A (clockwise bounding at most one edge of A).

Step I(a): Update process. We update the arrays of the middle sets bottom-up in post-order manner from the leaves of T to root r. During this updating process it is guaranteed that the local solutions for each minor associated with a middle set of the sc-decomposition are combined into a global solution for the overall graph G.

In each dynamic programming step, we compare the arrays of two middle sets  $\operatorname{mid}(e)$  and  $\operatorname{mid}(f)$  in order to create a new array assigned to the middle set  $\operatorname{mid}(g)$ , where e, f, and g have a vertex of T in common. From [15] we know that the combinatorial noose  $N_g$  is formed by the symmetric difference of the combinatorial nooses  $N_e$ ,  $N_f$  and that  $G_g = G_e \cup G_f$ . In other words, we are ensured that if two solutions on  $G_e$  and  $G_f$  bounded by  $N_e$  and  $N_f$  fit together, then they form a new solution on  $G_g$  bounded by  $N_g$ . We now determine when two solutions represented as tuples in the arrays  $\Psi_e$  and  $\Psi_f$  fit together. We update two triples  $\langle A^1, N_A^1, \gamma_1 \rangle \in \Psi_e$  and  $\langle A^2, N_A^2, \gamma_2 \rangle \in \Psi_f$  to a new triple in  $\Psi_g$  if

- $A^1 = A^2 =: A \in \mathcal{A}$  and every edge of A with no endpoint in  $V(N_A^1) \cup V(N_A^2)$  is an m-edge;
- for every  $x \in (V(N_e) \cup F(N_e)) \cap (V(N_f) \cup F(N_f))$ , we have  $\gamma_1(x) = \gamma_2(x)$ ; and
- for the subgraph  $A^1_{\text{sub}}$  of A separated by  $N^1_A$  and the subgraph  $A^2_{\text{sub}}$  of A separated by  $N^2_A$ , we have that  $E(A^1_{\text{sub}}) \cap E(A^2_{\text{sub}}) = \emptyset$  and  $V(A^1_{\text{sub}}) \cap V(A^2_{\text{sub}}) \subseteq \{\gamma(v) \mid v \in V(N_e) \cap V(N_f)\}$ .

That is, we only update solutions with the same graph A and with the two nooses  $N_A^1$  and  $N_A^2$  bounding (clockwise) two edge-disjoint parts of A and intersecting in a consecutive subsequence of both  $N_A^1$  and  $N_A^2$ . If the two solutions on  $N_e$  and  $N_f$  fit together, we get a valid mapping  $\gamma_3: N_g \to N_A^3$  to a noose  $N_A^3$  of A as follows:

- for every  $x \in (V(N_e) \cup F(N_e)) \cap (V(N_f) \cup F(N_f)) \cap (V(N_g) \cup F(N_g))$  we have  $\gamma_1(x) = \gamma_2(x) = \gamma_3(x)$ ;
- for every  $y \in (V(N_e) \cup F(N_e)) \setminus (V(N_f) \cup F(N_f))$  we have  $\gamma_1(y) = \gamma_3(y)$ ; and
- for every  $z \in (V(N_f) \cup F(N_f)) \setminus (V(N_e) \cup F(N_e))$  we have  $\gamma_2(z) = \gamma_3(z)$ .

We have that  $\gamma_3$  is a valid mapping from  $N_g$  to the combinatorial noose  $N_A^3$  that bounds subgraph  $A_{\text{sub}}^3 = A_{\text{sub}}^1 \cup A_{\text{sub}}^2$ .

Step I(b): Post-processing. Before adding a triple  $\langle A, N_A^3, \gamma_3 \rangle$  to array  $\Psi_g$ , we need to manipulate A so that (a) it does not grow too big and (b) it is suitable for future update operations. In A restricted to subgraph  $A_{\text{sub}}^3$ , we contract all c-edges with at least one end-vertex not in  $N_A^3$  in order to fulfill (a). Concerning (b), for every  $B \in \mathcal{A}$  we check for all its nooses  $N_B$  with  $|V(N_B)| = |V(N_A^3)|$  if there is a bijection  $\beta$  from  $N_A^3$  to  $N_B$  such that the following holds. If in a copy of B we contract those c-edges which (i) are in the subgraph counter-clockwise bounded by  $N_B$  and (ii) have at least one end-vertex not in  $N_B$ , then we obtain a edge-colored graph isomorphic to A. We define  $\delta = \gamma_3 \circ \beta$  and we replace  $\langle A, N_A^3, \gamma_3 \rangle$  in array  $\Psi_g$  by those triples  $\langle B, N_B, \delta \rangle$  which validate properties (i) and (ii).

Step 2: Termination. If, at some step, we have a solution where the entire minor H is formed, we terminate the algorithm accepting. That is the case, if for some triple we



have that  $H \leq A_{\text{sub}} \leq A$  and  $A_{\text{sub}}$  is bounded by  $N_A$ . We output model M of H in G represented by this A by reconstructing a solution top-down in  $\langle T, \mu, \pi \rangle$ . If at root r no  $A \in \mathcal{A}$  has been computed, we reject.

## 3.6 Correctness of the Algorithm

In Corollary 1, we already showed that the preprocessing correctly computes the collection  $\mathcal{A}$  of pairwise non-isomorphic edge-colored plane graphs A on at least h and at most  $1.5 \cdot \mathrm{bw}(G) + h$  vertices containing H as a minor. In the update process on the nooses  $N^1$ ,  $N^2$  of two incident edges of the sc-decomposition, we produce graphs with as many vertices since we have for candidate A of H in G that A intersects  $V(N^1) \cup V(N^2)$ , and by Remark 1 that  $|V(N^1) \cup V(N^2)| \leq 1.5 \cdot \mathrm{bw}(G)$  and up to h vertices of A might be outside  $N_A^1$  and  $N_A^2$ .

We have already seen how to map every combinatorial noose of G that identifies a separation of G via a valid mapping  $\gamma$  to a combinatorial noose of A determining a separation of A. Step 0 ensures that every edge of A is bounded by a combinatorial noose  $N_A$  of length two, which is determined by triple  $\langle A, N_A, \gamma \rangle$  in an array assigned to a leaf edge of T. We need to show that Step 1(a) and 1(b) compute a valid solution for  $N_g$  from  $N_e$  and  $N_f$ , given incident edges e, f, g. We note that the property that the symmetric difference of the combinatorial nooses  $N_e$  and  $N_f$  forms a new combinatorial noose  $N_g$  is passed on to the combinatorial nooses  $N_A^1, N_A^2$ , and  $N_A^3$  of A, too. If the two solutions fit together, then  $A_{\text{sub}}^1$  of A separated by  $N_A^1$  and subgraph  $A_{\text{sub}}^2$  of A separated by  $N_A^2$  only intersect in the image of  $V(N_e) \cap V(N_f)$ . We observe that  $N_A^1$  and  $N_A^2$  intersect in a consecutive alternating subsequence with order reversed to each other, i.e.,  $N_A^1 |_{N_e \cap N_f} = N_A^2 |_{N_e \cap N_f}$ , where  $\overline{N_A}$  means the reversed sequence  $N_A$ . Since every oriented  $N_A$  uniquely identifies a separation of E(A), we can easily decide whether two triples  $\langle A, N_A^1, \gamma \rangle \in \Psi_e$  and  $\langle A, N_A^2, \gamma \rangle \in \Psi_f$  fit together and bound a new subgraph  $A_{\mathrm{sub}}^3$  of A. In Step 1(b), we contract all c-edges in  $A_{\mathrm{sub}}^3$  with an end-vertex in  $V(A_{\mathrm{sub}}^3) \setminus V(N_A^3)$  such that  $A_{\mathrm{sub}}^3$ represents a subgraph of H. Next we blow up  $\overline{A_{\text{sub}}^3}$ , that is the subgraph of A on the other side which  $N_A^3$  bounds counter-clockwise. From A we compute every possible graph partitioned into one subgraph isomorphic to  $A_{\mathrm{sub}}^3$  and one subgraph major of  $\overline{A_{\text{sub}}^3}$  separated by a noose bijectively mapped from  $N_A^3$ . If there exists an  $A \in \mathcal{A}$  which is a minor of G, then at some point we will enter Step 2 and produce the entire model M.

## 3.7 Running Time of the Algorithm

By Corollary 1, computing the collection  $\mathcal{A}$  in the preprocessing step can be done in time  $2^{\mathcal{O}(\mathrm{bw}(G)+h)}$ . We now give an upper bound on the size of each array  $\Psi$ . For each  $A \in \mathcal{A}$ , the number of combinatorial nooses in  $\mathfrak{N}$  we are considering is bounded by the total number of combinatorial nooses in A, which is  $2^{\mathcal{O}(\mathrm{bw}(G)+h)}$  by Proposition 5. The number of partitions of vertices of any combinatorial noose N is bounded by  $2^{|V(N)|}$ . Since the order of both  $N_A$  and N is given, we only have  $2 \cdot |V(A)|$  possibilities to map vertices of N to  $N_A$ , once the vertices of N are partitioned. Thus, in



an array  $\Psi_e$  we may have up to  $2^{\mathcal{O}(|V(A)|)} \cdot 2^{|V(N)|} \cdot |V(A)|$  triples  $\langle A, N_A, \gamma \rangle$ . We first create all triples in the arrays assigned to the leaves. Since middle sets of leaves only consist of an edge in G, we get arrays of size  $\mathcal{O}(|V(A)|^2)$ , which we compute in the same asymptotic running time. When updating middle sets  $\mathrm{mid}(e)$ ,  $\mathrm{mid}(f)$ , we compare every triple of array  $\Psi_e$  to every triple in array  $\Psi_f$  to check if two triples fit together. We can compute the unique subgraph  $A^1_{\mathrm{sub}}$  (resp.  $A^2_{\mathrm{sub}}$ ) described by a triple in  $\Psi_e$  (resp.  $\Psi_f$ ), compare two triples in  $\Psi_e$  and  $\Psi_f$ , and create a new triple in time linear in the order of V(N) and V(H). For adding a new triple to  $\Psi_g$  in the post-processing, we apply the color coding technique [4] for computing each of the  $2^{\mathcal{O}(\mathrm{bw}(G)+h)}$  nooses in  $\mathfrak N$  in the same asymptotic running time.

This completes the proof of Lemma 2.

*Proof of Theorem 1* We put everything together by verifying Algorithm 3.1. We produce in time  $\mathcal{O}(n^2 \cdot \log n)$  a sc-decomposition of input graph G [19]. Next, either we can immediately compute a model of H in G in time  $\mathcal{O}(n^2 \cdot \log n + h^4)$  (Proposition 6) or we run our 2-step-algorithm: we produce all majors of the minor pattern (Lemma 1) with Algorithm 3.2 in time  $2^{\mathcal{O}(h)}$ , and run partially embedded dynamic programming in time  $2^{\mathcal{O}(h)} \cdot n$  (Lemma 2).

#### 4 Conclusions and Further Research

In this paper we showed that PLANAR H-MINOR CONTAINMENT is solvable in time  $\mathcal{O}(2^{\mathcal{O}(h)} \cdot n + n^2 \cdot \log n)$  for a host graph G on n vertices and a pattern H on h vertices. That is, we showed that the problem can be solved in *single-exponential* time in h, significantly improving all previously known algorithms. Similar to [13], we can enumerate and count the number of models within the same time bounds.

Let us discuss some interesting avenues for further research concerning minor containment problems. First, it seems possible to solve other variants of planar minor containment in single-exponential time using our approach, such as testing for a contraction minor, an induced minor, or a topological minor, as it has been recently done in [1] for general host graphs using completely different techniques. Also, it would be interesting to count the number of non-isomorphic models faster than just by enumerating models and removing isomorphic duplicates.

An important question is if, up to some assumption from complexity theory, the running time of our algorithm is tight. In other words, is there a  $2^{o(h)} \cdot n^{\mathcal{O}(1)}$  algorithm (i.e., a *subexponential* algorithm) solving PLANAR *H*-MINOR CONTAINMENT, or the existence of such an algorithm would imply the failure of, say, the Exponential Time Hypothesis? A first step could be to study the existence of subexponential algorithms when the pattern is further restricted to be a *k*-outerplanar graph for some constant *k*, or any other subclass of planar graphs.

Conversely, single-exponential algorithms may exist for host graphs more general than planar graphs. The natural candidates are host graphs embeddable in an arbitrary surface. One possible approach could be to use the framework recently introduced in [33] for performing dynamic programming for graphs on surfaces. The main ingredient of this framework is a new type of branch decomposition of graphs



on surfaces, called *surface cut decomposition*, which plays the role of sphere cut decompositions for planar graphs.

Finally, an interesting open question is whether the PLANAR DISJOINT PATHS problem, a problem closely related to minor testing (see [35]), can be solved with a similar running time.

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