Another Time-Complexity Analysis for
Maximal Clique Enumeration Algorithm CLIQUES

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Abstract We revisit the maximal clique enumeration algorithm CLIQUES that appeared in Theoretical Computer Science 2006. It is proved to work in $O(3^{n/3})$-time in the worst-case for an n vertex graph. In this note, we extend the time-complexity analysis with respect to the number of maximal cliques, an issue that was left as an open problem since TCS 2006.
Key words maximal clique, maximal independent set, enumeration, algorithm, time-complexity, delay

1. Introduction

A *clique* is defined to be a subgraph in which every pair of vertices are adjacent. In particular, it is *maximal* if it is not contained in a properly larger clique. Given a graph, the enumeration of all its maximal cliques is a fundamental and important problem in graph theory. In addition, it has many applications in practice as in clustering, data mining, bioinformatics, social networks, and many others. An *independent set* of a graph G is a clique of the complementary graph $\bar{G}$.

Tsukiyama et al. was the first to give an algorithm for enumerating all maximal independent sets with theoretical time-complexity analysis. Given a graph G with n vertices and m edges, their algorithm takes $\mu O(nm)$-time for enumerating all maximal independent sets, where μ is the number of maximal independent sets. This implies that their algorithm enumerates all maximal cliques in $\alpha O(n(n^2 - m))$-time for G, where α is the number of maximal cliques in G. Their approach is based on so-called *reverse search* [1] approach. On the other hand, Tomita et al. [20], [21] presented another algorithm CLIQUES for enumerating all maximal cliques. CLIQUES is independently developed based on *depth-first search* algorithms for finding a *maximum* clique

[11], [19], [23], but the pruning techniques in CLIQUES are found to be the same as in Bron and Kerbosch algorithms [4] for the same problem. CLIQUES is proved to enumerate all maximal cliques in $O(3^{n/3})$-time for an n-vertex graph in the worst-case. This is optimal as a function of n since there exist up to $3^{n/3}$ maximal cliques as in Moon-Moser graphs [17]. CLIQUES is the first algorithm with theoretical time-complexity analysis in this approach.

In the reverse search approach after Tsukiyama et al., steady improvements have been made by Chiba and Nishizeki [6], Johnson et al. [13], Makino and Uno [15], Chang et al. [5], and Conte et al. [7], [9]. One of Conte et al.'s algorithms requires $\alpha O(maxQ \cdot d\Delta)$-time for enumerating all maximal cliques, where $maxQ$ is the size of a maximum clique, d is the *degeneracy* of G, that is defined as the smallest number such that every subgraph of G contains a vertex of degree at most d. Based on CLIQUES, Eppstein et al. proposed an improved algorithm that runs in $O(dn3^{d/3})$-time, where *degeneracy d* is much smaller than n for sparse graphs. In general, it is experimentally confirmed that if the number of cliques is small then the running time for their enumeration is small. However, no theoretical time-complexity analysis with respect to the number of maximal cliques is

made after CLIQUES [21], where the problem is noted in [21] as an important open problem. This is in contrast to the time-complexity analysis in reverse search approach. Example runs by two approaches for the same problem can be found in [22]. For early histories and applications, see the excellent review articles [18], [2] where CLIQUES [20] is also included for review.

In this note, we are concerned with another time-complexity analysis of CLIQUES and others in the same approach other than that with respect to the number of vertices in the worst-case.

2. Definitions and notation

(1) We are concerned with a simple undirected *graph* $G = (V, E)$ with a finite set V of *vertices* and a finite set E of *unordered* pairs (v, w) of distinct vertices, called *edges*. A pair of vertices v and w are said to be *adjacent* if $(v, w) \in E$.

(2) For a vertex $v \in V$, let $\Gamma(v)$ be the set of all vertices that are adjacent to v in $G = (V, E)$, i.e., $\Gamma(v) = \{w \in V \mid (v, w) \in E\}$ ($\not\ni v$).

(3) For a subset $W \subseteqq V$ of vertices, $G(W) = (W, E(W))$ with $E(W) = \{(v, w) \in W \times W \mid (v, w) \in E\}$ is called a *subgraph* of $G = (V, E)$ *induced* by W. For a set W of vertices, $|W|$ denotes the number of elements in W.

(4) Given a subset $Q \subseteqq V$ of vertices, if $(v, w) \in E$ for all $v, w \in Q$ with $v \neq w$ then the induced subgraph $G(Q)$ is said to be a *clique*. In this case, we may simply say that Q is a clique. If a clique is not a proper subgraph of another clique then it is called a *maximal* clique.

3. Maximal clique enumeration algorithm CLIQUES

We revisit a depth-first search algorithm, CLIQUES [20], [21] for enumerating all maximal cliques of an undirected graph $G = (V, E)$ with $|V| = n$. All maximal cliques found are output in a tree-like form. The basic framework of CLIQUES is almost the same as that for finding a *maximum* clique *without the bounding condition* [11], [19], [23]. We maintain a global variable $Q = \{p_1, p_2, ..., p_h\}$ that consists of the vertices of a current clique, and let $SUBG = V \cap \Gamma(p_1) \cap \Gamma(p_2) \cap \cdots \cap \Gamma(p_h)$. We begin the algorithm by letting $Q \leftarrow \emptyset$ and $SUBG \leftarrow V$ (the set of all vertices). We select a certain vertex p from $SUBG$ and add p to Q. Then, we compute $SUBG_p = SUBG \cap \Gamma(p)$ as the new set of vertices

in question. In particular, the initially selected vertex $u \in SUBG$ is called a *pivot*. This CLIQUES() procedure is applied recursively while $SUBG_p \neq \emptyset$.

We describe two methods to prune unnecessary parts of the searching, which happen to be the same as in the Bron-Kerbosch algorithms [4]. We regard the set $SUBG$ ($= V$ at the beginning) as an *ordered* set of vertices, and we continue to enumerate maximal cliques from vertices in $SUBG$ step-by-step in this order

The first pruning method : Let $FINI$ be a subset of vertices of $SUBG$ that have already been processed by the algorithm ($FINI$ is short for <u>FINISHED</u>). Then we denote by $CAND$ the set of remaining <u>CANDIDATE</u> for expansion: $CAND = SUBG \setminus FINI$. Initially, $FINI \leftarrow \emptyset$ and $CAND \leftarrow SUBG = V$. In the subgraph $G(SUBG_p)$ with $SUBG_p = SUBG \cap \Gamma(p)$, compute
$$CAND_p = CAND \cap \Gamma(p),$$
$$FINI_p = FINI \cap \Gamma(p).$$
Then only the vertices in $CAND_p$ excluding $FINI_p$ can be candidates for expanding the clique $Q \cup \{p\}$ to find *new* larger cliques.

The second pruning method : For the initially selected *pivot* u in $SUBG$, any maximal clique Q' in $G(SUBG \cap \Gamma(u))$ is not maximal in $G(SUBG)$, since $Q' \cup \{u\}$ is a larger clique in $G(SUBG)$. Therefore, searching for maximal cliques from $SUBG \cap \Gamma(u)$ should be excluded.

Algorithm 1: Algorithm CLIQUES in [21]

Input : A graph $G = (V, E)$.
Output: All maximal cliques in G.

1   CLIQUES $(\emptyset, V, \emptyset)$
2   **Function** CLIQUES(Q, *CAND*, *FINI*)
3   **if** $CAND \cup FINI = \emptyset$ **then**
       // Q is a maximal clique
4   **print** (*"clique,"*)
5   **else**
6       $u \leftarrow$ a vertex in $CAND \cup FINI$
7   maximizing $|CAND \cap \Gamma(u)|$
8   **foreach** $p \in CAND \setminus \Gamma(u)$ **do**
9   **print** (*p*,",")
10   CLIQUES $(Q \cup \{p\}, CAND \cap \Gamma(p), FINI \cap \Gamma(p))$
11        $CAND \leftarrow CAND \setminus \{p\}$
12        $FINI \leftarrow FINI \cup \{p\}$
13   **print** (*"back,"*)

Taking the previously described pruning method into consideration, the only search subtrees to be expanded are from vertices in $(SUBG \setminus (SUBG \cap \Gamma(u))) \setminus FINI = CAND \setminus \Gamma(u)$. Here, to minimize $|CAND \setminus \Gamma(u)|$, we choose the pivot $u \in SUBG$ that *maximizes* $|CAND \cap \Gamma(u)|$, which is *crucial* to establish the *optimality* of the worst-case time-complexity of the algorithm.

The algorithm CLIQUES [20], [21] is shown as **Algorithm 1** in the previous page.

It enumerates all maximal cliques in $O(3^{n/3})$-time in the worst-case based upon the above methods, where all maximal cliques enumerated are presented in a tree-like form [20], [21]. We can easily obtain a tree representation of all the maximal cliques from the output sequence, where a dummy root is added to form a tree (Fig.4 of [21]). The tree-like output format is also important *practically*, since it saves space in the output file.

Search tree

The process of enumerating all maximal cliques of $G = (V, E)$ by CLIQUES $(\emptyset, V, \emptyset)$ is represented by a *search tree*:

• The root of the search tree is a newly introduced *dummy* root p_0 ($\notin V$) to form a tree.

• Every vertex in V is a child of the dummy root p_0. The vertex of V is called a *node* in the search tree.

• Assume we have a path from the dummy root p_0 to a certain node p_h in the search tree as a sequence of nodes $p_0, p_1, p_2, ..., p_h$, and let

$$SUBG_h = V \cap \Gamma(p_1) \cap \Gamma(p_2) \cap \cdots \cap \Gamma(p_h).$$

Then, every vertex in $SUBG_h$ is a child of p_h in the search tree.

Suppose u with the maximum degree is chosen as a pivot in $SUBG_h$ then every node in $\Gamma(u)$ is a *leaf* since it should not be expanded by CLIQUES according to the second pruning method. Such a leaf in $\Gamma(u)$ is called a *black node*.
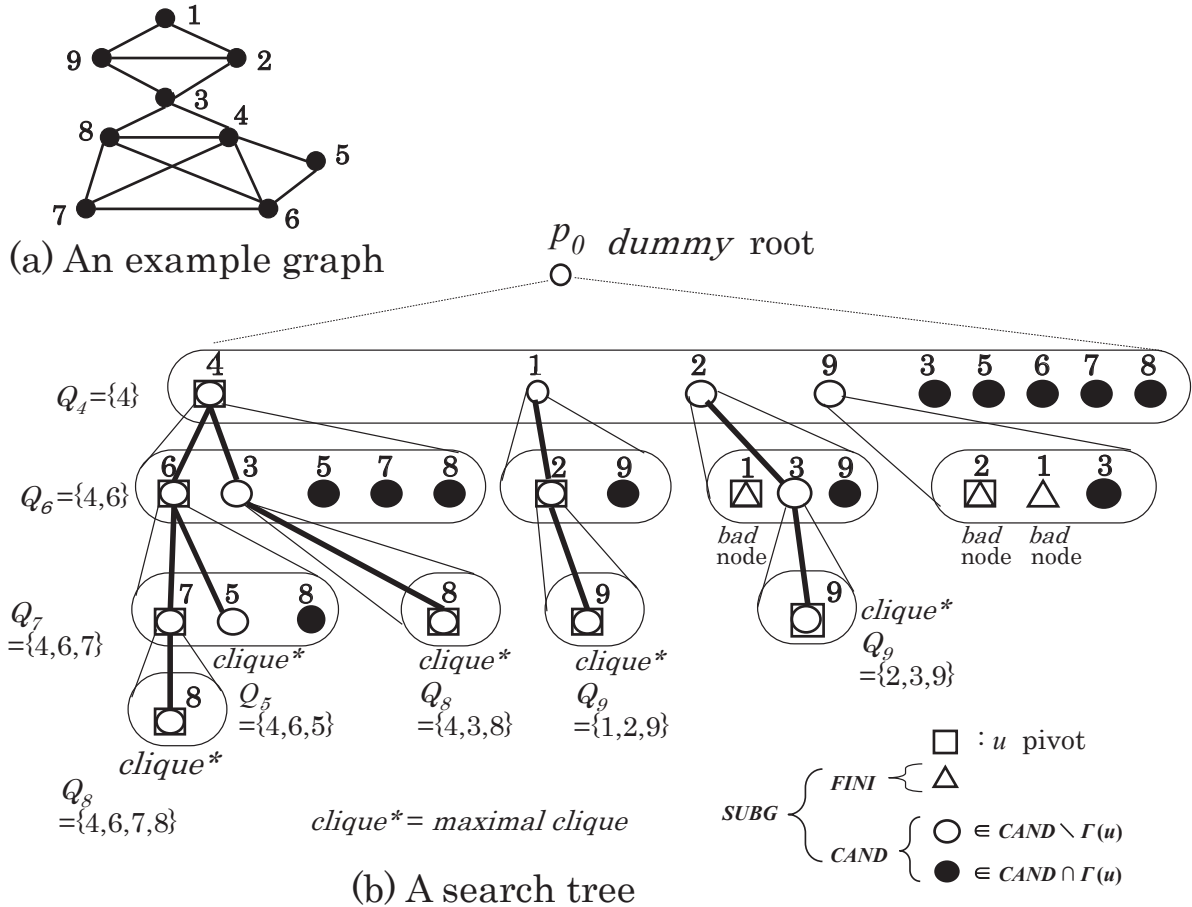

(a) An example graph

(b) A search tree

Figure 1   An example run of CLIQUES

Suppose

$$SUBG_h = FINI_h \cup CAND_h, \; q_i, q_j \in SUBG_h \setminus \Gamma(u),$$

where $i < j$ and q_i, q_j are adjacent with $q_i \in FINI_h$.

Then $q_j \in SUBG_h$ has a child q_i since q_j and q_i are adjacent in $SUBG_h$, but the child q_i should not be expanded by CLIQUES according to the first pruning method, and hence it is a leaf of the search tree. Such a leaf q_i is called a *bad node* or *bad leaf*. Note that if a bad node were expanded it could not lead to a *new* maximal clique.

- When the above $SUBG_h$ is a singleton $\{q_1\}$ then the q_1 is a leaf in the search tree.

The search tee consists of the nodes from $V \cup \{p_0\}$ and the parent-child relationship holds if and only if one of the above conditions holds.

Let q_i be a child of q_h, then the set $\{p_1, p_2, ..., p_h, p_i\}$ constitutes a *clique*. This clique is called an *accompanying clique* and is denoted by $Q_{p_0, p_1, ..., q_i}$, or simply Q_{q_i}, or Q when it is clear.

See Figure 1 of a search tree for an example run, and Figure 2 of a general search tree.

4. Bron-Kerbosch algorithms

Let us recall the antecedent algorithms by Bron and Kerbosch for enumerating maximal cliques [4].

The main difference between theirs (Bron-Kerbosch) and ours (CLIQUES) is how the pivot node is chosen on Line 7.

We call BK the original Bron-Kerbosch algorithm *without* pivoting, which corresponds to entirely removing Line 7, and replacing Line 8 with "**foreach** $v \in CAND$ **do**".

Next, we call BKP the original Bron-Kerbosch algorithm with pivoting, where no specific pivot choice is required, i.e., Line 7 can be replaced with "$u \leftarrow$ an arbitrary vertex in $CAND \cup FINI$."

In addition, CLIQUES outputs all maximal cliques in a *tree*-like format in order to avoid the time for outputting a maximal clique every time it is found that is proportional to the size of a maximal clique found.

The pivot selection of the maximum-degree and the tree-like outputting are crucial in CLIQUES so that it accomplishes the worst-case optimal $O(3^{n/3})$-time complexity.

The original Bron-Kerbosch algorithm BKP (with ar-

bitrary pivot choice) should take $O(n3^{n/3})$-time in the worst-case, while BK (without pivoting at all) runs in $\Theta(n2^n)$ on a complete graph, as it essentially generates all subsets of every clique.

5. The time-complexity of CLIQUES

We analyze the time-complexity of CLIQUES with respect to the output.

5.1 The overall time-complexity of CLIQUES with respect to the output

5.1.1 Size of the search tree

In this section we aim at bounding the number of nodes in the search tree with respect to the number of solutions.

Suppose we have a path from the dummy root p_0 to a certain node p_h in the search tree as a sequence of nodes $p_0, p_1, p_2, ..., p_h$. Then the set $\{p_1, p_2, ..., p_h\}$ constitutes an *accompanying clique*.

We are interested in the accompanying clique Q across different search tree nodes, and in particular let us observe the following:

Lemma 1. *The accompanying clique Q is distinct in any <u>internal</u> (non-leaf) node of a search tree of* CLIQUES.

Proof. Let x and y be any pair of non-root internal nodes in the search tree of CLIQUES, where x is generated before y in the search tree. Let the nearest common ancestor of x and y be p_h, and let the path from the dummy root p_0 to node p_h be $p_0, p_1, p_2, ..., p_h$ with the accompanying clique $Q_{p_h} = \{p_1, p_2, ..., p_h\}$. In addition,
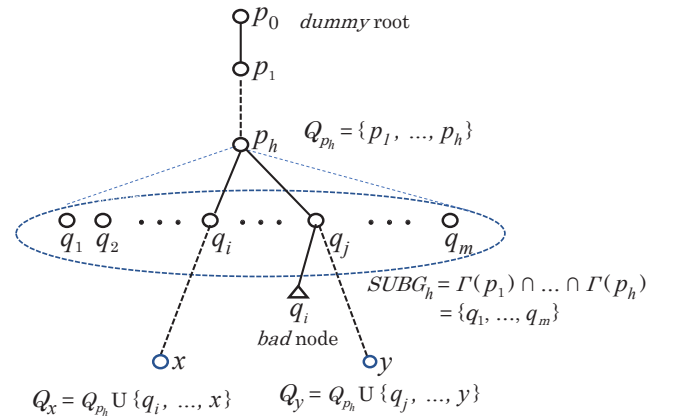

Figure 2   A Search tree

let the path from p_h to x be $p_h, q_i, ..., x$ and that from p_h to y be $p_h, q_j, ..., y$, respectively, and let $SUBG_h = \Gamma(p_1) \cap \Gamma(p_2) \cap ... \cap \Gamma(p_h) = \{q_1, q_2, ..., q_i, ..., q_j, ..., q_m\}$ $(i < j)$. See Figures 1 and 2. So, $Q_x = Q_{p_h} \cup \{q_i, ..., x\}$ and $Q_y = Q_{p_h} \cup \{q_j, ..., y\}$. When we visit node q_j we see that node q_i is in $FINI_h = SUBG_h \setminus CAND_h$ where $CAND_h = \{q_j, ..., q_m\}$ at that moment. If $q_i \notin \Gamma(q_j)$ then it clearly follows that $q_i \notin Q_y$. When $q_i \in \Gamma(q_j)$, q_i in the descendants of q_j is a *bad* node (leaf) since the previous node q_i is in $FINI_h$. Moreover, the *bad* node q_i is not in the path $q_j, ..., y$ since all nodes $q_j, ..., y$ are <u>internal</u> nodes by the assumption. Thus, $q_i \notin Q_y$. In any case, it follows that $q_i \in Q_x \setminus Q_y$ and the lemma is proved. $\qquad\square$

We observe that the accompanying clique Q is always a subset of a *maximal* clique. Let $maxQ$ be the size of a maximum clique and μ the number of maximal cliques in G.

Lemma 2. *The number of nodes of a search tree of* CLIQUES *is at most* $(1 + \Delta)\mu(2^{maxQ} - 1)$.

Proof. The above observation immediately implies that the number of internal nodes is bounded by the number of distinct non-maximal cliques of G. Each maximal clique has at most $2^{maxQ} - 1$ distinct subsets whose size is less than or equal to $maxQ - 1$. (The subset of size 0 corresponds to the dummy root p_0, and the subset of size $maxQ$ is excluded.) So, the number of internal nodes is at most $\mu(2^{maxQ} - 1)$ by **Lemma 1**. Accordingly, the number of leaves of the search tree is at most $\Delta\mu(2^{maxQ} - 1)$. Summarizing the above, the number of nodes of a search tree is at most $(1+\Delta)\mu(2^{maxQ}-1)$. $\quad\square$

5.1.2 The overall time-complexity of CLIQUES

Theorem 1. *Since each node can be executed in $O(n^2)$-time [21], the running time of* CLIQUES *is $O((1 + \Delta)\mu(2^{maxQ} - 1)poly(n))$ for some polynomial $poly(n)$.*
$\qquad\square$

Consider an algorithm $\mathcal{A}$ that enumerates all solutions in the given problem where the number of solutions in the problem is M. If algorithm $\mathcal{A}$ requires $O(M \cdot poly(n))$-time for some polynomial $poly(n)$ to enumerate all solutions then the complexity of algorithm $\mathcal{A}$ is called *amortized* polynomial time.

Corollary 1. *If the size of a maximum clique $maxQ =$*

$O(\lg n)$ *then the running time of* CLIQUES *is amortized polynomial.*

Proof. We have $maxQ \le c \lg n$ for some constant c from the assumption.
Therefore, $(1+\Delta)\mu(2^{maxQ}-1) \le (1+\Delta)\mu(2^{c \lg n}-1) = (1 + \Delta)(n^c - 1) \cdot \mu$.
Hence the result. $\qquad\square$

Note that the properties in this Section 5.1 also hold for BK and BKP.

5.2 Delay of CLIQUES

Next we consider the *delay* of the algorithm, that is, the maximum time which can elapse between two consecutive outputs of a solution.

Delay of BK and BKP is easy to analyze and is $\Omega(3^{n/3})$-time, that is described separately later in Section 6.1.

As for CLIQUES, the situation is challenging. The maximum-degree pivot choice is in practice very effective in pruning search subtrees that lead to no solutions, and we were unable to build ad-hoc counterexamples where this choice leads to an exponentially long time between two outputs.

We, however, manage to show a weaker, but essentially similar, result: we prove that if CLIQUES has polynomial delay, then $P = NP$.

It is worth observing that a recursive call of Algorithm 1 (ignoring the cost of its children calls) takes polynomial time, namely $O(n^2)$, so exponential delay means the algorithm must encounter an exponentially long sequence of consecutive recursive calls that perform no output. Furthermore, as the depth of the tree is $O(maxQ)$, a leaf is always reached in polynomial time, so exponential delay means the algorithm must encounter an exponentially long sequence of consecutive *leaves* that perform no output (which we call *bad leaves*).

5.2.1 Hardness of the extension problem

Firstly, let us define what is called the *extension problem* for maximal cliques, and prove that it is NP-complete.

Problem 1 (Extension Problem, EXT-P($G(V, E), X$)). *Given a graph $G(V, E)$ and $X \subset V$, does G have a maximal clique Q that does not intersect X?*

This problem is a common building block of enumeration algorithms based on binary partition, although typically challenging when dealing with maximal solutions [3], [8], [14]. Looking at a recursive call of CLIQUES, setting $X \equiv FINI$ and $V \setminus X \equiv CAND$, we can observe how it essentially corresponds to asking "will a maximal clique be output in its recursion subtree?" Answering this question would allow excellent pruning, as we could identify all useless recursive calls, resulting in no bad leaves and thus polynomial delay.

Unfortunately this problem is NP-complete, as we show by a reduction from CNFSAT.

Theorem 2. *The extension problem for maximal cliques* (EXT-P$(G(V,E), X)$) *is NP-complete.*

Proof. Let $\mathcal{F}$ be a CNF boolean formula on h variables $v_1, \ldots, v_h$ and l clauses $c_1, \ldots, c_l$, and let the positive and negative literals of the variable v_i be represented by v_i and $\neg v_i$.

We build a graph G, with a suitable vertex set X, such that $\mathcal{F}$ can be satisfied if and only if G has a maximal clique not intersecting X.

Let V be as follows:

- A vertex p_i for each positive literal v_i in $\mathcal{F}$
- A vertex n_i for each negative literal $\neg v_i$ in $\mathcal{F}$
- A vertex c_i for each clause of $\mathcal{F}$

Let $E(G)$ be as follows:

- For all distinct i and j, $E(G)$ contains (p_i, p_j), (p_i, n_j), (n_i, p_j), and (n_i, n_j), *i.e.*, all literals are connected to all others (positive and negative) except their own negation.

- $E(G)$ contains (c_i, p_i) if literal v_i does *not* appear in c_i. Similarly, $(c_i, n_i) \in E(G)$ if $\neg v_i$ does *not* appear in c_i, *i.e.*, each clause is connected to all literals *except* those that satisfy it.

An example is shown in Figure 3.

Now, let X be the set of all vertices c_i corresponding to clauses. Note that $V \setminus X$ is the set of all vertices corresponding to literals. Furthermore, observe how a clique cannot contain both p_i and n_i, and any set $S \subseteq V \setminus X$ is a clique if and only if it does *not* contain both a literal and its negation: it follows that cliques in $V \setminus X$ correspond exactly to valid truth assignments to the literals.

Furthermore, observe that if S does *not* contain any literal satisfying a clause c_i, then c_i is adjacent to all literals in S (due to how edges are placed in E), thus S
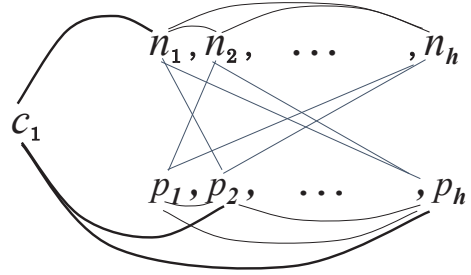

Figure 3   Example construction of G from the $\mathcal{F}$ formula for a clause $c_1 = (v_1, \neg v_2, \neg v_h)$

is not maximal because c_i can be added to it. So if the literals in S do not satisfy some clause c_i, then S is not maximal.

Conversely, if S is maximal and spans all variables (*i.e.*, either p_i or n_i is in S for any i), it means that each c_i is *not* adjacent to some vertex in S: this means c_i is satisfied by the corresponding literal, since clauses are adjacent to all literals that do not satisfy them. It follows that $S \subseteq V \setminus X$ is a maximal clique if and only if its corresponding literals are a satisfying assignment of $\mathcal{F}$ that assigns a truth value to all variables.

Finally, any CNF formula that has a satisfying assignment, trivially also has a satisfying assignment that assigns truth values to all variables (one can just choose an arbitrary value to each missing variable). It follows that EXT-P$(G(V,E), X)$ has positive answer if and only if $\mathcal{F}$ can be satisfied.

It only remains to show that EXT-P$(G(V,E), X)$ is NP: indeed testing a solution corresponds to verifying the maximality of a given clique, that can trivially be done in $O(n^2)$ time. The proof is complete. □

5.2.2 CLIQUES has polynomial delay only if P=NP

Now, given $G(V,E)$ and X, we build a graph $G'(V', E')$ such that, if CLIQUES runs on $G'(V', E')$ with polynomial delay, then we can solve the NP-complete problem (EXT-P$(G(V,E), X)$) in polynomial time.

$G'(V', E')$ is built as follows. In particular, $V' = Q \cup X \cup \{v\} \cup P$, where:

- $Q = \{q_1, \ldots, q_{|Q|}\}$ is a clique of size $|Q| > |X| + |V| + 1$ (large enough so that one vertex is chosen as a pivot in the first recursive call of CLIQUES)

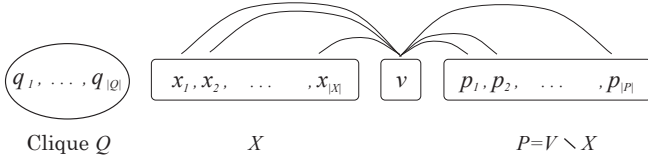- $X = \{x_1, \ldots, x_{|X|}\}$ are the vertices of X from

Figure 4   Graph $G'(V', E')$

EXT-P$(G(V, E), X)$.

- v is a vertex connected to all of X and P (but not to Q)

- $P = \{p_1, \ldots, p_{|P|}\}$ are the vertices of $V \setminus X$ from EXT-P$(G(V, E), X)$.

Where the vertices from $G(V, E)$ are connected as in G.

If we run CLIQUES on G', it will first choose q_1 as a pivot without loss of generality, and q_2, ..., $q_{|Q|}$ will become black nodes and be skipped.

The algorithm will find the maximal clique Q in $O(|Q|^3)$-time, then will start processing the remaining vertices, *i.e.*, $X \cup \{v\} \cup P$. By processing we mean they are considered in the **foreach** loop of the root recursive call of Algorithm 1. As the algorithm does *not* specify in which order the vertices are processed, we will assume this is in the same order as they appear in Figure 4, *i.e.*, first X, then v, then P.

Take the moment when v is processed:

We have that $CAND \cap \Gamma(v)$ is exactly P, while $FINI \cap \Gamma(v)$ is X because we already processed vertices of X.

If CLIQUES has polynomial delay, it must either find a new maximal clique or terminate, in polynomial time: as any maximal clique containing vertices of X has already been found, this process will output a new maximal clique if and only if there is a maximal clique in P that cannot be extended with vertices of X, *i.e.*, since P corresponds to $V \setminus X$, there is a maximal clique in $G(V, E)$ that does not intersect X.

Finally, since CLIQUES may spend exponential time before processing v, we want to skip this time: we can set up CLIQUES directly with arbitrary $CAND$ and $FINI$ by running

CLIQUES$(\{v\}, P, X)$

This way CLIQUES starts from the desired moment, and if it has polynomial delay, it will find the solution to EXT-P$(G(V, E), X)$ in polynomial time. More formally, we can state that:

Theorem 3. *If Algorithm* CLIQUES *[21] has polynomial delay, then $P = NP$.*   □

In addition, the following is a corollary to **Theorem 1**.

Corollary 2. *If $maxQ = O(\lg n)$ then the running time of* CLIQUES *is polynomial delay.*

Proof. Let p_l and $p_{l'}$ be any pair of consecutive outputs of *maximal* cliques in the search tree. (Consider Figure 2 where the internal node x is replaced by a leaf p_l and the internal node y by a leaf $p_{l'}$.) Then we can show that the number of nodes between p_l and $p_{l'}$ in the search tree is at most $2(1 + \Delta)(2^{maxQ} - 1)$ as in the proof of **Theorem 1.**

If $maxQ = O(\lg n)$ then $2(1 + \Delta)(2^{maxQ} - 1) \leqq 2(1 + \Delta)(n^c - 1)$.

Hence the result.   □

6.   Delay of the other algorithms

6.1   Delay of Bron-Kerbosch Algorithms

For BK and BKP the delay can be easily assessed. Consider a grap G built as follows: a clique A of constant size, a Moon-Moser graph M on k vertices [17], a vertex x adjacent to all of M, and another clique B of constant size. A graphical representation is given in Figure 5. Observe that the total number of vertices in G is $n = k + O(1)$. Furthermore, all maximal cliques in G except A and B contain x. Consider now BKP on G: in the root recursive call, assume a vertex a of A is chosen as the first pivot (thus recursion on all other vertices of A is prevented by the pivot rule), assume that vertices are processed in the following order (again, by processing we mean they are considered in the **foreach** loop): First a, then x, then vertices of M, then finally vertices of B. All maximal cliques involving x are found

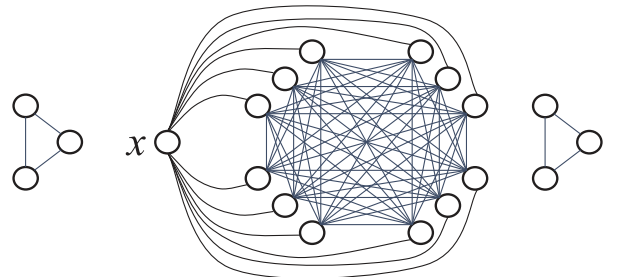Clique   Vertex x M-M graph   Clique

Figure 5   The graph built in Section 6.1.

in the recursion subtree when processing x. Once the algorithm backtracks and starts visiting vertices of M, assume that x is never selected as pivot (an acceptable assumption if the pivot choice is arbitrary). Then by [21] we know BKP will take $\Omega(3^{k/3}) = \Omega(3^{n/3})$ time to process the Moon-Moser graph M, but no new solution is found in this time, because all cliques in M can be extended with x, finally the algorithm processes vertices of B and outputs B, giving us a delay of $\Omega(3^{n/3})$-time. The same bound holds for BK, since its recursion tree includes that of BKP. We obtain the following:

Theorem 4. *The* BK *and* BKP *algorithms have delay* $\Omega(3^{n/3})$-*time.* □

6.2 Delay of Eppstein et al. algorithm

It is worth observing how Eppstein et al. [10] developed a variation of CLIQUES tailored for sparse graphs, running in $O(n3^{d/3})$ time, where d is the degeneracy of the input graph. The degeneracy d is typically small on real-world sparse graphs. This parameter is used to define a specific ordering of the vertices which creates smaller subgraphs to process (provided that d is small), using CLIQUES.

In a worst-case scenario, however, d can be $\Omega(n)$, and Conte et al. [9] proves, by building suitable ad-hoc graphs, how the delay of Eppstein et al. algorithm is $\Omega(3^{n/6})$.

Concluding remarks

In the midst of our present work, we noticed an article [16] claiming that "The Bron-Kerbosch algorithm with vertex ordering is output-sensitive." But Conte and Versari pointed out a bug in its main lemma (Lemma 11 of [16], confirmed by the author in private communications). Therefore the question of the delay is now settled, but that of the amortized cost per solution remains open.

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