An Exact Algorithm for Maximum k-Plexes in Massive Graphs

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Abstract

The maximum k-plex, a generalization of maximum clique, is used to cope with a great number of real-world problems. The aim of this paper is to propose a novel exact k-plex algorithm that can deal with large-scaled graphs with millions of vertices and edges. Specifically, we first propose several new graph reduction methods through a careful analyzing of structures of induced subgraphs. Afterwards, we present a preprocessing method to simplify initial graphs. Additionally, we present a branch-and-bound algorithm integrating the reduction methods as well as a new dynamic vertex selection mechanism. We perform intensive experiments to evaluate our algorithm, and show that the proposed strategies are effective and our algorithm outperforms state-of-the-art algorithms, especially for real-world massive graphs.

1 Introduction

With rapid growth of the information technology, real-world graphs have been growing quickly in the past decade. A significant challenge rises: traditional algorithms fail to deal with newly emerged massive graphs. Therefore, algorithms in massive graphs (or complex networks) have received much attention due to their great importance in real-world applications. There has been a huge amount of efforts devoted to solve various problems, such as maximum cliques [Batsyn *et al.*, 2014; Cai and Lin, 2016], and graph coloring [Rossi and Ahmed, 2014]. Among those problems, the maximum clique can settle various applications [Lakhlef, 2015; Conte *et al.*, 2016].

Due to the strict restrictions of cliques, it is always difficult to model problems directly, so the k-plex was proposed and investigated [Seidman and Foster, 1978; Balasundaram *et al.*, 2011; Conte *et al.*, 2017]. Unlike cliques that restrict a vertex should connect with all other vertices, the k-plex allows every vertex has several missing connections, and thus it is more useful than the clique. Recently, the k-plex has revealed great success in the social network detection [Xiao *et al.*, 2017; Miao and Balasundaram, 2017]. To

solve the maximum k-plex, several algorithms were developed, including exact algorithms [Balasundaram et al., 2011; Moser et al., 2012] as well as heuristic methods [Zhou and Hao, 2017]. Moreover, enumeration algorithms for finding all maximal k-plexes larger than a predefined integer were also developed [Berlowitz et al., 2015; Conte et al., 2017]. In recent years, massive graphs have attracted much attention. There are many maximum clique algorithms aiming at massive graphs [Jiang et al., 2017; Cai and Lin, 2016; Rossi et al., 2014]. For example, Jiang et al. [2017] proposed an exact algorithm able to deal with very large garphs. But we have reviewed much fewer algorithms for the maximum k-plex. This is partially because the k-plex is more difficult to solve than cliques due to the relaxation. Among maximum k-plex algorithms, we highlight an exact algorithm proposed by Xiao et al. [2017] very recently, which breaks the trivial exponential bound of 2^n for k > 3 and solves graphs containing 10 thousand vertices.

However, real-world graphs have grown extremely large in recent applications. The graphs may include millions of vertices and millions of edges. Those graphs are usually sparse and the degrees follow a power-law distribution. In this paper, we propose an exact algorithm to solve the maximum k-plex problem in massive sparse graphs. The algorithm is based on the framework of the branch-and-bound search and integrated with graph reduction methods and heuristic strategies. First, three reduction methods based on the lower bound are introduced, two of which are particularly developed for sparse graphs through a deep analysis of the structure of induced subgraphs. Moreover, a reduction method using the parameter k is also discussed as a further enhancement. Combinations of these methods are employed in two types of branches, i.e., selecting or discarding a vertex, to reduce search space. Also, reduction methods are performed in the preprocessing so as to simplify initial graphs. In addition, we employ heuristic methods to implement the dynamic vertex selection with the purpose of making the remaining graph as small as possible. Intensive experiments are performed on real-world graphs from Network Data Repository online [Rossi and Ahmed, 2015] and DIMACS benchmarks. First, it shows graphs can be reduced greatly through our preprocessing method. Also, comparative results are analyzed with varying k from 2 to 5, showing that our new algorithm solves graphs up to millions of vertices efficiently and outperforms

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the existing exact algorithms proposed recently. Finally, we give some conclusions.

2 Preliminaries

Graphs discussed in this paper are simple and undirected. Formally, an undirected graph is defined as G=(V,E), where V is a set of vertices and E is a set of edges. An edge in E is associated with two vertices in V, so an unordered pair of two vertices specifies an edge. We say a vertex u is a neighbor of a vertex v if there is an edge (v,u) in E. The set of the neighbors of a vertex v in V is denoted by $N_G(v)$. Specially, we define $N_G(v)$ as an empty set if v is not in V. The degree of a vertex v is defined as the number of its neighbors, denoted as $d_G(v) = |N_G(v)|$. Given a subset $S \subseteq V$, G[S] = (V', E') is the induced subgraph in G by S if V' = S and the edge set E' consists of all the edges in E whose both vertices are in S.

Given a graph G=(V,E) and a positive integer k, a subset $S\subseteq V$ is a k-plex if the degree of each vertex in the induced subgraph G[S] is at least |S|-k. We say a k-plex is maximal if any other k-plex cannot strictly contain it. The maximum k-plex problem is to find the k-plex with the largest size of a given graph. It is easy to see a 1-plex is a clique. It is proved that finding a k-plex with a fixed size is NP-complete for any constant positive integer k [Balasundaram et al., 2011].

3 Graph Reductions

In this section, we present some new theoretical results useful for graph reductions.

3.1 Properties

The first result is straightforward.

Property 1. Given a graph G = (V, E), and a positive integer LB, a vertex v cannot be included in a k-plex larger than LB if $d_G(v) + k \le LB$.

Obviously, Property 1 is correct due to the definition of the k-plex. Such vertex v can be removed from the graph when we search a k-plex larger than LB. However, the effect of this property is limited since it basically relies on vertex degrees.

To reduce more vertices, a complex situation is discussed below. We provide some notions first. Given a graph G=(V,E), a k-plex $S\subseteq V$, and a vertex u in V, we define the notation $r_{G,S}(u)$ as

$$r_{G,S}(u) = k - |S \setminus (N_G(u) \cup \{u\})| - 1.$$
 (1)

It indicates that a maximal k-plex containing $S \cup \{u\}$ can further include at most $r_{G,S}(u)$ vertices from $V \setminus (S \cup \{u\} \cup N_G(u))$. Note that u can be either an element in S or in $V \setminus S$. The notation $c_{G,S}(u,v)$ is defined as

 $c_{G,S}(u,v) = min(d_{G'}(u) + r_{G,S}(u) + 1, |V'|) + r_{G,S}(v), (2)$ where $G'(V', E') = G[(N_G(v) \setminus S) \cup \{v\}], v$ is a vertex in V, u is a vertex in $V' \setminus \{v\}$, and function min returns the smaller value between two parameters.

Property 2. Given a graph G = (V, E) and a k-plex $S \subseteq V$, the maximum k-plex including $S \cup \{u, v\}$ is no larger than $c_{G,S}(u,v) + |S \setminus \{v\}|$, where v is a vertex in V and u is a vertex in $N_G(v) \setminus S$.

Proof. Let $G'(V',E')=G[(N_G(v)\setminus S)\cup \{v\}]$. We divide vertices from $V\setminus S$ into 2 parts: the first one is $V\setminus (V'\cup S)$, from which we can choose $r_{G,S}(v)$ vertices at most due to the restriction on non-neighbors of v; the second one is V'. After u and v are selected, due to the restriction on non-neighbors of u, at most $d_{G'}(u)-|\{v\}|+r_{G,S}(u)$ extra vertices can be selected in V' if there are enough vertices. Otherwise, all vertices in |V'| can be selected legally. So $min(d_{G'}(u)+r_{G,S}(u)+1,|V'|)$ is an upper bound of the vertices including u and v from V', and thus the the maximum k-plex including $S\cup\{u,v\}$ is no larger than $c_{G,S}(u,v)+|S\setminus\{v\}|$. □

Property 3. Given a graph G = (V, E), a k-plex $S \subseteq V$, a vertex v in S, and a vertex u in $N_G(v) \setminus S$, $S \cup \{u\}$ cannot be involved in a k-plex larger than a positive integer LB if $c_{G,S}(u,v) + |S| - 1 \le LB$.

It is easy to see Property 3 is correct according to Property 2. Furthermore, we can also apply Property 2 to make a further reduction. When we try to add a vertex v into S, we can check each vertex u in $G' = G[(N_G(v) \setminus S) \cup \{v\}]$, and remove the vertex u ($u \neq v$) satisfying Property 2 under a given lower bound LB temporarily. We use V^* to denote the remaining vertex set of G'. After that, we can count the vertices in V^* to determine whether v can be removed or added into S. Thereafter, we have the following property.

Property 4. Given a graph G=(V,E), a k-plex $S\subseteq V$ and a vertex v in $V\setminus S$, $S\cup \{v\}$ cannot be included in a k-plex larger than a positive integer LB if there exists a remaining vertex set V^* such that $|S|+|V^*|+r_{G,S}(v)\leq LB$.

Proof. It is obvious that we can choose at most $r_{G,S}(v)$ vertices from $V \setminus (S \cup V^*)$. With the case that all vertices in V^* can be selected, we obtain $|S| + |V^*| + r_{G,S}(v)$ as an upper bound of the k-plex. Thereafter, v cannot be included in a k-plex larger than LB.

To explain Property 4, we provide an example shown in Figure 1, where $S = \{s_1\}$. Then we check v, and have $G' = G[(N_G(v) \setminus S) \cup \{v\}]$ composed of u, v and w_3 . In the case k=2 and LB=4, we have $r_{G,S}(v)=1$, $r_{G,S}(u)=0$ and $d_{G'}(u)=1$, and thus $c_{G,S}(u,v)+|S|=4$ equal to LB. Therefore, u is removed from G' temporarily, and thus $V^* = \{v, w_3\}$. Then, we have $|S|+|V^*|+r_{G,S}(v)=4$, so v should be removed when searching a k-plex larger than 4.

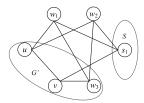


Figure 1: An example for the graph reduction

Property 4 is useful for further reduction on remaining vertices that cannot be deleted by Property 1. The sparser the graph is, the more vertices in the induced graph can be deleted. Therefore, efforts will be promoted for sparse graphs.

3.2 Graph Reduction Algorithms

With the graph reduction properties presented above, we propose several reduction algorithms. The first one, we call vertex-reduction, is based on Property 1. The detailed description of vertex-reduction is shown in Algorithm 1.

The second one employs Property 4 to reduce the graph by counting allowed vertices in induced subgraphs, so we call it subgraph-reduction, shown in Algorithm 2. Rather than vertex-reduction that reduces the graph to minimal, subgraph-reduction only checks each vertex once. When it finds a vertex that can be removed, it reduces the graph and updates vertex degrees immediately and then checks the next vertex. Subgraph-reduction runs in $O(|V| \times (|V| + |E|))$ time, so the efficiency may be increased if the graph is sparse.

Algorithm 2: function subgraph-reduction(G, S, LB, k, C)

```
Input: G = (V, E), S a k-plex in G(S \subseteq V), k the k-plex
            parameter, LB a lower bound, C the vertex set to be
            checked satisfying C \cap S = \emptyset
   Output: G the reduced graph
1 foreach vertex v in C do
        compute r_{G,S}(u) by Eq. (1), for \forall u \in (N_G(v) \setminus S) \cup \{v\};
2
         G''(V'',E'') \leftarrow G, and Q \leftarrow \emptyset;
3
        foreach u in N_{G''}(v) \setminus S do
4
             compute c_{G'',S}(u,v) by Eq. (2);
5
             if c_{G'',S}(u,v) \leq LB - |S| then add u into Q;
6
7
         while Q is not empty do
             u \leftarrow \text{pop a vertex in } Q;
8
             H \leftarrow (N_{G''}(u) \cap N_{G''}(v)) \setminus S;
             if u is in V'' then remove u from G'';
10
             foreach w in H do
11
                  update c_{G'',S}(w,v) by Eq. (2);
12
                  if c_{G'',S}(w,v) \leq LB - |S| then add w into Q;
13
             if |(N_{G''}(v) \cup \{v\}) \setminus S| + r_{G,S}(v) \le LB - |S| then
14
                  remove v from G;
15
                  break;
16
17 return G;
```

Similarly, according to Property 3, we present the reduction method called v-reduction in Algorithm 3. It can reduce vertices in the induced subgraph after a vertex \boldsymbol{v} is selected. It works similar to subgraph-reduction, but it removes vertices

from G directly, since vertices not satisfying Property 3 no longer makes efforts.

```
Algorithm 3: function v-reduction(G, S, LB, k, v)
   Input: G = (V, E), S a k-plex in G (S \subseteq V), k the k-plex
           parameter, LB a lower bound, v a vertex in S
   Output: G the reduced graph
1 compute r_{G,S}(u) by Eq. (1), for \forall u \in (N_G(v) \setminus S) \cup \{v\};
2 Q \leftarrow \emptyset;
3 foreach u in N_G(v) \setminus S do
        compute c_{G,S}(u,v) by Eq. (2);
4
       if c_{G,S}(u,v) \leq LB - |S| + 1 then add u into Q;
5
   while Q is not empty do
        u \leftarrow pop a vertex in Q;
        H \leftarrow (N_G(u) \cap N_G(v)) \setminus S;
8
        if u is in V then remove u from G;
9
        foreach w in H do
10
            update c_{G,S}(w,v) by Eq. (2);
11
            if c_{G,S}(w,v) \leq LB - |S| + 1 then add w into Q;
12
13 return G;
```

Beside those methods, a strategy to deal with vertices violating the k-plex constraint is introduced here. If we aim at finding a k-plex including S, two kinds of vertices should be removed: 1) if there is a vertex u in S satisfying $|S\setminus N_G(u)|+1=k$, then all vertices in $V\setminus (N_G(u)\cup S)$ should be removed; 2) if there is a vertex u in $V\setminus S$ satisfying $|S\setminus N_G(u)|+1>k$, then u should be removed. Algorithm 4 formally describes how to reduce the graph using the strategy above.

```
Algorithm 4: function k-reduction(G, S, k)

Input: G = (V, E), S a k-plex in G (S \subseteq V), k the k-plex parameter, LB a lower bound

Output: G the reduced graph

1 foreach u in S satisfying |S \setminus N_G(u)| + 1 = k do

2 \lfloor remove any w in V \setminus (N_G(u) \cup S) from G;

3 foreach u in V \setminus S satisfying |S \setminus N_G(u)| + 1 > k do

4 \lfloor remove u from G;

5 return G;
```

4 Preprocessing

As we know, massive graphs may have numerous nodes, which may cause the execution of an algorithm out of memory. So preprocessing is a necessary step, which can usually reduce the input graphs to a reasonable size. The more nodes in the original graphs are removed, the less branches are exploited in the searching algorithm. Algorithm 5 shows the procedure of the preprocessing step. We employ a heuristic method to obtain a lower bound, which is an extension of the method in [Jiang *et al.*, 2017]. Based on the lower bound, vertex-reduction can reduce graphs efficiently due to its linear computation time for checking a vertex. However, the effectiveness, unfortunately, is not satisfactory, if the heuristic method fails to find a lower bound close to the optimal value.

It is not surprise that a heuristic sometimes finds a solution far from the best one. As a result, subgraph-reduction is also used in preprocessing for a further reduction. With the same lower bound, it can remove more vertices, though it needs higher computation complexity. Two reduction methods are preformed alternatively until no vertex can be removed. During iterations, the lower bound is updated before each execution of the vertex-reduction, since the heuristic may find new lower bounds when the graph is reduced.

Algorithm 5: function reduce-graph(G,k)

```
Input: G = (V, E), k the k-plex parameter.
   Output: LB a lower bound, and G the reduced graph
1 m \leftarrow |V| + 1, LB \leftarrow 0;
2 while |V| < m do
        m \leftarrow |V|;
3
        if LB-heuristic (G, k) > LB then
         LB \leftarrow LB-heuristic(G, k);
5
        G \leftarrow \text{vertex-reduction}(G, LB, k);
        G \leftarrow \text{subgraph-reduction}(G, \emptyset, LB, k, V);
8 return (LB, G):
9 function LB-heuristic(G, k);
10 for i \leftarrow 1 to |V| do
        v \leftarrow the vertex with the smallest d_G(v) in V;
11
        if d_G(v) + k \ge |V| then break;
12
        remove v from G;
14 return |V|;
```

5 The Branch-and-Bound Algorithm

In this section, we present our branch-and-bound algorithm. The entire framework of the algorithm is indicated in Algorithm 6.

The algorithm starts with the preprocessing algorithm to reduce the input graphs, and then it performs the branch-and-bound algorithm. Search tree nodes are expanded recursively by the function BB. It returns the size of S if there is no vertex left, or the best lower bound found in its branch. There are some crucial components in the function BB to be explained.

The first one is the reduction methods. The branching rule here is binary, *i.e.*, selecting a vertex and adding it into the target k-plex or discarding the vertex. After that, reductions are performed to simplify the remaining graph. Two combinations of reduction methods are employed after selecting and discarding, respectively.

In a selecting branch, we take three reduction methods. They are performed in the ordering of k-reduction, vertex-reduction and v-reduction. The reason of taking this ordering is under the following observation: vertex-reduction cannot make efforts if performed first as it is irrelevant with S, but k-reduction can delete all conflict vertices directly. Therefore k-reduction is the first, and then vertex-reduction and v-reduction. Different from the preprocessing step, v-reduction is employed instead of the subgraph-reduction due to the efficiency. In a discarding branch, vertex-reduction and subgraph-reduction are used. Obviously, k-reduction

Algorithm 6: The Branch-and-Bound Algorithm (BnB)

```
Input: G = (V, E), k the k-plex parameter.
    Output: the size of the maximum k-plex
 1 S \leftarrow \emptyset;
 2 (LB, G) \leftarrow \text{reduce-graph}(G, k);
 \mathbf{3} \stackrel{.}{L}B \leftarrow \mathbf{BB}(G, S, k, LB);
 4 return LB;
 5 function BB(G, S, k, LB);
 6 if V \setminus S is empty then return |S|;
 7 select a vertex v in V \setminus S, and S \leftarrow S \cup \{v\};
 8 store G as G_r;
 9 G \leftarrow \text{k-reduction}(G, S, k);
10 G \leftarrow \text{vertex-reduction}(G, LB, k);
11 G \leftarrow \text{v-reduction}(G, S, LB, k, v);
12 if |\{u|u \text{ in } V \setminus S \text{ and } d_G(u) + k > LB\}| + |S| > LB then
         lower \leftarrow BB(G, S, k, LB);
         if LB < lower then LB \leftarrow lower;
15 restore G_r to G;
16 S \leftarrow S \setminus \{v\}, and remove v from G;
17 G \leftarrow \text{vertex-reduction}(G, LB, k);
18 G \leftarrow \text{subgraph-reduction}(G, S, LB, k, V \setminus S);
19 if |\{u|u \text{ in } V \setminus S \text{ and } d_G(u) + k > LB\}| + |S| > LB then
20
         lower \leftarrow BB(G, S, k, LB);
         if LB < lower then LB \leftarrow lower;
21
22 return LB;
```

could not work in this branch, since S does not change after discarding a vertex so as to $r_{G,S}(v)$. However, the fact that degrees of discarded vertex's neighbors decrease makes vertex-reduction work well. To further enhance the reduction, subgraph-reduction is performed after vertex-reduction. Different from reducing graphs iteratively, all reductions are performed only once after selecting or discarding.

The second component discussed is the computation of upper bounds. After reductions, the algorithm should compute upper bounds to determine selecting a new vertex or trying another branch. Since the graph is reduced deeply, a simple but efficient method that counts the number of remaining vertices is developed for the upper bound, but the actual number of remaining vertices is not used as the bound directly in our algorithm. This is because the reduction methods during branch-and-bound searching are not performed iteratively, as we described above, and thus some vertices in a reduced graph may still satisfy the vertex-reduction rule. Therefore, we have to scan the vertices again to calculate the bound. To speed up this task, we employ a lazy method to determine whether the current branch has an upper bound bigger than the lower bound, rather than the method computing the exact value of the upper bound. This is achieved by counting remaining vertex whose degree exceeds the lower bound plus k. If it finds more than LB such vertices, the algorithm executes the function BB; otherwise, it tries the other branch or backtracks to the previous level.

The last one is the vertex selection heuristic, which specifies how to select a vertex on each node of the search tree. As we know the ordering of vertex selection may impact greatly on the search tree size. To make the selection mechanism effective, a dynamic vertex ordering strategy is employed in our algorithm. It selects a vertex according to the current branch and the remaining graph. The purpose of the dynamic vertex ordering strategy is to delete vertices as many as possible in the reduction after a selection. As the reduction algorithms are mainly based on vertex degrees, it is clear that selecting the vertex with the maximum degree from the remaining graph is a basic strategy that can decrease vertex degrees in full measure. Furthermore, our reduction method k-reduction is according to the value $r_{G,S}(v)$, the maximal allowed nonneighbors from the remaining graph, and thus if a vertex v with $r_{G,S}(v)=0$ is selected, all its non-neighbors can be removed. So a compound measurement is proposed, where $r_{G,S}(v)=0$ is the primary condition while the degree is the secondary condition used to break ties.

With the discussion above, two ordering rules are considered in our algorithm.

Rule 1. Select the vertex v with the maximum $d_{G[V\setminus S]}(v)$. **Rule 2.** Select the vertex v with the maximum $d_{G[V\setminus S]}(v)$ from $V_0 = \{v|k = |S\setminus N_G(v)|+1\}$ if V_0 is not empty; Select the vertex according to Rule 1 otherwise.

If there is more than one vertex as candidates, select the first one in lexicographical order. We call the BnB with Rule 1 as the vertex selection strategy BnBd, and the BnB with Rule 2 BnBk. In the following section, we will show comparative results of the exact algorithm we proposed.

6 Experiments

In this section, we performed intensive experiments to evaluate the algorithm we proposed. Two benchmarks were considered. One is selected instances from DIMACS benchmark, tested in [Xiao *et al.*, 2017]. The other is real-world massive graphs, originally from the Network Data Repository online [Rossi and Ahmed, 2015], where 139 massive sparse graphs are selected ¹. Those benchmarks are popular on evaluating various massive graph algorithms [Rossi and Ahmed, 2014; Cai, 2015; Lin *et al.*, 2017].

Our algorithms were implemented in C++ language and compiled by g++ version 4.7 with -O3 option. Benchmarks were solved parallel on a workstation with an Intel Xeon E5-1650v4 (3.6GHz) CPU, and 16GB RAM, running Ubuntu Linux 16.04, with the cutoff time 10000s. There are several exact algorithms proposed for the maximum *k*-plex, such as IPBC [Balasundaram *et al.*, 2011], GuidedBranching [Moser *et al.*, 2012] and BS [Xiao *et al.*, 2017].

We first compare our algorithms with the three algorithms on DIMACS benchmarks fixing k=2. Table 1 indicates runtimes, where "—" means time out or out of memory. BS and our algorithms are computed by our workstation whereas data of others come from [Xiao $et\ al.$, 2017], which were computed by slower machines. From Table 1, it can be seen that IPBC is faster than others on hamming and johson instances, but those instances are relatively small. BS, BnBd and BnBk are good at solving instances with more than 200 vertices, such as c-fat500, brock and p_hat instances. This result is compatible with [Xiao $et\ al.$, 2017], in which it is

instance	IPBC	GuideBranching	BS	BnBd	BnBk
c-fat200-1	148.9	1.1	0.112	0.008	0.004
c-fat200-2	19.1	3.53	0.132	0.004	0
c-fat200-5	2.1	22.44	0.158	0.036	0.028
c-fat500-1	1356	11.01	0.408	0.008	0.016
c-fat500-2	605.3	50.21	0.536	0.068	0.024
c-fat500-5	141.5	350.6	0.352	0.132	0.108
c-fat500-10	76.5	1547	0.47	0.388	0.424
hamming6-2	0	1.77	32.59	110.9	13.82
hamming6-4	0.3	0.24	0.066	0.024	0.02
hamming8-4	8115	_	_	_	_
johnson8-2-4	0	0.02	0.012	0.008	0.004
johnson8-4-4	4.4	40.7	228.9	20.92	8.836
MANN_a9	0	0.09	4.218	16.66	11.26
brock200_2	_	606.1	409.8	268.7	55
brock200_4	_	9691	_	—	4121
p_hat300-1	_	502.5	31.84	12.91	9.368
p_hat700-1	_	_	_	5527	1970

Table 1: Runtimes of DIMACS benchmarks in seconds

reported that BS outperforms others on large instances (e.g. social networks). Among BS, BnBd and BnBk, BnBk performs best on most instances.

Consequently, we compare our algorithms with BS on massive graphs in the followings. It is noted that benchmarks from the Network Data Repository online includes many graphs with millions of vertices, which are much larger than the social networks (10 thousand vertices at maximum) used in [Xiao *et al.*, 2017].

Before analyzing runtimes on massive graphs, we show the effect made by our subgraph-reduction. We count the remaining vertices in the preprocessing step compared to the method with vertex-reduction only (removing subgraph-reduction from function reduce-graph). Detailed results of the facebook networks with k=2 and k=5 are listed in Table 2. It shows that using subgraph-reduction can remove much more vertices compared with using vertex-reduction only. In fact, the preprocessing function reduce-graph can reduce to 0.15% scale of the original graphs on average of 139 instances when k=2, about 1/20 size of the vertex-reduction results. Surprisingly, it can reduce 68 out of 139 graphs to empty and thus achieved the maximum solutions without the branch-and-bound search. Moreover, similar results can be achieved when k=3,4,5.

Facebook	original	k =	2	k = 5		
networks	size	vertex	reduce-	vertex	reduce-	
		reduction	graph	reduction	graph	
socfb-A-anon	3097165	390144	2239	298641	819	
socfb-B-anon	2937612	489129	37564	432416	34767	
socfb-Berkeley13	22900	12322	831	10136	292	
socfb-CMU	6621	3616	289	3209	54	
socfb-Duke14	9885	7090	685	6536	399	
socfb-Indiana	29732	19088	2316	15859	976	
socfb-MIT	6402	3871	280	3578	60	
socfb-OR	63392	10506	294	7599	180	
socfb-Penn94	41536	20353	0	17080	59	
socfb-Stanford3	11586	7759	3021	7109	1388	
socfb-Texas84	36364	16086	452	9918	111	
socfb-uci-uni	58790782	499697	0	26159	0	
socfb-UCLA	20453	7631	66	6877	65	
socfb-UConn	17206	4574	63	899	63	
socfb-UCSB37	14917	1050	148	814	0	
socfb-UF	35111	14613	1527	8526	714	
socfb-Uillinois	30795	24838	6629	21421	3938	
socfb-Wisconsin87	23831	12709	266	10431	132	

Table 2: Reduction results in preprocessing

¹ available at: http://lcs.ios.ac.cn/~caisw/Resource/realworld%20 graphs.tar.gz

Table 3 indicates statistic results grouped by families with k=2,3,4,5, respectively, by counting the successfully solved instances of three algorithms: BS, BnBk and BnBd. We use family abbreviations in the table, and full names are biological networks, collaboration networks, Facebook networks, interaction networks, infrastructure networks, amazon recommend networks, retweet networks, scientific computation networks, social networks, technological networks, temporal reachability networks and web link networks.

Families	k = 2		k = 3		k = 4			k = 5				
(#instance)	BS	BnBd	BnBk	BS	BnBd	BnBk	BS	BnBd	BnBk	BS	BnBd	BnBk
bio (4)	4	4	4	4	4	4	4	4	4	4	4	4
col (13)	12	13	13	12	13	13	12	13	13	12	13	13
fb (18)	3	15	16	2	14	14	2	14	14	1	15	15
inf (4)	2	4	4	2	4	4	2	2	2	2	1	1
int (9)	8	9	9	7	9	9	7	9	9	6	9	9
rec (1)	0	1	1	0	1	1	0	0	0	0	0	0
ret (3)	2	3	3	2	3	3	2	3	3	2	3	3
sci (8)	2	5	5	2	2	2	1	1	1	2	1	1
soc (23)	10	16	17	9	17	18	7	16	16	6	16	16
tec (7)	7	5	6	6	6	7	4	6	6	4	7	7
tem (37)	37	37	37	37	37	37	37	36	36	37	35	35
web (12)	12	12	12	11	12	12	11	11	11	11	10	10
total (139)	99	124	127	94	122	124	89	115	115	87	114	114

Table 3: Comparison results of instances solved

From Table 3, it is easy to see that BnBk solves the most amount of instances among the three algorithms, much more than BS. For each k, BnBk also performs best. Indeed, most instances solved by BS can also be solved by BnBk successfully, except only 7 instances partly due to unsatisfactory lower bounds obtained. In contrast, BS fails to solve some instances that can be solved by BnBk very easily, or saying within 100s. Also, BnBd reveals better performance than B-S, but solves a bit fewer instances than BnBk. Moreover, the number of solved instances decreases with k growing for each algorithm, because most instances become harder to solve.

We then analyze runtimes. We only show solved instances with more than 50000 vertices, and thus select out 44 instances. Table 4 indicates runtimes with k=2,5, respectively, where "—" means time out or out of memory. It shows that BnBd and BnBk are much faster than BS on most instances; BnBd is slightly better than BnBk if instances can be solved in a short time; but BnBk shows a distinct advantage when instances are hard to solve, *i.e.*, larger than 100s.

We also compare our BnBk with BS on the reduced instances. Specifically, we reduce benchmark instances with our preprocessing method, and then solve the reduced instances by both algorithms. Runtimes excluding preprocessing time are listed in Table 5, where instances solved within 10s and unsolved instances are omitted. From it, we can see BS is greatly improved when combining our preprocessing method as more instances can be solved than the original BS. Among 27 instances, BnBk performs better on 18 instances whereas BS on 9 instances. In total, BnBk is also better than BS when BS is associated with our preprocessing method.

7 Conclusions

To solve maximum k-plexes exactly, we have proposed a branch-and-bound algorithm for massive sparse graphs.

	1	-, -			_,	
		k=2			k=5	
instance	BS	BnBd	BnBk	BS	BnBd	BnBk
ca-citeseer	0.692	0.424	0.524	0.512	0.496	0.448
ca-coauthors-dblp	12.41	9.064	9.6	11.34	10.65	12.75
ca-dblp-2010	0.448	0.436	0.352	0.348	0.38	0.428
ca-dblp-2012	1.245	0.656	0.52	0.853	0.672	0.756
ca-hollywood-2009	_	31.59	38.86	_	34.14	37.83
ca-MathSciNet	0.965	0.404	0.68	0.823	0.592	0.688
socfb-A-anon	_	153.2	154.9	—	207.4	180.9
socfb-B-anon	_	840.1	1012	—	_	_
socfb-OR	_	26.004	8.036	—	7.7	5.276
socfb-uci-uni	_	86.14	112.9	—	82.99	90.08
inf-roadNet-CA	8.056	1.824	2.012	8.401	_	_
inf-roadNet-PA	_	0.988	1.044	—	_	_
inf-road-usa	_	22.2	24.42	—	_	_
ia-wiki-Talk	_	18.09	11.85	_	2779	938.5
rec-amazon	_	0.112	0.072	_	_	_
rt-retweet-crawl	_	2.536	2.664	l —	2.092	2.208
sc-nasasrb	_	332.1	340.4	_	_	_
sc-pkustk11	_	3.528	3.776	_	3838	3705
sc-pkustk13	_	275.8	313.8	_	_	_
sc-shipsec1	600.1	1.428	1.388	675.2	_	_
sc-shipsec5	1144	30.77	32.08	1064	_	_
soc-brightkite	1.185	0.136	0.244	345.5	0.168	0.156
soc-delicious	43.34	1.192	1.204	l —	1.236	1.276
soc-douban	83.29	0.352	0.572	78.09	0.376	0.384
soc-flixster	<u> </u>	_	465.7	l —	_	_
soc-FourSquare	_	_	_	_	411.5	395.6
soc-gowalla	_	24.99	11.96	l —	279.1	106.7
soc-lastfm	_	165.5	108.4	l —	7039	1896
soc-livejouranl	46.44	17.43	20.01	46.23	22.21	19.91
soc-LiveMocha	_	313.3	220.1	l —	_	_
soc-pokec	_	74.54	75.73	l —	65.99	63.64
soc-slashdot	96.19	252.6	17.05	l —	440.1	266.8
soc-twitter-follows	4683	4.52	4.68	l —	25.01	21.76
soc-youtube	_	11.01	11.65	l —	212.6	109.9
soc-youtube-snap	_	17.61	15.88	l —	612.1	322.7
tech-as-skitter	3845	_	1426	l —	170.3	77.22
tech-p2p-gnutella	67.23	0.084	0.108	352.2	0.1	0.092
tech-RL-caida	68.38	1.524	1.592	l —	1.224	1.352
scc_retweet-crawl	0.109	0.14	0.076	0.098	0.084	0.128
web-arabic-2005	0.579	0.872	0.82	0.68	0.896	0.72
web-it-2004	76.59	8.576	9.444	77.89	2817	2859
web-sk-2005	0.22	0.256	0.18	0.325	16.98	17.14
web-uk-2005	59.78	19.19	19.88	60.86	_	_
web-wikipedia2009	914.2	43.29	41.7	l —	_	_

Table 4: Runtimes of massive graphs in seconds

instance	k:	=2	instance	k=2		
ilistance	BS	BnBk	ilistance	BS	BnBk	
socfb-A-anon	1012	28.99	soc-flixster	91.80	455.0	
socfb-B-anon	_	755.4	soc-gowalla	115.0	7.328	
socfb-Duke14	958.1	841.2	soc-lastfm	2663	90.68	
socfb-Indiana	565.1	417.4	soc-LiveMocha	_	179.2	
socfb-MIT	10.72	36.52	soc-slashdot	6.162	16.50	
socfb-Stanford3	6704	179.4	soc-twitter-follows	2073	1.504	
socfb-Texas84	108.3	_	soc-youtube	158.4	3.252	
socfb-UF	584.4	_	soc-youtube-snap	782.7	5.516	
socfb-Uillinois	2083	429.6	tech-as-skitter	483.3	1414	
ia-enron-large	21.88	0.920	tech-WHOIS	50.56	_	
ia-wiki-Talk	136.7	10.08	scc_fb-forum	0.180	394.0	
sc-nasasrb	l —	338.0	scc_twitter-copen	0.336	94.26	
sc-pkustk13	—	272.9	web-wikipedia2009	764.8	35.21	
sc-shipsec5	148.8	29.97				

Table 5: Searching time in seconds after the proposed preprocessing step

Three main components have been described in detail. Special graph reduction methods for sparse graphs have been proposed to enhance the efficiency of the algorithm, where analyzing of induced subgraph structures makes significant effects. Moreover, a dynamic vertex selection mechanism has been proposed to speed up searching. Additionally, a lazy

computation of the upper bound has been introduced. Besides, a preprocessing method with our reduction strategies has been introduced in the algorithm, which may always reduce massive graphs to tractable sizes. We have shown our reduction strategies work effectively through computational experiments, and our algorithm outperforms the state-of-theart exact algorithms. It is also worth to mention that the graph with more than 50 million vertices can be solved in a very short time by our algorithm. In the future, parallelization of the proposed methods and algorithms for maximum weighted k-plex problems are valuable research directions.

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