# Frequency-driven tabu search for the maximum s-plex problem

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#### Abstract

The maximum s-plex problem is an important model for social network analysis and other studies. In this study, we present an effective frequency-driven multineighborhood tabu search algorithm (FD-TS) to solve the problem on very large networks. The proposed FD-TS algorithm relies on two transformation operators (Add and Swap) to locate high-quality solutions, and a frequency-driven perturbation operator (Press) to escape and search beyond the identified local optimum traps. We report computational results for 47 massive real-life (sparse) graphs from the SNAP Collection and the 10th DIMACS Challenge, as well as 52 (dense) graphs from the 2nd DIMACS Challenge (results for 48 more graphs are also provided in the Appendix). We demonstrate the effectiveness of our approach by presenting comparisons with the current best-performing algorithms.

**Keywords**: Clique relaxation; Heuristic; Massive network; s-plex.

#### 1 Introduction

Given a simple undirected graph G = (V, E) with a set of vertices V and a set of edges E, let N(v) denote the set of vertices adjacent to v in G. Then, an s-plex for a given integer  $s \geq 1$  ( $s \in \mathbb{Z}^+$ ) is a subset of vertices  $C \subseteq V$  that satisfies the following condition:  $\forall v \in C$ ,  $|N(v) \cap C| \geq |C| - s$ . Thus, each vertex of an s-plex C must be adjacent to at least |C| - s vertices in the subgraph  $G[C] = (C, E \cap (C \times C))$  induced by C.

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The maximum s-plex problem involves finding, for a fixed value of s, an s-plex of maximum cardinality among all possible s-plexes of a given graph. As indicated in [3], the maximum s-plex problem can be formulated as a binary linear program as follows:

$$\max \qquad \omega_s(G) = \sum_{i \in V} x_i$$
s.t. 
$$\sum_{j \in V \setminus (N(i) \cup \{i\})} x_j \le (s-1)x_i + \bar{d}_i(1-x_i), \forall i \in V,$$

$$x_i \in \{0,1\}, \forall i \in V$$

$$(1)$$

where  $x_i$  is the binary variable associated with vertex i, such that  $x_i = 1$  if vertex i is in an s-splex,  $x_i = 0$  otherwise. Also,  $\bar{d}_i = |V \setminus N(i)| - 1$  denotes the degree of vertex i in the complement graph  $\bar{G} = (V, \bar{E})$ . Note that  $i \notin N(i)$  by definition.

The s-plex concept was first introduced for graph-theoretic social network studies [27]. The decision version of the maximum s-plex problem with any fixed positive integer s is known to be NP-complete [3]. When s equals 1, the maximum s-plex problem reduces to the popular maximum clique problem, the decision version of which was among Karp's 21 NP-complete problems [14]. The maximum s-plex problem is often referred to as a clique relaxation model [23, 24]. Other clique relaxation models include s-defective clique [37], quasi-clique [8,21,22], and k-club [7], which are defined by relaxing the edge number, the edge density, and the pairwise distance of vertices in an induced subgraph, respectively. In addition to studies of social networks, the maximum s-plex problem has also been investigated in other contexts [5, 6, 10]. For instance, an interesting application of the maximum s-plex model was described in [6], where the maximum s-plex algorithm of [29] was used to find profitable diversified portfolios on the stock market.

Similar to a clique, an s-plex C has the heredity property, which means that every subset of vertices  $C' \subset C$  remains an s-plex, i.e., the subgraph induced by C' always has the property of an s-plex [29]. The most successful combinatorial algorithms for s-plex essentially rely on the heredity property and a polynomial feasibility verification procedure. For example, a powerful exact algorithmic framework was introduced in [29] for detecting optimal hereditary structures (s-plex and s-defective clique), which is based on the maximum clique algorithm proposed in [20]. This algorithm performed well on the maximum s-plex problem for graphs in the 2nd DIMACS Challenge and popular large-scale social networks. Other exact algorithms for the s-plex problem include the following. A branch-and-cut algorithm was introduced in [3] based on a polyhedral study of the s-plex problem. Two branch-and-bound algorithms were presented in [16], which are based on popular exact algorithms for the maximum clique problem [9,20]. In [19], exact combinatorial algorithms

were investigated using methods from parameterized algorithmics. Finally, a parallel algorithm for listing all the maximal s-plexes was introduced in [36].

However, given the computational complexity of the maximum s-plex problem, any exact algorithm is expected to require an exponential computational time to determine the optimal solution in the general case. Thus, it is useful to investigate heuristic approaches, which aim to provide satisfactory solutions within an acceptable time frame, but without a provable optimal guarantee for the solutions obtained. However, our literature review found only two heuristics for the maximum s-plex problem [12,17], which are based on the general GRASP method [26]. This situation contrasts sharply with the huge body of heuristics for the conventional maximum clique problem [34] and other clique relaxation problems [24]. We note that exact and heuristic approaches may complement each other, and together they can enlarge the classes of problem instances that can be solved effectively. Moreover, they can even be combined within a hybrid approach, as exemplified in [17] where the GRASP heuristic was used to enhance the exact algorithm proposed in [3] to solve very large social network instances.

In this study, we aim to partially fill the gap in terms of heuristic methods for solving the maximum s-plex problem by introducing an effective heuristic approach. The main contributions of this study can be summarized as follows.

- From an algorithmic perspective, this is the first study to employ the tabu search metaheuristic [11] to solve the maximum s-plex problem (Section 2). Thus, the proposed frequency-driven tabu search algorithm (FD-TS) integrates several original components. First, FD-TS jointly employs three dedicated move operators called Add, Swap, and Press, two of which (Swap and Press) are applied for the first time to the maximum s-plex problem. Second, we introduce a frequency-based mechanism for perturbation and constructing initial solutions, which is proven to be more effective than a random mechanism. We also apply a peeling procedure to dynamically reduce the graph with the best identified lower bound. Finally, specific design decisions are made in order to handle very large networks with thousands and even millions of vertices.
- From a computational perspective, our experimental results indicate that the proposed algorithm performs very well with both sparse and dense graphs (Section 4). For 47 very large networks from the SNAP collection and the 10th DIMACS Challenge benchmark set, our algorithm successfully obtained or improved the best-known results from previous studies for s=2,3,4,5. Our algorithm even proved the optimality of many instances for the first time using the peeling procedure. For 52 dense graphs from the collection used in the 2nd DIMACS Challenge, our algorithm also obtained or improved the best-known results for s=2,3,4,5. To comprehensively assess the performance of our algorithm, we compared FD-TS with several

cutting edge algorithms, including the commercial CPLEX solver (version 12.6.1). Results of 48 additional graphs for the s-plex problem are also presented for the first time in the Appendix.

The remainder of this paper is organized as follows. Section 2 presents the FD-TS algorithm. Section 3 discusses the implementation and complexity issues related to FD-TS. Section 4 presents the computational results obtained on benchmark instances and provides comparisons with state-of-the-art algorithms. In the final section, we give our conclusions and discuss future research.

# 2 FD-TS algorithm for the maximum s-plex problem

## 2.1 General procedure

The general scheme of the proposed FD-TS algorithm is shown in Algorithm 1. FD-TS starts from an initial feasible solution (s-plex) built using the Init\_Solution() procedure (Section 2.4), before entering the main multineighborhood local search procedure, Freq\_Tabu\_Search(), to improve the initial solution (Section 2.5). A vector freq, which records the number of times each vertex is moved in the last round of the Freq\_Tabu\_Search() procedure, is initialized as a null vector (Algorithm 1, line 3). This vector is used by the Init\_Solution() procedure as well as the perturbation method explained in Section 2.5.3. If the solution returned by tabu search is better than the current best solution  $C^*$ ,  $C^*$  is updated (Algorithm 1, lines 7–8). The new lower bound  $|C^*|$  is then given to the Peel() procedure (Section 2.6) to reduce the current graph (Algorithm 1, line 9). If Peel() returns a reduced subgraph with fewer vertices than  $|C^*|$ , then  $C^*$  must be an optimal solution and the overall algorithm stops. Otherwise, the algorithm enters a new round of search to build a new starting solution with  $Init\_Solution()$ , before improving the new starting solution with Freq\_Tabu\_Search() and reducing the graph with Peel() if this is possible. The algorithm continues until a given stopping condition (e.g., a cut-off time limit) is met.

# 2.2 Preliminary definitions

Given G = (V, E),  $s \in \mathbb{Z}^+$ , let  $C \subseteq V$  be a subset of vertices and N(v) the set of vertices adjacent to v. The following definitions are provided, which are useful for the description of our algorithm.

We say that C is a (feasible) solution or an s-plex if  $\forall v \in C, |N(v) \cap C| \geq$ 

# Algorithm 1: Main framework of Frequency Driven Tabu Search

**Input**: Problem instance (G, s), predefined sample size q, maximum allowed iterations in tabu search L.

Output: The largest s-plex ever found

```
1 begin
 \mathbf{2}
       C^* \leftarrow \emptyset;
                                          /* the best solution found so far */
       freq(v) \leftarrow 0 \text{ for all } v \in V; /* frequency count of vertex moves */
 3
       while the stopping condition is not met do
 4
           \{C, freq\} \leftarrow Init\_Solution(G, s, freq, q);
                                                                              /* §2.4 */
 5
           \{C, freq\} \leftarrow Freq\_Tabu\_Search(G, s, C, freq, L);
                                                                              /* §2.5 */
 6
           if |C| > |C^*| then
               C^* \leftarrow C;
 8
               G \leftarrow Peel(G, s, |C^*|);
                                                                              /* §2.6 */
 9
               if |V| < |C^*| then
10
                   return C^*;
                                          /* return the best solution found */
11
12 end
13 return C^*
```

|C|-s; otherwise, C is an infeasible solution (i.e.,  $\exists v \in C, |N(v) \cap C| < |C|-s$ ). For a vertex  $v \in C$ , we say that v is saturated (first introduced in [29]) if  $|N(v) \cap C| = |C|-s$ . If  $|N(v) \cap C| < |C|-s$ , v is deficient. Obviously, whenever a deficient vertex exists in C, C is an infeasible solution. The saturated set S of set C is defined as the set of all saturated vertices in C, i.e.,  $S = \{v \in C : |N(v) \cap C| = |C|-s\}$ . We can see that if C is a 1-plex (i.e., a clique), then all the vertices in C are saturated. The search space C0 of C1 includes all C2 - C3. For brevity, we also use C4 to denote the set of vertices in C5 with at least one adjacent vertex in C5, C6.

In Figures 1 and 2, we provide examples of S and N(C), as well as their uses in the definitions of the Add and Swap operators introduced in the next section.

Finally, the quality of any candidate solution (s-plex)  $C \in \Omega$  is evaluated by its cardinality |C|. Thus, given two candidate solutions C' and C, C' is better than C if |C'| > |C|.

#### 2.3 Move operators

Our FD-TS algorithm explores the search space  $\Omega$  by jointly applying three move (or transformation) operators, Add, Swap, and Press, to generate new solutions in  $\Omega$  from the current solution (or s-plex). If we let C be the incumbent solution, then each move operator transfers one vertex  $v \in N(C)$  inside

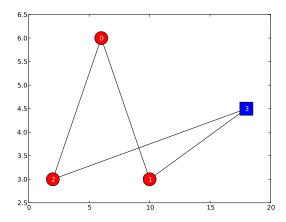


Fig. 1. Suppose s=2,  $C=\{0,1,2\}$  is the incumbent solution,  $S=\{1,2\}$  is the saturated set of C, then  $M_1=\{3\}$  and  $C\cup\{3\}$  is an extended s-plex.

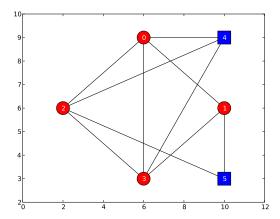


Fig. 2. Suppose s=2,  $C=\{0,1,2,3\}$  is the incumbent solution,  $S=\{1,2\}$  is the saturated set of C,  $N(C)=\{4,5\}$ . Then for the Swap operator, we get  $A=\{4\}$  and exchangeable pair <4,1>;  $B=\{5\}$  and exchangeable pairs <5,0> and <5,3>.

C and eliminates zero, one, or more vertices from C to keep C feasible. If we let OP be a move operator, then we use  $C' \leftarrow C \oplus OP(v, X)$  to denote the new (neighboring) solution obtained by applying OP to C (X represents the subset of vertices eliminated from C, which can possibly be empty). The details of these operators are described as follows. For simplicity, when the subset of eliminated vertices X is empty or a singleton, the set notation of X is ignored.

(1) Add(v): This operator extends the incumbent solution C by including a new vertex from N(C). Clearly, each application of this operator will increase the cardinality of the solution by one, which always leads to a better solution. However, we must take special care to ensure that the extended solution remains an s-plex. Thus, we identify the following

vertex subset  $M_1 \subseteq N(C)$ , which has the required feasibility property (first introduced in [29]):

$$M_1 = \{ v \in N(C) : |N(v) \cap C| \ge |C| - s + 1, S \setminus N(v) = \emptyset \}$$
 (2)

By definition (2), if a vertex v in N(C) is adjacent to at least |C|-s+1 vertices in C and adjacent to all the saturated vertices of C, then adding v to C yields a new solution C', the size of which is increased by one (see Fig. 1 for an example).

The set of neighboring solutions of C induced by Add(v) is then given by:

$$\mathcal{N}_{Add} = \{ C' : C' \leftarrow C \oplus Add(v), v \in M_1 \}$$
 (3)

The Add operator is used by the search algorithm to improve the quality of the incumbent solution.

- (2) Swap(v, u): This operator exchanges a vertex  $v \in N(C)$  with another vertex u in the incumbent solution C ( $u \in C$ ), while keeping the quality of the solution unchanged. Similar to Add, to ensure the feasibility of the transformed solutions, we need to identify the set of suitable candidate pairs  $\langle v, u \rangle \in N(C) \times C$ . Considering the definition of s-plex, a pair of vertices  $\langle v, u \rangle$  is eligible for exchange only if it satisfies one of the following two conditions.
  - First, v is adjacent to at least |C| s vertices in C and u is the unique saturated vertex that is not adjacent to v (i.e.,  $S \setminus N(v) = \{u\}$ ).
  - Second, v is adjacent to exactly |C| s vertices in C and these |C| s vertices must include all the saturated vertices (i.e.,  $S \setminus N(v) = \emptyset$ ), and u is an arbitrary vertex from  $C \setminus N(v)$ .

We use sets A and B to denote the candidate sets of v that satisfy the two conditions above, respectively, (see Fig. 2 for an example of these two types of vertices):

$$A = \{v \in N(C) : |N(v) \cap C| \ge |C| - s, |S \setminus N(v)| = 1\}$$

$$B = \{v \in N(C) : |N(v) \cap C| = |C| - s, S \setminus N(v) = \emptyset\}$$

$$(4)$$

The set of neighboring solutions induced by Swap(v, u) is then given by:

$$\mathcal{N}_{Swap} = \{ C' : C' \leftarrow C \oplus Swap(v, u), (v \in A, u \in S \setminus N(v)) \lor (v \in B, u \in C \setminus N(v)) \}$$
(5)

If we let  $M_2 = A \cup B$ , a practical method for generating a suitable pair  $\langle v, u \rangle$  is to first build the set  $M_2$ , then pick a vertex v from  $M_2$ , and finally determine an appropriate vertex u from  $S \setminus N(v)$  or  $C \setminus N(v)$ . The Swap operator is used by the search algorithm to visit neighboring

- solutions of equal quality, so a transition with Swap is also called a sidewalk move.
- (3) Press(v, X): If we let  $M_3 = N(C) \setminus (M_1 \cup M_2)$ , this operator adds one vertex  $v \in M_3$  to C and then eliminates two or more vertices from  $C \setminus N(v)$  so the s-plex structure of the new solution is maintained. Obviously, given a vertex  $v \notin M_1$ , the set  $C' = C \cup \{v\}$  is an infeasible solution because v or the vertices in  $S \setminus N(v)$  become deficient (see Section 2.2) in G[C']. Therefore, to restore the feasibility of the transformed solution, this operator iteratively eliminates vertices from  $(C' \setminus \{v\}) \setminus N(v)$  (i.e.,  $C \setminus N(v)$ ) until the solution becomes an s-plex. Preference is given to the deficient vertices in  $C \setminus N(v)$  and if no deficient vertex exists in  $C \setminus N(v)$ , the vertices to be eliminated are selected randomly from C. All of the eliminated vertices are collected in X. The set of neighboring solutions induced by the Press(v, X) operator is given by:

$$\mathcal{N}_{Press} = \{ C' : C' \leftarrow C \oplus Press(v, X), v \in V \setminus C, X \subseteq C \setminus N(v) \}$$
 (6)

By definition, Press eliminates at least two vertices from the solution (i.e., |X| > 1). Thus, the application of Press always degrades the quality of the current solution. Hence, this operator is only used by the perturbation procedure when the Add and Swap operators are no longer applicable, or when the search stagnates in local optima.

Next, we provide some additional comments about these operators and discuss some implementation issues.

- Add and Swap have been used in several algorithms for the maximum clique problem and the equivalent maximum independent set problem [4,13,25,35]. However, given the generality of the maximum s-plex problem, the definitions of these operators are different and more complex in the present study; in particular, the saturated set S must be involved. The Press operator can be treated as a dedicated application to the s-plex problem of the PUSH operator introduced in [38] for the maximum weight clique problem.
- It should be noted that traditional maximum clique benchmark graphs, such as those from the 2nd DIMACS Challenge, include many dense graphs. Thus, most maximum clique algorithms typically operate on the complement graph  $\bar{G}$  to search for the maximum independent sets in terms of runtime efficiency [13, 25, 33]. The FD-TS algorithm proposed in the present study is designed to handle very large real-world networks, which are typically sparse graphs. Consequently, FD-TS operates directly on the original graph based on its adjacency-list representation. In addition, we restrict the candidate vertices considered by the three move operators to N(C) instead of  $V \setminus C$  because N(C) is much smaller than  $V \setminus C$  for very large networks. Indeed, given the size of the networks considered (up to millions of ver-

tices), even a simple operation such as scanning all the vertices of  $V \setminus C$  becomes too expensive, and it can slow down the algorithm considerably. Thus, special care is taken to avoid ineffective or unpromising examinations.

• When Swap and Press are applied, each dropped vertex is forbidden from rejoining the solution during a number of subsequent iterations to avoid revisiting previously examined solutions. This is achieved by using a tabu list (see Section 2.5.4).

# 2.4 Constructing the initial solutions

Each round of the FD-TS algorithm requires a starting solution (see Algorithm 1, line 5). In general, the starting solutions can be generated by any method that ensures the s-plex property. In our method, we employ the following construction procedure, which applies the Add operator while considering the frequency information for vertex moves.

From a random sample of q vertices  $(q \in [50, 150])$ , we use the vertex with the minimum frequency (ties are broken randomly) to create a singleton set C. From the singleton s-plex C, the procedure repeats the following three steps to extend the current solution: 1) generate the  $M_1$  set from N(C), 2) select one vertex with the minimum frequency from  $M_1$  (ties are broken randomly), and 3) add the selected vertex to C. We repeat this process until  $M_1$  becomes empty. The final s-plex C is returned as the initial solution.

The intuitive assumption that less frequently moved vertices are preferred is intended to make the initial solution as diverse as possible. Moreover, using a sample of q vertices instead of the whole set V for seeding the solution helps to reduce the computational overheads in the initialization procedure. This is particularly true for massive graphs because scanning all the vertices of the graph can be time consuming in this case. In Section 4.2, we discuss the calibration of the parameter q.

# 2.5 FD-TS

# 2.5.1 General procedure

The key search procedure employed in the FD-TS algorithm (see Algorithm 2) combines a double-neighborhood search procedure (we refer to this procedure as TS<sup>2</sup>; Algorithm 2, lines 11–24) to facilitate intensification (to obtain local optima) and a frequency-based perturbation procedure (we refer to this procedure as PERTURB; Algorithm 2, lines 25–30) for diversification (to escape from local optima).

# Algorithm 2: Frequency driven tabu search

```
Input: Problem instance (G, s), current solution C, max. allowed iterations L
     Output: The largest s-plex found C_{best}
 1 begin
  2
           C_{best} \leftarrow C;
  3
           tabu\_list \leftarrow \emptyset;
  4
           l \leftarrow 0;
                                                                          /* Counter of cycles of TS2 + PERTURB runs */
           fix \leftarrow \emptyset;
                                                                          /* The vertices that are forbidden to drop */
 5
           freq(v) \leftarrow 0 \text{ for all } v \in V ;
  6
                                                                                                 /* Reset frequency records */
                                                               /* Counter of consecutive non-improving iterations */
           // Run {
m TS}^2 + PERTURB a maximum of L iterations
           while l \le L do
  8
                 // Start the two neighborhood tabu search procedure - {
m TS}^2
  9
                 Updating the saturated subset S:
                                                                                                                     /* Sect. 2.2 */
                 Decompose set N(C) into M'_1 = \{v \in M_1 : v \notin tabu\_list \lor |C| + 1 > |C_{best}|\},
10
                 M_2' = \{v \in M_2 : v \notin tabu\_list\}, M_3' = \{v \in M_3 : v \notin tabu\_list\};
                                                                                                                     /* Sect. 2.3 */
11
                 if M_1' \neq \emptyset then
                       v \leftarrow a random vertex from M_1';
12
13
                       C \leftarrow C \oplus Add(v);
14
                       freq(v) \leftarrow freq(v) + 1;
                 else if M_2' \neq \emptyset then
15
                       (v, \overline{u}) \leftarrow \text{two random exchangeable vertices from } M'_2 \times N(C); /* See Sect. 2.5.2 for
16
                       selection rule */
                       C \leftarrow C \oplus Swap(v, u);
17
18
                       Add u to tabu\_list with tabu tenure T_u;
                                                                                                                  /* Sect. 2.5.4 */
19
                       freq(v) \leftarrow freq(v) + 1, freq(u) \leftarrow freq(u) + 1;
20
                 if |C| > |C_{best}| then
21
                       C_{best} \leftarrow C;
22
                       \lambda \leftarrow 0;
23
                 else
                  \  \  \, \bigsqcup \  \  \, \lambda \leftarrow \lambda + 1;
                 // Run the PERTURB procedure
                 if (No feasible Add and Swap operation) \vee (\lambda > s * |C_{best}|) \wedge M_3' \neq \emptyset) then v \leftarrow a vertex with maximum freq(v) from M_3', break ties randomly;
25
26
27
                       C \leftarrow C \oplus Press(v,X); /* X collects the removed vertices from C, Sect. 2.3 */
                       fix \leftarrow \{v\}, \lambda \leftarrow 0;
28
29
                       freq(v) \leftarrow 0, freq(u) \leftarrow freq(u) + 1 \text{ for all } u \text{ in } X;
                       Add each u \in X to tabu_list with tabu tenure T_u;
30
                                                                                                                  /* Sect. 2.5.4 */
31
32 end
33 Return C_{best}, freq;
```

Based on a given initial solution,  $TS^2$  uses Add and Swap to improve the current solution until search stagnation occurs. PERTURB then applies Press to modify (perturb) the current local optimum and passes the modified solution to  $TS^2$  for further improvement. FD-TS iterates this  $TS^2$ +PERTURB process a maximum of L times and then starts the next round of its search procedure.

In addition to the three move operators (Add, Swap, and Press), FD-TS employs a tabu mechanism (see Section 2.5.4) [11] and a frequency technique to ensure the effective exploration of the search space. A tabu list ( $tabu\_list$ ) is used to mark the vertices that are forbidden from joining the current solution during a specific number of iterations. Information related to the move frequency of each vertex v is collected, where freq(v) records the number of times that v is operated upon by a move operator in the recent history. Initially,  $tabu\_list$  is empty and the frequency of each vertex is set to 0. The

fix set (a singleton set used by the PERTURB procedure) records an added vertex, which is forbidden from moving out of the current solution in TS<sup>2</sup> and the counter l counts the completed iterations, the upper limit of which is given by the parameter L. Moreover, in order to avoid being trapped by local optima, a counter  $\lambda$  records the consecutive iterations that have passed since the last improvement of the current solution. Each time that the counter  $\lambda$  reaches a threshold, the perturbation procedure is triggered to modify the current solution using the Press operator.

At the beginning of each iteration, set C is checked to identify the saturated subset S. Next, N(C) is decomposed into three disjoint subsets:  $M'_1$ ,  $M'_2$ , and  $M'_3$ . These sets correspond to  $M_1$ ,  $M_2$ , and  $M_3$  (defined in Section 2.3), respectively, but they exclude the vertices in the  $tabu\_list$ . However, a vertex  $v \in M_1$  is always retained in  $M'_1$  if adding v to C leads to a new solution that is better than the best solution found previously, i.e.,  $|C| + 1 > |C_{best}|$ , regardless of the tabu status of the vertex.

#### 2.5.2 Solution improvement with Add and Swap

The current solution is transformed successively by applying Add and Swap. Preference is given to the Add operator. Thus, whenever  $M'_1$  is not empty, Add is applied to improve the current solution by adding one vertex of  $M'_1$  to the solution (Algorithm 2, lines 11–14). If no vertex can be added to the solution  $(M'_1 = \emptyset)$ , but  $M'_2$  is not empty, then the search continues with the Swap(v, u) operator by visiting solutions of equal quality (Algorithm 2, lines 15–19).

To provide Swap(v, u) with an appropriate exchangeable pair  $\langle v, u \rangle$ , v is first selected randomly from  $M'_2$ , and u is then selected from the candidate set defined by the rules given in Section 2.3, while excluding the vertex recorded in fix. In particular, if v is a vertex of type (set) A, then u is selected from  $S \setminus (N(v) \cup fix)$  (which is a trivial set with zero or one vertex); and if v is a vertex of type (set) B, u is selected randomly from  $C \setminus (N(v) \cup fix)$ . If there is no eligible candidate for u ( $S \setminus N(v) = fix$  or  $C \setminus N(v) = fix$ ), then we simply give up attempting to apply Swap(v, u) and move on to the PERTURB procedure (Algorithm 2, line 25).

#### 2.5.3 Perturbation with Press

Inevitably, at a certain search stage, no candidate vertex v or candidate pair  $\langle v, u \rangle$  is available for the Add(v) or Swap(v, u) operator (i.e., both  $M'_1$  and  $M'_2$  are empty or no eligible vertex can be found for u), or the search stagnates on the Swap(v, u) operator. In the latter case, search stagnation occurs when the current solution has not been improved for  $s*|C_{best}|$  consecutive iterations. The self-adaptive threshold,  $s*|C_{best}|$ , is identified based on the assumption

that if the current s-plex cannot be improved even after the replacement of each of its vertices at least s times by the Swap operator, then no better solution can be found in the current search region. To continue the search, the algorithm triggers the PERTURB procedure in order to move the search to a distant new region. This procedure applies the Press(v, X) operator, where v is the vertex from  $M'_3$  with the largest freq value and X collects the dropped vertices from  $C \setminus N(v)$  to recover a feasible solution (Algorithm 2, lines 26–27). It is important to reset the frequency of vertex v (Algorithm 2, line 29), or the accumulated frequency of this vertex could dominate other vertices in subsequent cycles. The dropped vertices in X are added to the tabu list (Algorithm 2, line 30) and they will not be considered during the forbidden period, as explained in the next section.

# 2.5.4 Tabu tenure and management

As mentioned above, to avoid revisiting recently examined solutions, we use a tabu list to record the vertices dropped from the current solution in order to exclude them from consideration during a number of consecutive iterations. According to the definitions in Section 2.3, each application of Swap(v, u) or Press(v, X) removes one or more vertices from the current s-plex. Each dropped vertex u will be kept in the tabu list for the next  $T_u$  iterations (the tabu tenure), which is set according to the following two rules:

$$\begin{cases}
T_u = 10 + random(0, |M_2|), & u \text{ is dropped by } Swap(v, u) \\
T_u = 7, & u \in X \text{ is dropped by } Press(v, X)
\end{cases}$$
(7)

where random(0, I) is a random integer in  $\{0, \ldots, I\}$ . These rules are based on previous studies of the related maximum clique problem [13,35]. The first rule estimates the forbidden period for vertices for side-walk moves with the Swap operator, which ensures that a dropped vertex will not be reconsidered for at least 10 iterations, and the second rule for the Press operator prevents any dropped vertex from being reconsidered for a small number of iterations (seven in this case). Based on experiments, we observed that other values around these tabu tenures obtained similar performance. Thus, we selected the values used in [13,35].

Finally, the tabu list is more useful when the number of candidate vertices for Swap or Press is limited, because a dropped vertex u will have a high probability of being added again to the solution if it is not prohibited. However, if numerous candidate vertices exist (such as in massive graphs), there is little chance of a dropped vertex being re-selected immediately. Thus, the tabu mechanism is more useful for graphs of limited size than massive graphs.

Given a graph G = (V, E) and a parameter s, suppose that after tabu search (Algorithm 1, line 6), the current best s-plex has cardinality  $|C^*|$  (lower bound of the maximum s-plex of G). Clearly, to further improve  $C^*$ , considering any vertex in V with a degree smaller than or equal to  $|C^*| - s$  would not be beneficial because such a vertex cannot extend  $C^*$ . Thus, these vertices can be safely removed from the graph [1]. In FD-TS, we explore this strategy using the  $Peel(G, s, |C^*|)$  procedure (Algorithm 1, line 9), which recursively deletes the vertices (and their incident edges) with a degree less than or equal to  $|C^*| - s$  until no such vertex exists. Finally, if the subgraphs obtained after  $Peel(G, s, |C^*|)$  have fewer vertices than  $|C^*|$ , then  $C^*$  must be an optimal solution because no better solution can exist.

For very dense graphs, the Peel procedure may not reduce the graph size greatly because the degrees of most vertices will remain larger than  $|C^*| - s$ . However, this technique is highly effective when it is applied to large sparse graphs such as massive real-world complex networks. As shown in Section 4.3, by using the high-quality lower bound  $|C^*|$  provided by our tabu search procedure, this pruning technique can effectively reduce large sparse graphs to very small graphs (even the null graph).

We note that the idea of removing unpromising vertices was used previously in a GRASP heuristic for detecting dense subgraphs (quasi-cliques) in massive sparse graphs [1], as well as in several exact algorithms for the maximum clique and s-plex problems [3, 29, 32].

#### 3 Implementation and time complexity

To effectively implement FD-TS, we maintain two structures: the vector  $deg_C[v]$  (i.e.,  $deg_C[v] = |N(v) \cup C|, v \in V$ ) and the set N(C) (i.e.,  $N(C) = \bigcup_{v \in C} N(v) \setminus C$ ), which are updated whenever the current solution C changes. Thus, each time a vertex (say u) is added to or removed from C by a move operator, we increase or decrease  $deg_C[v]$  by one for each  $v \in N(u)$ . Since the set N(C) must only contain vertices with  $deg_C[v] > 0$ , it is also adjusted when  $deg_C[v]$  changes.

Next, we discuss the time complexity of the main components of the proposed algorithm. First, we consider the procedure for constructing initial solutions (Section 2.4). In each iteration, we need to build the subset  $M_1$  from N(C) and update  $deg_C[v]$  and N(C) after adding a vertex, which can be achieved in  $O(|N(C)| + \Delta)$  ( $\Delta = \max_{v \in V} \{|N(v)|\}$ ).

The efficiency of  $TS^2$  (Section 2.5) and PERTURB (Section 2.5.3) is closely related to the method used for building sets  $M'_1$ ,  $M'_2$ , and  $M'_3$  from N(C) from scratch in each iteration. In our implementation, we build the saturated set S (Algorithm 2, line 9) from C in a time of O(|C|) at the very beginning of each iteration. Then, for each vertex in N(C) (e.g., u), we count the number of saturated vertices in the set of vertices adjacent to u, i.e.,  $|S \cap N(u)|$  (the saturated connectivity of u). Obviously, if the saturated connectivity of u is  $0, S \setminus N(u) = \emptyset$ ; and if the saturated connectivity of u is  $1, |S \setminus N(u)| = 1$ . According to the definitions of  $M_1$ ,  $M_2$ , and  $M_3$ , once the saturated connectivity of u and  $deg_C[u]$  is known, it is trivial to identify u as an element of  $M'_1$ ,  $M'_2$  or  $M'_3$ . Consequently, decomposing N(C) (Algorithm 2, line 10) can be achieved in  $O(|N(C)| * \Delta)$ .

Next, we consider the time complexity when employing the move operators. First, for the Add operator, we only need to update the vector  $deg_C[v]$  and set N(C) after reallocating a vertex v, which can be achieved in  $O(\Delta)$ . Second, to apply Swap with a vertex  $v \in N(C)$ , we first need to identify the other vertex  $u \in C$ . According to the rule defined in Section 2.5.2, the sets  $S \setminus N(v)$  and  $C \setminus N(v)$  can be identified by traversing sets C and N(v) respectively. Thus, the time required to identify u is bounded by  $O(|C|+\Delta) = O(2*\Delta+s) = O(\Delta+s)$  (because  $|C| \leq \Delta + s$ ), while updating  $deg_C[v]$  and N(C) is bounded by  $O(2*\Delta) = O(\Delta)$  because two vertices are displaced during each application of Swap. Finally, each application of the Press operator can be achieved in  $O(\Delta^2)$ .

Overall, one operator (Add, Swap or Press) is applied during one iteration, so the total time complexity of  $TS^2$  and PERTURB for each iteration is bounded by  $O(|C|+|N(C)|*\Delta+\Delta^2)$ . We note that for sparse graphs, |C|, N(C), and  $\Delta$  are usually extremely small compared with the number of vertices in a graph.

## 4 Computational assessment

#### 4.1 Benchmarks

In this section, we present computational evaluations of the proposed FD-TS algorithm for the maximum s-plex problem based on the following three sets of 79 (19, 17, and 43, respectively) benchmark graphs.

• Stanford Large Network Dataset Collection (SNAP) <sup>1</sup>. SNAP provides a large range of large-scale social and information networks [15], in-

<sup>&</sup>lt;sup>1</sup> http://snap.stanford.edu/data/

cluding graphs retrieved from social networks, communication networks, citation networks, web graphs, product co-purchasing networks, Internet peer-to-peer networks, and Wikipedia networks. Some of these networks are directed graphs, so we simply ignored the direction of each edge and eliminate self-looping and duplicated edges.

- The 10th DIMACS Implementation Challenge Benchmark (10th DIMACS)<sup>2</sup>. This testbed contains many large networks, including artificial and real-world graphs from different applications. This benchmark set is popular for testing graph clustering and partitioning algorithms. More information about the graphs can be obtained from [2].
- The 2nd DIMACS Implementation Challenge Benchmark (2nd DIMACS)<sup>3</sup>. This set is from the 2nd DIMACS Implementation Challenge for the maximum clique problem. These instances cover real-world problems (e.g., coding theory, fault diagnosis, and the Steiner triple problem) and random graphs. The instances range from small graphs (50 vertices and 1,000 edges) to large graphs (4,000 vertices and 5,000,000 edges). These DIMACS graphs are very popular and they are generally used as a testbed for evaluating clique and s-plex algorithms. Unlike the instances in the two first benchmark sets, most of these instances are dense graphs.

# 4.2 Experimental protocol and parameter tuning

The proposed FD-TS algorithm was implemented in C++ <sup>4</sup> and compiled by g++ with optimization option '-O3'. All experiments were conducted on a computer with an AMD Opteron 4184 processor (2.8 GHz and 2 GB RAM) running CentOS 6.5. When we solved the DIMACS machine benchmarking program fdmax.c <sup>5</sup> without compilation optimization flag, the run time on our machine was 0.40, 2.50, and 9.55 seconds for graphs r300.5, r400.5, and r500.5, respectively.

An interesting feature of FD-TS is that it has very few parameters. In addition to the tabu tenure discussed in Section 2.5.4, the parameter q (the sample number of vertices in Section 2.4) was set to 100. As indicated in Section 2.4, q is not a sensitive parameter so we simply fixed it to the middle value in the range of [50,150]. However, according to our experiments, the best value of parameter L, which specifies the maximum number of iterations in each round of tabu search, was highly dependent on the instance considered. We compared different values in  $\{10, 100, 1000, 5000\}$  for L with s = 2, 3, 4, 5 for six selected instances from the 2nd DIMACS benchmark set  $(MANN_a27, brock400_a2, brock400_$ 

 $<sup>^2\ \</sup>mathrm{http://www.cc.gatech.edu/dimacs10/downloads.shtml}$ 

<sup>3</sup> http://www.cs.hbg.psu.edu/txn131/clique.html

 $<sup>^4</sup>$  We will make our program available.

<sup>&</sup>lt;sup>5</sup> ftp://dimacs.rutgers.edu/pub/dsj/clique/

brock800\_2, C1000.9, keller6, p\_hat1500-3). As a trade-off, we retained L = 1000 because the algorithm achieved the best average solution quality in most cases. We also observed that fine tuning L for each pair (instances, s) could improve the performance. However, to report our computational results, we used the fixed setting L = 1000, which allowed the algorithm to obtain highly competitive results. Given the stochastic nature of our algorithm, we ran FD-TS to solve each instance 20 times for each given s, where each run was limited to a maximum of 180 CPU seconds (3 minutes).

# 4.3 Computational results for very large networks from SNAP and the 10th DIMACS Challenge

Table 1 shows the performance of FD-TS on 47 instances taken from the SNAP and 10th DIMACS benchmarks, namely, all 37 instances tested in [29] for the s-plex problem, as well as 10 instances from the recent literature [32]. Note that in [32], only the best clique size was reported, which is just a lower bound of the maximum s-plex for s = 2, 3, 4, 5. To be complete, we also report our results for the remaining 20 instances from [32] in the Appendix (Table A.1).

For each instance, the columns "Instance," "|V|," and "|E|" indicate basic information for the name, number of vertices, and number of edges, respectively. For each s=2,3,4,5, column "BKV" denotes the best-known objective values collected from [29] (for the first 37 instances) and [32] (for the last 10 instances). For the items in this column, an additional symbol "\*" indicates that this objective value was proved to be optimal in [29], and the symbol  $\omega$ shows that this value is the maximum clique size mentioned in [32] (which is a lower bound of the maximum s-plex). The "max" column shows the best objective value found by FD-TS among its 20 trials and the "time" column indicates the average time (in seconds) required for runs to obtain the best objective value (excluding the time spent reading the graph). The "|V'|" column shows the number of remaining vertices in the reduced subgraph after executing the Peel(G, s, max) procedure (see Section 2.6). As mentioned earlier, if the number of vertices in the reduced subgraph (column "|V'|") is less than or equal to the lower bound (column "max"), then the latter is guaranteed to be the optimal solution. In these cases, we put a "\*" beside the value.

Moreover, for instances where the optimum could not be determined in either [29] or FD-TS, we conducted an additional experiment with the CPLEX solver (version 12.6.1). First, we reduced the original graph by applying the Peel(G, s, max) procedure. Then, for the reduced subgraph and a given  $s \in \{2, 3, 4, 5\}$ , a cutoff time of one CPU hour was used when running CPLEX with the mathematical model (1) presented in Section 1. Experiments were conducted using the same machine employed for running FD-TS. The best

objective values found by CPLEX are listed in the "cplex" column, where "\*" indicates the optimal values. If CPLEX was unable to load the model, the entry is marked by "N/A."

Table 1 shows that FD-TS always obtained the same or better objective values compared with the current best-known results (BKV values) (better objective values are highlighted with a bold font). In particular, FD-TS improved the best-known results for more instances as s increased (FD-TS found better solutions for 7, 15, 19, 21 instances with s = 2, 3, 4, 5, respectively). This observation indicates that the instances become more challenging for exact algorithms with a larger s. Using the Peel procedure, FD-TS was also provably optimum for the first time for 6, 6, 6, 10 instances and s = 2, 3, 4, 5, respectively. For the cases where the number of vertices in the reduced subgraph remained larger than the lower bound given by FD-TS, the majority of these cases were manageable when the subgraph had less than 10,000 vertices (exceptions include the instances 333SP, cage15, cit-Pattens, and wiki-Talk for s=2). In terms of the computational time, FD-TS was able to obtain the best solutions in less than one second in most cases, including instances with millions of vertices. The average time required by FD-TS on cit-Patents for s=3 was the longest but still less than one minute. Unfortunately, and similarly to CPLEX, it could not determine any solution better than that found by FD-TS in one hour when given the reduced subgraph. However, for the instances rgg\_n\_2\_17\_s0 with s = 5, and rgg\_n\_2\_20\_s0 with s = 4 and 5, optimal solutions were also obtained by CPLEX. Finally, we note that for the instances coPapersCiteseer, coPapersDBLP, and cond-mat-2005, the size of the maximum clique was the same as the size of the maximum s-plex for s = 2, 3, 4, 5.

Table 1 Computational results of FD-TS on 47 large networks from the SNAP Collection and the 10th DIMACS Implementation Challenge.

Computational results of FD-13 on 41 large networks from the SNAP	resuits		7 110 6	71 19T	36 1100	WOLDS	HOTH	TIC DII	- 1	OILCO.	CONCOUNT WITH THE		TOOT D	77777	1-11-1 ()	DIMITO IIII DICIII CII CII AII CII SCII SCII SCII SCI	LUCAULCE	1 C11&1	201121			
Instance	Ā	E	BKV	max	s=2	1,/1	colex		max	s=3	//	colex	BKV	max	s=4	1//1	colex	BKV	max	s=5	1,1	colex
adinonn	112	425	***	9	0.00	102			×	- 1	102		**		- 1	- 68		10*			102	
celegans_metabolic	453	2025	10*	10	0.00	313	,	11*	11	0.00	429	,	13*	13	0.00	240	,	14*	14	0.00	429	,
dolphins	62	159	*9	9	0.00	53		*-	7		36	_	<b>*</b>			45	,	*6			45	
email	1133	5451	12*	12	0.01	238		12*	12		349	_	12*			349	,	13*			848	
football	115	613	10*	10	0.01	114		11*	11		115	,	12*			115	,	12*			115	
jazz	198	2742	30*	30	0.00	130		30*	30		164	,	30*			127	,	30*			127	-
karate	34	78	*9	9	0.00	22		*9	9		10	,	**			10	,	$11^a$			0	
netscience	1589	2742	*02	20	0.00	20		*02	20		137	,	*02			158	,	20			20	
polblogs	1490	16715	23*	23	0.02	541		27*	27		894	,	*62			489	,	32			293	32*
polbooks	105	441	*_	7	0.00	103		*6	6		103	_	10*			105		$11^{*}$			105	
power	4941	6594	*9	9	0.00	36		*9	9		231	_	9			12	*	8			12	*6
PGPgiantcompo	10680	24316	*62	29	0.03	115		31*	31		43		33*			41		35*			41	
as-22july06	22963	48436	19*	19	0.03	117		21*	21		110	,	22*			110		24*			104	
astro-ph	16706	121251	27*	22	90.0	113	,	22*	57		113	,	27*			113	,	27*			165	,
caidaRouterLevel	192244	990609	20*	20	0.31	2039		22	23		1290	13	23			1290	12	23			860	14
cnr-2000	325557	2738969	*28	80 22 80	6.75	0		*98	98	_	0	,	*98			98		*98			98	
coAuthorsCiteseer	227320	814134	*42	*28	0.21	87		*28	*48		87	,	*2*			87	,	*2*			87	,
coAuthorsDBLP	299067	92926	115*	115*	0.27	115		115*	115*		115	,	115*			115	,	115*			115	,
cond-mat-2002	40421	175691	30*	30*	80.0	30	-	30*	*08		30	,	30*			30		30*			57	
memplus	17758	54196	*26	*46	80.0	26	,	*26	*26		26	,	*26			26		*26			26	1
rgg_n_2_17_s0	131072	728753	16*	16*	0.15	0		17*	17*		0		18*			0		18			34	18*
rgg_n_2_19_s0	524288	3269766	19*	19*	69.0	0		19*	19*		19	,	*02			19		20			19	
rgg_n_2_20_s0	1048576	6891620	18*	18	1.42	29		19*	19		59	,	19			59		19			172	20*
cit-HepPh	34546	420877	24*	24	0.37	4768		25	27		3498		25			2344	11	25			.804	22
cit-HepTh	27770	352285	<b>58</b> *	28	1.20	4815		31	31		3999		31			2922	9	31			.726	23
email-EuAll	265214	364481	19*	19	0.15	1357		22	22		1198	20	22			1055	23	27			886	26
p2p-Gnutella04	10876	39994	* *	ю	0.22	6809		rc.	4		5433		9			4857	ю	7			1857	N/A
p2p-Gnutella24	26518	62369	۳0 *	ю	0.07	9271		ы	9		9271	N/A	22			7480	N/A	22			480	N/A
p2p-Gnutella25	22687	54705	ж *	ю	0.03	7892		വ	9		7892		ಬ			6091	ro	വ			0	
soc-Epinions1	75879	405740	58* 58*	28	0.09	4156	12	28	32		3791	12	28			3280	12	28			8178	10
soc-Slashdot0811	77360	469180	31*	31	0.02	3665		33	34		3188	9	36			2596	-	36			3416	7
soc-Slashdot0902	82168	504230	32*	32	0.02	3669		32	35		3185	9	36			2435	6	36			287	7
web-BerkStan	685230	6649470	202*	202	4.19	392		202*	202		392	,	205 <sub>*</sub>			392		205 <sub>*</sub>			392	,
web-Google	875713	4322051	46*	46*	1.13	0		*47*	*7*		0		48*			0		*8*			48	
web-NotreDame	325729	1090108	155*	155	0.74	1367		155*	155		1367		155*			1367		155*			367	
web-Stanford	281903	1992636	* * *	64	3.86	741		. 64*	64		985	, ,	64			985	65	64			985	65
w1k1-Vote	7115	100762	21.	7.7	0.10	2098		7.74	24		2002	+	526		- 1	1925	16	7.7			925	17
333SP	3712815	11108633	3	ທີ	1.19	2261408	N/A	4(ω)	9		261408		$4(\omega)$			2261408	N/A	$\frac{4}{3}$			61408	N/A
belgium.osm	1441295	1549970	3(6)	* •	99.0	0		$\frac{3(\kappa)}{2}$	* •		22	, ;	$3(\omega)$			22	, ;	$\frac{3(\kappa)}{2}$			22	
cage15	5154859	47022346	(ω) (ω)		2.48	5135355	A/A	(α) (π)	20		134115		$6(\omega)$			5091619	A/A	6(ω)			91619	N/A
coPapersCiteseer	434102	16036720	845(ω)		7.92	845		$845(\omega)$	845*		845		$845(\omega)$			845	,	$845(\omega)$			845	,
coPapersDBLP	540486	15245729	$337(\omega)$	•	2.83	337		337(w)	337*		337		$337(\omega)$			337		$337(\omega)$			337	
amazon0312	400727	2349869	- 11(ε)		0.54	0		$11(\omega)$	13*		0		$11(\omega)$			0	,	$11(\omega)$			0	
amazon0505	410236	2439437	$11(\omega)$		0.50	0		$11(\omega)$	13*		0		$11(\omega)$			0	,	$11(\omega)$			0	,
amazon0601	403394	2443408	$11(\omega)$	12*	0.75	0		$11(\omega)$	13*		0		$11(\omega)$			0	,	$11(\omega)$			0	,
cit-Patents	3774768	16518947	$11(\omega)$		21.25	29585	N/A	$11(\omega)$	21	51.02	14717	N/A	$11(\omega)$			6293	N/A	$11(\omega)$			8604	2
wiki-Talk	2394385	4659565	$26(\omega)$	32	1.18	11327	N/A	$26(\omega)$	36	3.60	9814	_	$26(\omega)$			8376	N/A	$26(\omega)$			7119	N/A
Note $a$ : The value of 11 reported in [29] for 'karate' is wrongly claimed to be the	of 11 report	ed in [29] for	· 'karate	i is wron	gly clain	ned to be		optimal solution.	on.			_										

Table 2 shows the computational results obtained by FD-TS for 52 classical hard instances from the 2nd DIMACS set, with s=2,3,4,5. The first group contains all 36 instances that have been studied by four state-of-the-art s-plex algorithms [3, 16, 19, 29]. The second group includes 16 additional large 2nd DIMACS instances with at least 800 vertices, which have not been tested previously by any existing s-plex algorithm. For the sake of completeness, we also tested the remaining 28 instances of this benchmark set, for which no previous s-plex result is available. These results are reported in Table A.2 of the Appendix.

For each instance, the following information is included. The "|V|" column shows the number of vertices in the original graph. The " $\omega$ " column indicates the best-known maximum clique size reported previously [34] (lower bounds for the maximum s-plex). The "BKV" column indicates the best-known objective values obtained by the algorithms in [3, 16, 19, 29] (proven optima are indicated by "\*"). The letters between parentheses following each "BKV" value indicate the algorithm(s) that obtained the BKV value.

- "B" A branch-and-cut algorithm [3] based on polyhedral analysis of the convex hull of the maximum s-plex problem. This algorithm was evaluated based on instances from the 2nd DIMACS set with s=1,2. Each instance was solved within a maximum of 3 hours on a machine with a 2.66 GHz XEON® processor, 3 GB RAM, and 120 GB HDD.
- "M" A branch-and-bound algorithm [16] adapted from the classical maximum clique algorithm [20]. Results were obtained based on 2nd DIMACS instances with s=2,3,4. The experiments were conducted on a machine with a 2.2 GHz Dual-Core AMD Opteron processor and 3 GB RAM. A time limit of one hour was allowed to solve each instance.
- "T" A generalized algorithm framework used to detect optimal hereditary structures in graphs [29]. For the maximum s-plex problem, this approach was tested based on instances from the 2nd DIMACS, 10th DIMACS and SNAP benchmark sets, for s=2,3,4,5. Experiments were conducted with a Dell Optiplex GX620 computer with an Intel Core<sup>TM</sup>2 Quad 3 GHz processor and 4 GB RAM with a time limit of 3 hours for each instance.
- "H" Exact combinatorial algorithms based on methods from parameterized algorithmics [19]. Results were reported for a subset of the 2nd DIMACS instances (s = 1, 2) with a time limit of 3 hours on a machine with an AMD Athlon 64 3700+ 2.2 GHz CPU, 3 GB RAM, and 1M L2 cache.

For instances where the optimal solution has not been proven by any of the algorithms mentioned above (i.e., no "\*" is indicated for "BKV"), we used the CPLEX solver to solve these instances (i.e., their reduced subgraphs after

applying the Peel procedure, see Section 2.6) with a time limit of one CPU hour on our computer. The "cplex" column shows the best feasible solutions attained by CPLEX. The "max(ave)" column reports the maximum value achieved by FD-TS in 20 runs and the average value (in parentheses) if the 20 best values were not the same. The "time" column shows the average time (in seconds) required by the runs that obtained the best value among the 20 runs. Obviously, the total time allowed to FD-TS in 20 runs (180\*20 = 3600 s) was exactly one hour.

Table 2 shows that the FD-TS algorithm matched or improved (highlighted in bold font) the current best-known results with s = 2, 3, 4, 5. The average objective values obtained by our algorithm were even better than the bestknown values based on these instances for different s (except for MANN\_a27 and MANN\_a45 with s=2). In terms of the stability of the best solution, for most of the small instances ( $|V| \leq 400$ ) in the first group, the best solution could be obtained in each of the 20 runs (except for MANN\_a27, brock400\_4 and  $\sin 200 \cdot 0.7 \cdot 2$  with s = 2, brock  $400 \cdot 1$  with s = 5). The larger graphs in the second group of the table (which were not reported previously), such as brock800\_X, CXXX.X, hamming10-4, and keller6, represent the most challenging cases for FD-TS because the best solution could not be found in every run. Moreover, for instances such as brock400\_1, brock800\_2, brock800\_3, keller6, and p\_hat1500-2, FD-TS failed to achieve a 100% success rate as sincreased. However, for instances such as MANN\_a27 and brock800\_4, there was no correlation between the success rate and the value of s. In terms of the computational time, FD-TS achieved its best values rather quickly, since it rarely exceeded one minute, whereas CPLEX failed to solve these instances (in fact, the reduced subgraphs) to optimality for any s within one hour. Nevertheless, for the cases where the optimal value is still unknown, the lower bounds obtained by CPLEX were competitive compared with the four other reference algorithms [3, 16, 19, 29].

Table 2 Computational results of FD-TS on 52 benchmark instances of the 2nd DIMACS Implementation Challenge

Computational results of FD-13 on 32 benchmark instances of the 2nd DIMACS implementation Chanenge	Hai Ici	Surve	OT _ T T TO	CFT		~*** XT TY	COLUMN OF	7110	ZIM DIMI		OICTIOTTO	1000	STIGHTON I					
90	1771		DIVI		Ω.	17	1271				77.77		s=4	7.7	757.0			7
Instance	_	3	BK V	cpiex		time	BKV	cplex		time	BK V	cplex	max(ave)	time	BKV	cplex		time
MANN_a9 MANN a27	45 378	126	26*(BMTH) 236*(H)		26 236(235 90)	0.00	36*(MT)		351	0.00	36*(MT)	1 85 125 1	36 351	0.00	44(T) <sup>2</sup> 351(T)	351	45 351	0.00
MANN-a45	1035	345	662(BH)	662*		5.46	990(M)	*066	066	7.40	990(M)	*066	066	7.48	(T) TOG	*066	066	7.44
$brock200_{-1}$	200	21	25(B)	56		0.15	24(M)	29	30	0.05	27(M)	33	35	3.59	27(T)	38	39	2.36
$brock200_{-2}$	200	12	13*(MTH)	,	13	0.01	16*(T)	,	16	0.27	17(TM)	17	18	90.0	17(TM)	20	20	0.04
$brock200_{-3}$	200	15	17*(T)	1	17	0.02	19(T)	19	20	0.02	19(T)	22	23	0.03	19(T)	25	26	0.09
$brock200_{-4}$	200	17	20*(TH)	ı	20	90.0	20(T)	22	23	0.07	21(M)	26	26	0.03	21(M)	28	30	0.20
brock400_1	400	27	23(T)	29	30	0.42	23(T)	34	36	7.24	23(T)	37	41	41.37	23(T)	43	<b>46</b> (45.50)	36.57
brock400_2	400	29	27(B)	8 0	30	0.51	27(M)		36	15.37	29(M)	37	41	33.51	29(M)	41	45	2.14
brock400_4	400		27(B)	28	<b>33</b> (31.20)	64.18	27(B)	34	36	4.60	30(M)	36	41	1.76	30(M)	44	46	25.67
c-fat200-1	200	12	12*(BMTH)	1	12	0.00	12*(MT)		12	0.00	$12^*(MT)$		12	0.00	$14^{*}(T)$		14	0.00
c-fat200-2	200	24	24*(BMTH)	,	24	0.01	24*(MT)		24	0.01	24*(MT)		24	0.02	$24^{*}(T)$	,	24	0.01
c-fat200-5	200	то 80	58*(BMTH)	í	28	0.01	58*(MT)	í	28	0.01	58*(MT)	,	28	0.00	58*(T)	í	28	0.00
c-fat500-1	200	14	14*(BMTH)		14	0.00	14*(MT)		14	0.00	14*(MT)		14	0.00	15*(T)		15	0.00
c-fat500-2	200	56	26*(BMTH)	,	26	0.00	26*(MT)	,	26	0.00	26*(MT)	,	26	0.00	26*(T)	,	26	0.00
c-fat500-5	200	64	64*(BMTH)	1	64	0.02	64*(MT)		64	0.02	64*(MT)		64	0.03	64*(T)	1	64	0.04
c-fat500-10	200	126	126*(BMTH)	-	126	0.08	126*(MT)	,	126	0.22	126*(MT)		126	0.23	126*(T)	,	126	0.16
hamming6-2	64	32	32*(BMTH)	í	32	0.00	32*(MT)		32	0.00	40*(MT)		40	0.00	48*(T)		48	0.00
hamming6-4	64	4	6*(BMTH)	1	9	0.00	8*(MT)		∞	0.00	10*(MT)		10	0.00	12*(T)		12	0.00
hamming8-2	256	128	128*(BMT)	,	128	0.09	128*(MT)	,	128	60.0	128(MT)	128	129	22.41	128(MT)		152	0.97
hamming8-4	256	16	16*(BMT)	í	16	0.01	20(T)	20	20	0.01	20(T)	24	25	0.28	20(T)	32	32	0.02
hamming 10-2	1024	512	512*(M)	,	512	8.97	512(M)	512	512	4.66	512(M)	512	512	8.62	512(M)	512	513(512.15)	16.82
johnson8-2-4	28	4	5*(BMTH)	1	rO	0.00	8*(MT)	,	×	0.00	9*(MT)	,	6	0.00	12*(T)	,	12	0.00
johnson8-4-4	20	14	14*(BMTH)	í	14	0.00	18*(TH)	,	18	00.00	22*(T)		22	0.00	24(T)	24	28	0.00
johnson16-2-4	120	∞	10*(T)	,	10	0.00	16(T)	16	16	0.00	19(T)	19	19	0.00	21(T)	24	24	0.00
keller4	171	11	15*(BMTH)	,	15	0.00	21*(T)	,	21	0.09	22(T)	23	23	90.0	22(T)	28	28	0.02
p-hat300-1	300	∞	10*(MTH)	1	10	0.00	12*(MT)		12	0.00	$14^{*}(T)$	,	14	0.01	14(T)	15	16	0.02
p-hat300-2	300	22	30*(T)		30	0.01	30(T)	36	36	0.02	33(M)	41	41	90.0	33(M)	46	46	0.03
p-hat300-3	300	36	43(B)	43	44	0.07	43(B)	25	52	0.06	43(B)	59	59	0.09	43(B)	65	65	0.11
p-hat500-1	200	o ;	$12^*(T)$	1	12	0.06	14*(T)	, ,	14	0.20	14(T)	16	16	0.11	14(T)	17	20 0	0.15
p-hat700-1	700	11	13*(M)	, [	13 10	0.06	13(M)	14	15	0.13	13(M)	15	17	0.45	13(M)	x 1	19	2.00
p-hat700-2	700	44	50(B)	51	22	0.06	50(B)	09	62	1.35	50(B)	× 1	70	0.26	50(B)	75	79	9.42
p-hat/00-3	000	7.0	73(B)	90	76	0.54	73(B)	1 00 0 00	0.00	2.13	(3(B)	97	100	1.46	(3(B)	106	109	1.35
san200_0.7_2 san200_0_9_1	200	10	24 (M) 90 (M)	*00	26(25.4U) 90	9.74	36(M) 125*(M)	, ,	37 195	0.02	48(M) 125(M)	105*	195	1.90	48(M) 125(M)	1.05*	125	0.02
hamming 10-4	1024	40	41(BM)	44	82	1.53	46(M)	53	64	1.13	51(M)	64	68(67.20)	20.64	51(M)	73	79(78,05)	34.37
brock800_1	800	23		21	25	10.90	-	26	29	12.35		29	34(33.20)	24.43	-	33	37	27.12
$brock800_{-2}$	800	24	,	22	25	11.36	,	27	30(29.30)	31.20	,	30	34(33.15)	26.40	,	33	<b>38</b> (37.15)	32.46
brock800_3	800	25	1	24	25	12.68	1	25	30(29.20)	11.71	1	30	34(33.15)	28.46	1	32	38(37.10)	36.78
$brock800_{-4}$	800	26	,	22	26(25.55)	56.21	,	26	29	14.35		31	33	32.04		33	37	33.58
C1000.9	1000	89		71	81(80.55)	39.65	,	84	95(93.75)	58.59		26	107(106.00)	48.91	1	108	119(118.15)	60.39
C2000.5	2000	16		14	<b>19</b> (18.95)	27.96	,	18	22(21.90)	62.35		21	25(24.50)	22.79	1	23	28(27.15)	11.72
C2000.9	2000	80	1	92	06(88)06	65.21	1	88	105(103.40)	69.41	1	94	118(116.80)	63.85	1	107	132(129.65)	26.66
C4000.5	4000	18		12	20	39.23		14	23	69.37		13	26(25.55)	53.64	1	16	29(28.20)	26.93
keller6	3361	59		25	63	3.60		65	90(87.80)	66.21		7.7	<b>107</b> (103.45)	67.88	1	06	<b>125</b> (123.20)	73.91
p-hat1000-1	1000	10		11	13	0.28		13	15	0.15		15	18	4.09		18	50	7.17
p-hat1000-2	1000	46		52	56	0.43	1	64	67	0.81		71	92	28.89	1	82	84	1.30
p-hat1000-3	1000	8 (		7.7	82	0.33		95	80 1	2.19		109	111	3.50		117	122	29.84
p-hat1500-1	1500	12		13	14	1.46		14	17	22.47		16	19	3.73	1	16	21	1.43
p-hat1500-2	1500	65		71	80	1.75	1	90	93	0.41		66	107(106.50)	29.46	1	113	<b>117</b> (116.55)	41.90
p-nat1500-3	1000	94		108	114	0.01	,	120	133	17.83		141	150	3.90	,	154	164	48.70
saniooo	Tonni	CI CI		1 ,	1,1	9.59	<u>.</u>	67	67	0.73		22	99	T. (0	_	4.1	41	0.39
Note a: 1	he value	of 44 r.	eported in [29]	tor M.	Note $a$ : The value of 44 reported in [29] for MANN-a9 is wrongly claimed to be the optimal solution.	ly claime	d to be the o	ptımaı	solution.	_								

#### 4.5 Impact of frequency information

As described in Sections 2.4 and 2.5, the construction procedure and perturbation procedure are guided by frequency information. In this section, we evaluate the effectiveness of this frequency strategy. We compared the original FD-TS algorithm with a variant, FD-TS-R, in which the frequency-based vertex selection rule was replaced by a random selection rule. In particular, to create a new solution, FD-TS-R randomly adds a vertex from  $M_1$  to the current solution (Section 2.4) and randomly selects a vertex from  $M'_3$  for perturbation (Algorithm 2, line 31).

To better differentiate FD-TS and FD-TS-R, we selected 27 instances from the three benchmark sets, such that the selected instances cover different characteristics (random vs real-world, dense vs sparse) and are sufficiently challenging based on the search effort required to attain the best solutions according to the results of Tables 1 and 2. For this experiment, both FD-TS and FD-TS-R were run 20 times to solve each instance, each run being limited to 20 seconds for the 2nd DIMACS instances and 180 seconds for the other (larger) instances. We compared the average objective values reached by both algorithms ("ave" columns), the average time required to first obtain the best objective value ("time" columns), and the improvement in the average objective value achieved by FD-TS as a percentage ("ave\_imp" column).

Table 3 shows the results achieved for s=2,3,4,5 respectively. A difference in the average solution quality obtained by the two algorithms was only observed with the 2nd DIMACS instances (the first 12 instances). For the large instances, both FD-TS and FD-TS-R converged so fast that the best solution was found quite early (the average time required to first obtain the best solution was less than one second in most cases). For the 2nd DIMACS instances, FD-TS achieved better solutions than FD-TS-R for 5, 6, 8, 8 instances with s=2,3,4,5, respectively (marked in bold font). In addition, for 4,4,4,2 instances with s=2,3,4,5, respectively, the average objective values found by FD-TS were worse than those found by FD-TS-R (marked in italic font). In general, there was a slight advantage when using the frequency mechanism, and it increased with s. This experiment confirms that the frequency mechanism is helpful for solving hard dense graphs that require persistent search efforts.

Table 3 Impact of frequency information - comparison between FD-TS and FD-TS-R.

			s=2					s=3			s=4			s=5	
instance	FD-,	FD-TS-R	FD-TS	SI	dai ove	FD-TS-R	-R	FD-TS	dai erre	FD-TS-R	FD-TS	ami eve	FD-TS-R	FD-TS	dani eve
	ave	time	ave	time	ave_mp	ave ti	time a	ave time		ave time	ave time	ave_imp	ave time	ave time	ave_mp
C1000.9	79.40	8.59	79.70	7.45	0.38%	92.70 7.	7.95	92.80 7.89	0.11%	104.95 5.18	105.10 6.68	0.14%	116.85 10.40	117.30 9.88	0.38%
C2000.5	18.45	3.25	18.50	5.05	0.27%	21.05 1.	1.29   2	21.10 2.71	0.24%	24.15 3.35	24.10 4.06	-0.21%	26.85 6.27	26.95 9.35	0.37%
C2000.9	87.80	6.97	87.85	9.92	0.06%	102.20 9.	9.50 1	101.70 7.97	-0.49%	115.05 8.65	114.85 9.61	-0.17%	128.25 9.76	127.70 8.95	-0.43%
C4000.5	19.60	4.80	19.25	2.75	-1.82%	22.35 3.	3.96 2	22.25 3.86	-0.45%	25.05 3.99	25.20 4.18	0.60%	27.90 4.81	27.90 6.42	0.0%
brock800_1	24.85	4.98	24.70	4.30	-0.61%	28.90 7.	7.04 2	28.95 5.38	0.17%	32.70 4.57	33.15 5.91	1.36%	36.40 3.93	36.50 4.00	0.27%
brock800_2	24.85	6.20	24.90	7.63	0.20%	29.00 7	7.62 2	28.95 5.08	-0.17%	32.80 5.74	32.75 5.48	-0.15%	36.45 5.56	36.65 6.22	0.55%
brock800_3	24.85	8.74	24.80	5.31	-0.20%	28.75 5	5.06   2	29.00 7.19	898.0	32.80 7.74	32.95 6.84	0.46%	36.30 3.60	36.45 5.27	0.41%
brock800_4	24.70	6.84	24.85	7.28	0.60%	28.65 5	$5.76 \mid 2$	28.70 5.99	0.17%	32.40 5.43	32.55 5.49	0.46%	36.25 4.71	36.30 3.07	0.14%
hamming10-4	48.00	1.36	48.00	1.16	0.0%	64.00 0.	0.98   6	64.00 1.12	0.0%	66.75 5.73	66.85 3.98	0.15%	77.50 3.92	77.45 3.05	-0.06%
keller6	63.00	2.98	63.00	3.33	0.0%	84.70 8	8.68	85.70 8.46	1.17%	99.50 9.00	98.95 11.69	-0.56%	121.00 9.89	121.40 14.53	0.33%
p_hat1500-2	80.00	1.66	80.00	2.02	0.0%	93.00 0	$0.47 \mid 9$	93.00 0.48	0.0%	106.00 1.40	106.05 1.50	0.05%	116.00 1.49	116.10 2.78	0.09%
san1000	16.80	4.73	16.75	2.50	-0.30%	25.00 3.	$3.30 \mid 2$	24.95 2.46	-0.20%	32.95 3.35	33.00 1.29	0.15%	41.00 3.59	41.00 3.82	0.0%
333SP	5.00	1.21	5.00	1.19	%0.0	0.00	0.71 6	6.00 1.04	0.0%	7.00 1.03	7.00 0.51	%0.0	8.00 0.96	8.00 0.46	0.0%
cage15	00.9	2.41	00.9	2.48	0.0%	8.00.8	8.89	8.00 5.68	0.0%	10.00 5.84	10.00 6.02	%0.0	11.00 34.77	11.00 26.71	0.0%
caidaRouterLevel	20.00	0.55	20.00	0.31	0.0%	23.00 0.	0.94   2	23.00 0.51	0.0%	24.00 0.81	24.00 0.74	0.0%	26.00 0.52	26.00 0.82	0.0%
cit-HepPh	24.00	0.32	24.00	0.37	0.0%	27.00 0.	$0.46 \mid 2$	27.00 0.27	0.0%	30.00 0.37	30.00 0.19	0.0%	32.00 0.34	32.00 0.30	0.0%
cit-HepTh	28.00	1.31	28.00	1.20	0.0%	31.00 1	1.51	31.00 0.70	0.0%	34.00 1.02	34.00 0.81	%0.0	37.00 1.39	37.00 0.72	0.0%
cit-Patents	17.00	15.56	17.00	21.25	0.0%	21.00 13	8.90  2	21.00 51.02	%0.0	26.00 9.55	26.00 16.23	%0.0	31.00 10.13	31.00 11.09	0.0%
email-EuAll	19.00	0.13	19.00	0.15	0.0%	22.00 0.	0.19   2	22.00 0.22	0.0%	25.00 0.17	25.00 0.22	0.0%	27.00 0.15	27.00 0.26	0.0%
p2p-Gnutella04	5.00	0.03	5.00	0.03	0.0%	7.00 0.7	0.08	60.0 00.2	0.0%	9.00 0.12	9.00 0.10	0.0%	10.00 0.00	10.00 0.01	0.0%
p2p-Gnutella24	5.00	0.04	5.00	0.07	0.0%	0.00	0.01   6	6.00 0.02	0.0%	8.00 0.12	8.00 0.10	0.0%	9.00 0.32	9.00 0.18	0.0%
p2p-Gnutella25	5.00	0.03	5.00	0.03	0.0%	0.00	0.01   6	6.00 0.01	0.0%	8.00 0.01	8.00 0.01	0.0%	10.00 0.08	10.00 0.10	0.0%
soc-Slashdot0811	31.00	90.0	31.00	0.02	0.0%	34.00 0	0.06	34.00 0.06	0.0%	38.00 0.13	38.00 0.12	0.0%	40.00 0.08	40.00 0.10	0.0%
soc-Slashdot0902	32.00	0.06	32.00	0.05	0.0%	35.00 0.	0.07	35.00 0.06	0.0%	40.00 0.11	40.00 0.10	0.0%	42.00 0.12	42.00 0.10	0.0%
web-NotreDame	155.00 1.21	1.21	155.00 0.74	0.74	0.0%	$155.00\ 1.30$		$155.00\ 0.71$	0.0%	155.00 1.22	155.00 0.84	0.0%	155.00 1.18	155.00 0.99	0.0%
wiki-Talk	32.00 1.40	1.40	32.00 1.18	1.18	0.0%	36.00 3	$3.41 \mid 3$	36.00 3.60	0.0%	41.00 2.68	41.00 1.51	0.0%	44.00 3.82	44.00 3.72	0.0%
wiki-Vote	21.00	0.07	21.00 0.10	0.10	0.0%	24.00 0.	0.13 2	24.00 0.11	0.0%	27.00 0.16	27.00 0.12	%0.0	28.00 0.08	28.00 0.13	0.0%

#### 5 Conclusions and perspectives

The NP-hard maximum s-plex problem is significant in both theory and practice. In this study, we proposed an effective local search algorithm for solving this problem heuristically based on the general tabu search method. To ensure its efficiency, the proposed algorithm combines a multi-neighborhood search procedure with vertex-moving frequency, where the search process is driven by two intensification oriented operators (Add and Swap) and one diversification operator (Press). Dedicated rules are defined to explore the neighborhoods introduced by these operators. Information regarding vertex moves is collected and used to guide the construction of the starting solutions and the perturbation process. A graph peeling technique is also integrated to dynamically reduce large sparse graphs.

We assessed the performance of the proposed algorithm using three popular benchmark sets: 47 instances from the Stanford Large Network Dataset Collection and the 10th DIMACS Implementation Challenge, and 52 dense graphs from the 2nd DIMACS Implementation Challenge. For the SNAP and 10th DIMACS benchmarks, FD-TS obtained improved solutions (new lower bounds) for 7, 15, 19, 20 instances when s=2,3,4,5, respectively. Moreover, many of these solutions were proved to be optimal using the *Peel* procedure. FD-TS also performed very well on the instances from the 2nd DIMACS benchmark set. The *Peel* procedure was no longer effective for these dense graphs, but FD-TS still obtained the current best-known results for all of the instances and discovered better solutions for most instances compared with four recent reference algorithms and the powerful CPLEX solver. Additional results for 48 more graphs from the above benchmark sets showed in the Appendix further demonstrated the performance of the proposed algorithm.

Several areas of research require further investigation. First, to achieve high search robustness across a large range of problem instances with very different characteristics, it would be useful to develop adaptive and learning techniques to help the algorithm to adjust its search strategies dynamically. Second, it would be interesting to explore other ways of employing frequency information to improve the performance of the algorithm. For instance, we could investigate frequency information in new selection rules for the transformation operators as well as other guided perturbation mechanisms such as that proposed in [4]. Finally, it would be interesting to adapt the ideas introduced in this work to design search algorithms for other clique-relaxations, such as s-defective clique [37], quasi-clique [1,22,31] and k-club [7,18,28,30].

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# A Computational results on additional instances

This Appendix includes additional results of our FD-TS algorithm and CPLEX for 48 instances from the three benchmark sets. Note that these instances have not been tested previously by any s-plex algorithm. Only lower bounds (from the best-known maximum clique sizes [32,34]) are available. Table A.1 contains the 20 large SNAP and 10th DIMACS instances while Table A.2 includes the remaining 28 instances of the 2nd DIMACS Challenge.

cplex 10 10\* Computational results of FD-TS on 20 large networks from the SNAP Collection and the 10th DIMACS Implementation Challenge 929820 952203 99982 119747 98961 491 19 20 22 75.77 0.10 0.00 24.9511.50time 0.02 0.04 0.25 1.38 0.81 0.00 0.02 0.09 2.53 5.29  $\frac{11*}{152}$  $\frac{18}{393*}$ 491\*222 \* \* 222 \* \* 252 \* \* 252 944\* 8 10\* 10 47 45 45 45 47 N/A 391\*20\* 875349 95220399982 119747 16174 0 98961 1115 20 22 944 12.16 42.96 47.05 0.02 0.24 21.8490.0 0.00 0.72 7.57 7.54 2.62 5.38 0.02 8 10\* 148 **16** 391 491\*20 22\* 23\* 24\* 944\* 9 6\* 111 21 21 6\* N/A 390\* N/A\*61 0155074 936015 952203 1000000 9779 99982 0 99734  $0 \\ 1115 \\ 20 \\ 22$ 944 0 34.83 25.33 44.23 11.17 0.03 0.02 0.53 0.92 0.00 0.85 0.01 0.540.02 2.82 5.64 6.909\* 144 491\*
19
21\*
22\*
23\* **14** 390 944\* ო \* 9 39 21 4 ω \* N/A N/A N/A 388\* N/AN/AN/A N/A 937779 952203 |V'| 12097 19765 119747 99982 0 98961 6731 944 12.5524.9445.00 0.04 0.24 0.89 2.49 4.73 3.51 2.83 5.67 0.00 3.95 0.02 0.020.681.08 5 5 8\* 140 13
388
490\*
19\*
20\*
22\*
22\*
944\* ა \* 4 7 136 10 387 489 119 20 21 21 21 31 6 5 2 5 36 2 2 2  $\frac{1156647}{16138468}$ 132557200 22785136 261787258 30359198 63501393  $\begin{array}{c} 147892 \\ 899792 \\ 2456071 \end{array}$ 13591473 14487995 501198 499985 38354076 10593311998000 119666499998  $18520486 \\ 100000 \\ 100000$ 8388608 16777216 943695 952203 1000000 268495 1382908 262111 65536 862664 2097152 4194304 156317 100000 114599kron\_g500-simple-logn16 citationCiteseer luxembourg.osm p2p-Gnutella30 p2p-Gnutella31 preferential Attachment G\_n\_pin\_pout Amazon0302 rgg-n-2-21-s0 rgg-n-2-23-s0 rgg\_n\_2\_24\_s0 rgg-n-2-22-s0 smallworld uk-2002audikw1 eu-2005in-2004

ecology1

Table A.2 Computational results of FD-TS on 28 benchmark instances of the 2nd DIMACS Implementation Challenge

	14.71			s=2			s=3			s=4			s=2	
instance	<u>-</u>	$\omega$ [34]	cplex	max(ave)	time	cplex	max(ave)	time	cplex	max(ave)	time	cplex	max(ave)	time
C125.9	125	34	43*	43	00.0	51*	51	1.89	58	58	0.07	65*	65	0.38
C250.9	250	44	53	55	8.43	63	65	22.29	75	75	4.81	84	84	4.72
C500.9	200	57	63	69	10.67	74	81(80.95)	57.72	84	<b>92</b> (91.75)	55.11	26	103(102.25)	36.65
DSJC1000-5	1000	15	15	18	26.61	18	21	25.35	21	<b>24</b> (23.05)	5.63	24	27(26.25)	32.14
DSJC500_5	200	13	15	16	0.29	17	19	2.81	20	21	0.12	23	24	1.07
MANN_a81	3321	1100	2162*	2162(2113.90)	139.35	3240*	3240(3125.35)	138.36	3240*	3240(2788.70)	147.51	3240*	3135(2660.75)	190.01
brock400_3	400	31	28	30	0.35	33	36	6.59	38	41	5.90	43	<b>46</b> (45.90)	55.95
gen200-p0.9-44	200	44	53	53	0.14	99	99	0.00	92	92	0.03	84	84	0.05
gen200-p0.9-55	200	55	22	57	0.02	64	64	0.14	73	73	0.12	80	80	0.19
gen400-p0.9-55	400	55	64	68(67.70)	65.76	82	87	26.62	108	112	0.19	124	124	0.16
gen400-p0.9-65	400	65	73	73(71.60)	28.01	100	101(100.45)	10.68	132	132	0.25	138	138	0.15
gen400-p0.9-75	400	75	78	79(78.05)	30.62	112	114	0.35	136	136	0.10	136	136	0.12
johnson32-2-4	496	16	21	21	0.02	32	32	0.02	38	38	0.38	48	48	0.09
keller5	922	27	31	31	0.09	41	45	8.19	46	<b>53</b> (52.75)	53.67	28	61	5.04
p_hat500-2	200	36	42	42	0.02	49	50	0.12	22	57	0.07	62	62	0.25
p-hat500-3	200	20	09	62	0.18	7.1	72	1.24	80	81	1.94	88	88	1.73
san200_0.7_1	200	30	31	31	0.59	46	46(45.70)	1.51	09	09	0.01	75*	75	0.01
san200-0.9-2	200	09	7.1	71	2.47	105*	105	0.03	105*	105	0.02	105*	105	0.03
san200_0.9_3	200	44	53	54(53.95)	72.31	73	73	7.48	*96	96	80.0	100*	100	0.03
san400-0.5-1	400	13	15	15	1.24	22	22	3.61	29	29	4.92	35	36(35.70)	55.40
san400_0.7_1	400	40	41	41	0.20	61	61	2.49	80	81(80.45)	17.54	100	100	90.0
san400_0.7_2	400	30	32	32	7.22	46	47(46.10)	0.46	61	61	0.57	92	92	10.34
san400_0.7_3	400	22	27	27(26.30)	12.27	38	38	11.17	50	50(49.45)	24.71	61	61	0.29
san400_0.9_1	400	100	102	102(101.30)	60.6	150	150	0.09	200*	200	0.12	*002	200	0.15
sanr200_0.7	200	18	22	22	0.01	25	26	0.03	30	30	0.14	33	33	0.02
sanr200-0.9	200	42	51	51	0.61	09	61	2.25	69	69	0.07	92	2.2	5.74
sanr400_0.5	400	13	14	15	0.02	18	18	0.09	20	21	0.56	23	24	1.63
sanr400_0.7	400	21	25	26	1.03	28	30	0.45	32	35	5.31	35	39	31.01