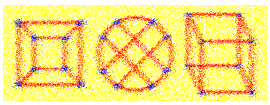
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Decycling Bipartite Graphs

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Abstract. Let $G = (V, E)$ be a graph and let $S \subseteq V$ be a subset of its vertices. If the subgraph of G induced by $V \setminus S$ is acyclic, then S is said to be a decycling set of G. The size of a smallest decycling set of G is called the decycling number of G. Determining the decycling number of a graph G is NP-hard, even if G is bipartite. We describe a tabu search procedure that generates decycling sets of small size for arbitrary bipartite graphs. Tests on challenging families of graphs show that the proposed algorithm improves many best-known solutions, thus closing or narrowing the gap to the best-known lower bounds.

1 Introduction

We consider the problem of eliminating all cycles from a graph by means of deletion of vertices. Determining a decycling set (i.e., a set of vertices whose removal eliminates all cycles) of minimum size is an NP-hard problem, even for bipartite graphs. We describe a heuristic procedure that produces small decycling sets in arbitrary bipartite graphs. Experiments performed on challenging families of bipartite graphs will demonstrate the efficiency of the proposed procedure. We first need to fix some notation to describe the problem more precisely.

Let $G = (V, E)$ be a simple undirected graph with vertex set V and edge set E. The *order* of G is its number $|V|$ of vertices. For a subset $W \subseteq V$ of vertices, we write $E(W)$ for the set of edges of G with both endpoints in W, while $G[W] = (W, E(W))$ is the subgraph of G induced by W. Also, we write $N_G(v)$ for the set of vertices adjacent to v in G and $d_G(v) = |N_G(v)|$ is the *degree* of v. A *stable* set of G is a set of pairwise non-adjacent vertices, while a *clique* of G is a set of pairwise adjacent vertices. A *forest* is an acyclic graph. A graph $G = (V, E)$ is bipartite if there is a partition (V_1, V_2) of V so that all edges of E have one endpoint in V_1 and the other in V_2, and we also write $G = (V_1, V_2, E)$. For other basic notions of graph theory that are not defined here, we refer to Diestel [10].

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Let W be a subset of vertices of a graph $G = (V, E)$. If the induced subgraph $G[V \setminus W]$ of G is acyclic, then W is said to be a decycling set of G. The smallest size of a decycling set of G is the *decycling number* (or feedback vertex number) of G and is denoted by $\phi(G)$. Let $\varphi(G)$ denote the largest order of an induced forest of $G = (V, E)$. Clearly, determining $\phi(G)$ is equivalent to computing $\varphi(G)$ since a subset $F \subseteq V$ of vertices induces a forest in G if and only if $V \setminus F$ is a decycling set of G, which implies $\phi(G) + \varphi(G) = |V|$.

Determining the decycling number of a graph has various applications, including deadlock recovery [27], synchronous distributed systems [22], VLSI design [12], constraint satisfaction and Bayesian inference [4]. Karp [17] has shown that the problem is NP-hard, even when restricted to planar graphs, bipartite graphs and perfect graphs. On the other hand, the problem is known to be solvable in polynomial time for various other families of graphs, including cubic graphs [19, 28], cocomparability graphs and convex bipartite graph [20]. A 2-approximation (i.e., a polynomial time algorithm that generates a decycling set of cardinality at most $2\phi(G)$) is described in [3], while a branch-and-cut algorithm for the exact solution of the problem is given in [8]. Local search algorithms are proposed in [8, 9, 23, 26] for determining an upper bound on the decycling number of arbitrary graphs. More results on the decycling number can be found in [5].

In a seminal paper on the topic, Beineke and Vandell [7] have bounded the decycling number of hypercubes. Improving these bounds for hypercubes was continued in [6, 13, 25]. Other families of graphs have then also been investigated, including Fibonacci cubes [11], bubble sort graphs [30] and star graphs [29]. All these graphs are bipartite and have nice topological properties that provide attractive interconnection schemes for massively parallel systems [1, 2]. The problem of avoiding deadlocks when using these topologies for exchanging data between processors can be reformulated as a decycling problem, and it turns out that computing the decycling number of these bipartite graphs is a real challenge. Indeed, there is a gap between the best-known lower and upper bounds on $\phi(G)$ for such graphs G with only 120 vertices. We aim to decrease these gaps by generating decycling sets of small size.

In the next section, we describe an algorithm that determines small decycling sets in arbitrary bipartite graphs. We first design in Section 2.1 a procedure that extends an induced forest $G[F]$ of G to a larger one $G[F'{=}F{\cup}S]$ with a stable set S of G. We then show in Section 2.2 how to embed such forest extensions into a tabu search. In Section 3, we demonstrate the efficiency of our algorithm by applying it on various challenging families of bipartite graphs. The proposed procedures offer flexibility at different levels, and we will discuss this in Section 4.

2 The proposed algorithm

As a first observation, note that if a vertex set induces a forest in G, then all its subsets also induce a forest in G. As a particular case, assume that F is the vertex set of an induced forest of G and that S is obtained from F by removing all vertices of degree at least 2 in $G[F]$ as well as one vertex per connected component of $G[F]$ with exactly 2 vertices. Then S is a stable set of G, and $G[F{\setminus}S]$ is a forest of G obtained from $G[F]$ by removing all isolated vertices as well as some of its leaves. We show in Section 2.1 how to use this property the other way around. In other words, given a subset F that induces a forest in G, we will extend it to a larger set $F'{=}F{\cup}S$ so that $G[F']$ also induces a forest in G, while imposing that the set S of added vertices is a stable set of G.

Given a forest $G[F]$ of G, finding a set F' of maximum cardinality so that $F' \supseteq F$ and $G[F']$ is also a forest of G is a difficult task. Indeed, for $F = \emptyset$, the problem is equivalent to determining $\varphi(G)$. We will focus on bipartite graphs G and on supersets F' of F obtained by adding a stable set S of G (i.e., $F' = F \cup S$). These assumptions make the problem a little easier. Indeed, while

finding a stable set of maximum cardinality in an arbitrary graph is an NP-hard problem, the famous Hungarian method [18] can be used when restricted to bipartite graphs. Hence, given a vertex set F that induces a forest $G[F]$ in a bipartite graph $G = (V, E)$, the following problem can be solved in polynomial time: find a stable set $S \subseteq V \setminus F$ of maximum cardinality such that every vertex in S is adjacent to at most one vertex in F. With such a stable set S we deduce that $G[F' = F \cup S]$ is a forest of G. It is obtained by adding leaves and isolated vertices to $G[F]$. We will go one step further by considering stable sets S of G that possibly contain vertices with more than one neighbor in F.

The next subsection gives more details on the proposed procedure that extends an induced forest $G[F]$ of G to a larger one $G[F' = F \cup S]$ with a stable set S of G. The next step will be to embed these forest extensions into a tabu seach. Roughly speaking, we are going to explore the space that contains all the induced forests of G that have fixed numbers of vertices in X and in Y, and for each considered forest, we will try to extend it by adding a stable set of G. This will be explained in detail in Section 2.2.

2.1 Forest extensions with stable sets

Let F be the vertex set of an induced forest of a graph $G = (V, E)$. Assume $G[F]$ has r connected components with vertex sets $C_1, C_2, \ldots, C_r$. We denote by V_F the set of vertices $v \in V \setminus F$ with at most one neighbor in every C_i:

$$V_F = \{v \in V \setminus F \text{ with } |N_G(v) \cap C_i| \leq 1 \text{ for } i = 1, 2, \ldots, r\}.$$

Clearly, adding a vertex $v \notin V_F$ to F creates a cycle. We therefore extend F by choosing vertices in V_F. We consider the partition $(V_F^{\leq 1}, V_F^{\geq 2})$ of V_F where $V_F^{\leq 1}$ is the subset of vertices in V_F with at most one neighbor in F and $V_F^{\geq 2}$ is the subset of vertices in V_F with at least two neighbors in F:

$$V_F^{\leq 1} = \{v \in V_F \text{ with } |N_G(v) \cap F| \leq 1\}, \quad \text{and}$$
$$V_F^{\geq 2} = \{v \in V_F \text{ with } |N_G(v) \cap F| \geq 2\}.$$

For illustration, consider the bipartite graph G at the top left of Figure 1 and assume $F = \{a, f, g, h, j, n, p, r\}$, which gives the forest $G[F]$ of G at the bottom left, with four connected components, the vertices of F becoming colored in black. Vertex b does not belong to V_F since its neighbors a and g belong to the same connected component of $G[F]$. Vertices d and k have no neighbor in F, while vertices c and e have one neighbor in F. Also, vertices l, m, o and q have a neighbor in two different connected components of $G[F]$, while vertex i has a neighbor in three connected components of $G[F]$. Hence, $V_F = \{c, d, e, i, k, l, m, o, q\}$ with $V_F^{\leq 1} = \{c, d, e, k\}$ and $V_F^{\geq 2} = \{i, l, m, o, q\}$.

We have observed that if F induces a forest in G and S is any stable set in $V_F^{\leq 1}$, then $F \cup S$ also induces a forest in G. Such an extension is similar to adding leaves and isolated vertices to $G[F]$. Adding one vertex $v \in V_F^{\geq 2}$ to F does not create any cycle in $G[F \cup \{v\}]$, while it reduces the number of connected components by $|N_G(v) \cap F| - 1$ units. Note however that adding more than one vertex of $V_F^{\geq 2}$ to F may create a cycle, even if the added vertices form a stable set of G. For example, if $G = (V, E)$ is a cycle on four vertices a, b, c, d and F contains two non-adjacent vertices of G, say a and c, then the two other vertices b and d are also non-adjacent and they both belong to $V_F^{\geq 2}$. Hence $G[F = \{a, c\}]$ is a forest and $S = \{b, d\}$ is a stable set of G, while $G[F \cup S] = G$ is not a forest of G.

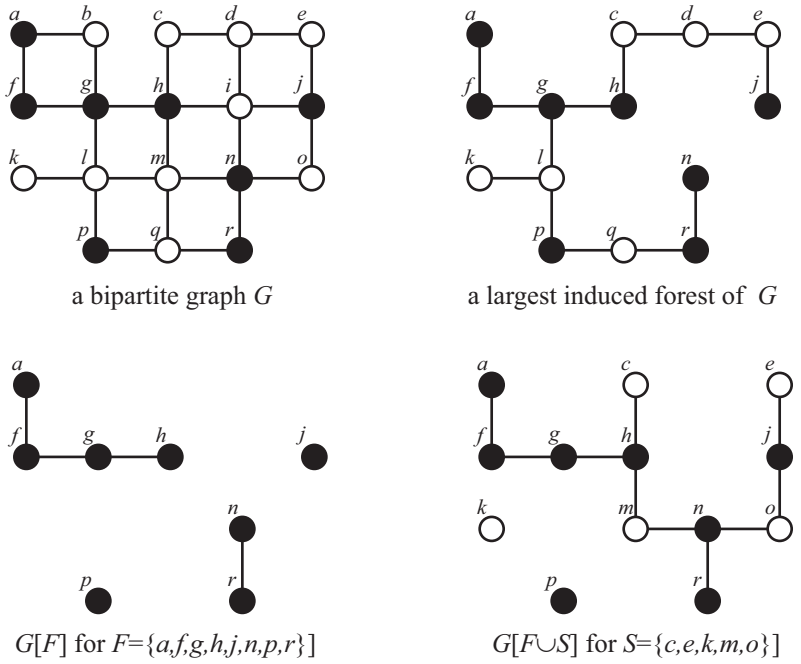

Figure 1: A bipartite graph $G = (X, Y, E)$ and three of its induced forests.

Consider again the example of Figure 1 with $F = \{a, f, g, h, j, n, p, r\}$. Clearly, $S = \{c, e, k\}$ is the only maximum stable set in $V_F^{\leq 1}$, which means that the largest forest $G[F \cup S]$ with $S \subseteq V_F^{\leq 1}$ has order 11. As shown at the bottom right of Figure 1, a forest $G[F \cup S]$ of order 13 can be obtained by choosing the stable set $S = \{c, e, k, m, o\}$ with c, e, k in $V_F^{\leq 1}$ and m, o in $V_F^{\geq 2}$. Observe also that $\{c, e, i, k, m, o, q\}$ is the only stable set of V_F with 6 vertices, but it contains vertices i and o, and the addition of these two vertices to F creates the cycle (i, j, o, n). The next Property indicates which stable sets S in V_F ensure that $G[F \cup S]$ is a forest of G.

Property 2.1 *Let F be the vertex set of an induced forest of a graph G, let S be a stable set in V_F, and let $S' = S \cap V_F^{\geq 2}$. Then $G[F \cup S]$ is a forest of G if and only if $G[F \cup S']$ is a forest of G.*

Proof: Since $G[F \cup S']$ is a subgraph of $G[F \cup S]$, we only have to prove that if $G[F \cup S']$ is a forest of G, then the same is true for $G[F \cup S]$. So assume $G[F \cup S']$ is a forest of G, and let $F' = F \cup S'$ and $S'' = S \cap V_F^{\leq 1} = S \setminus S'$. Then $V_{F'}^{\leq 1} = V_F^{\leq 1}$ since the vertices in S'' have no neighbor in S'. Hence, no vertex in S'' has more than one neighbor in F' and we have observed that this implies that $G[F' \cup S''] = G[F \cup S]$ is a forest of G. □

In the example of Figure 1, $G[F \cup \{c, e, k, m, o\}]$ is a forest of G since the same is true for $G[F \cup \{m, o\}]$. Also, $G[F \cup \{c, e, i, k, m, o\}]$ is not a forest of G since this is also not the case for $G[F \cup \{i, m, o\}]$.

Property 2.1 shows that when looking for a stable set S that extends a forest $G[F]$ of G, we may determine a stable set S' in $V_F^{\geq 2}$ such that $G[F \cup S']$ is a forest of G. We can then easily

determine a maximum stable set S'' among the vertices of $V_F^{\leq 1}$ that are not adjacent to any vertex of S', and this will give a stable set $S = S' \cup S''$ so that $G[F \cup S]$ is a forest of G.

We will restict our attention to stable sets S with $S \subseteq X$ or $S \subseteq Y$. So, let $G = (X, Y, E)$ be a bipartite graph. We first show how to determine a stable set S_X in $V_X = V_F^{\geq 2} \cap X$ so that $G[F \cup S_X]$ is a forest of G. As mentioned above, we assume that $G[F]$ has r connected components with vertex sets $C_1, C_2, \ldots, C_r$. We set $A = \{1, 2, \ldots, r\}$ and construct a bipartite graph $H_X(F)$ with bipartition (V_X, A) of its vertex set and with edge set E_X, where a vertex $v \in V_X$ is linked to a vertex $a \in A$ if and only if v has a neighbor in the connected component C_a of $G[F]$. We then look for a large subset S_X of V_X so that $S_X \cup A$ is a forest of $H_X(F)$. Clearly such a subset S_X is a stable set in $V_F^{\geq 2}$ so that $G[F \cup S_X]$ is a forest of G.

For the forest $G[F]$ of Figure 1, assuming that $a \in X$, $C_1 = \{a, f, g, h\}$, $C_2 = \{j\}$, $C_3 = \{n, r\}$, and $C_4 = \{p\}$, we have $H_X(F)$ equal to the graph at the top left of Figure 2. The removal of i from V_X deletes all cycles in $H_X(F)$. Hence, by setting $S_X = \{m, o\}$, we can conclude that $A \cup S_X$ induces a forest in $H_X(F)$, which means that $G[F \cup S_X]$ is a forest of G, as shown at the bottom left of Figure 2.

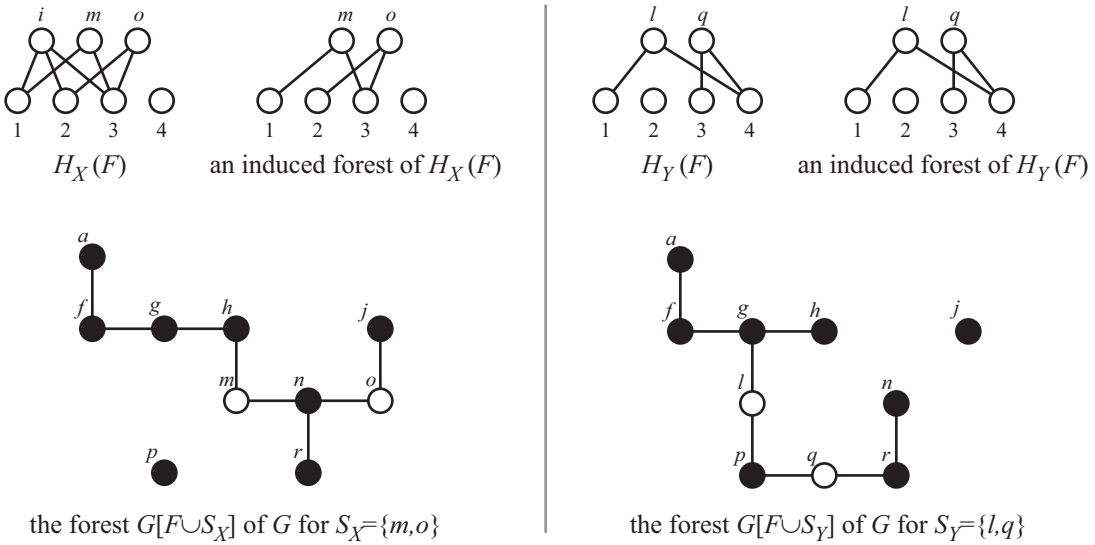

Figure 2: Two extensions of the forest $G[F]$ of Figure 1, with $a \in X$, $C_1 = \{a, f, g, h\}$, $C_2 = \{j\}$, $C_3 = \{n, r\}$, and $C_4 = \{p\}$.

Finding a stable set $S_X \subseteq V_X$ of maximum cardinality so that $S_X \cup A$ induces a forest in $H_X(F)$ is a difficult problem since it is equivalent to determining the smallest decycling set D of $H_X(F)$ with the additional constraint that $D \subseteq V_X$. Since the tabu search described in the next section will perform this task many times, we have decided to use a greedy algorithm that works as follows. We start by setting S_X equal to the empty set, and then iteratively increase S_X ensuring that $S_X \cup A$ induces a forest in $H_X(F)$. Let W be the set of vertices v of V_X so that $S_X \cup A \cup \{v\}$ induces a forest in $H_X(F)$. At each step of the iterative process, we choose a vertex $v \in W$ with minimum degree $d_{H_X(F)}(v)$, add it to S_X and update W. Hence, W is initially equal to V_X, and the algorithm stops when W is empty. Algorithm `GreedyStable` describes the general steps of this

algorithm. Its input is a bipartite graph $H = (X, Y, E)$ and its output is a subset $S \subseteq X$ so that $H[S \cup Y]$ is a forest of H. Hence, we will run it on $H_X(F) = (V_X, A, E_X)$.

Algorithm `GreedyStable`
Input : a bipartite graph $H = (X, Y, E)$
Output: a set $S \subseteq X$ so that $H[S \cup Y]$ is a forest of H

1: Set $S \leftarrow \emptyset$ and $W \leftarrow X$.
2: **while** $W \neq \emptyset$ **do**
3: Choose a vertex v of minimum degree $d_H(v)$ in W.
4: Set $S \leftarrow S \cup \{v\}$.
5: Remove v from W as well as all vertices u that belong to a cycle of $H[S \cup Y \cup \{u\}]$.
6: **end while**

When applied to the graph $H_X(F)$ of Figure 2, Algorithm `GreedyStable` determines S_X by first choosing m or o in $W = \{i, m, o\}$, say m. Then m is removed from W as well as i that belongs to the cycle $(1, i, 3, m)$. Vertex o is then added to the stable set and W becomes empty. The output is therefore $S_X = \{m, o\}$.

The same process can be repeated with Y instead of X. More precisely, we can construct the graph $H_Y(F)$ with vertex set $V_Y \cup A$ for $V_Y = V_F^{\geq 2} \cap Y$, and with edge set E_Y where a vertex $v \in V_Y$ is linked to a vertex $a \in A$ if and only if v has a neighbor in the connected component C_a of $G[F]$. We can then run Algorithm `GreedyStable` on $H_Y(F) = (V_Y, A, E_Y)$ to produce a stable set $S_Y \subseteq V_Y$ so that $G[F \cup S_Y]$ is a forest of G. As shown at the top right of Figure 2, $V_Y = \{l, q\}$ and $H_Y(F)$ is a forest, which means that $G[F \cup S_Y]$ is a forest of G for $S_Y = \{l, q\}$.

Note that $S_X \cup (V_F^{\leq 1} \cap V_X)$ is a stable set of $G = (X, Y, E)$ since all vertices of this set belong to the same part X of the bipartition (X, Y) of the vertex set of G. Hence, $G[F \cup S_X \cup (V_F^{\leq 1} \cap V_X)]$ is a forest of G. Similarly, $G[F \cup S]$ is a forest of G for the stable set $S = S_Y \cup (V_F^{\leq 1} \cap V_Y)$. For the forest $G[F]$ of Figure 1, we have observed that $S_X = \{m, o\}$ and $S_Y = \{l, q\}$. Also, $V_F^{\leq 1} \cap X = \{c, e, k\}$ and $V_F^{\leq 1} \cap Y = \{d\}$, which means that $G[F \cup \{c, e, k, m, o\}]$ and $G[F \cup \{d, l, q\}]$ are forests of G obtained from $G[F]$ by adding a stable set.

We are now ready to explain how we extend an induced forest $G[F]$ of a bipartite graph $G = (X, Y, E)$ by adding a stable set S so that $G[F \cup S]$ is also a forest of G. We first construct $H_X(F)$ and $H_Y(F)$ and then apply Algorithm `GreedyStable` to both bipartite graphs to obtain sets S_X and S_Y such that $G[F \cup S_X]$ and $G[F \cup S_Y]$ are forests of G. We then compare $n_X = |S_X| + |V_F^{\leq 1} \cap X|$ with $n_Y = |S_Y| + |V_F^{\leq 1} \cap Y|$. If $n_X \geq n_Y$, we set $F' = F \cup S_X \cup (V_F^{\leq 1} \cap X)$, otherwise we set $F' = F \cup S_Y \cup (V_F^{\leq 1} \cap Y)$. It follows from Property 2.1 that $G[F']$ is a forest of G. In our example, we have seen that $S_X = \{m, o\}$, $S_Y = \{l, q\}$, $V_F^{\leq 1} \cap X = \{c, e, k\}$ and $V_F^{\leq 1} \cap Y = \{d\}$. Hence $n_X = 2 + 3 = 5$ and $n_Y = 2 + 1 = 3$. We therefore add $\{c, e, k, m, n\}$ to F to obtain the forest drawn at the bottom right of Figure 1. Algorithm `ForestExtension` describes the general steps of this procedure.

Several remarks are worth making at this stage. Algorithm `ForestExtension` determines a stable set S with $S \subseteq X$ or $S \subseteq Y$ while a larger stable set S' could exist in $X \cup Y$ so that $G[F \cup S']$ is a forest of G. Also, we only look for forest extensions $G[F \cup S]$ of $G[F]$ obtained by adding a stable set S, while larger forests $G[F \cup W]$ of G might exist for non-stable sets W. For the example of Figure 1 a forest with 14 vertices is depicted at the top right. It is obtained by adding the non-stable set $\{c, d, e, k, l, q\}$ to F. But finding the largest set W so that $G[F \cup W]$ is a forest of G is an NP-hard problem since it is equivalent to determining $\varphi(G)$ when $F = \emptyset$.

Algorithm `ForestExtension`
Input : a bipartite graph $G = (X, Y, E)$ and a vertex set $F \subseteq X \cup Y$ so that $G[F]$ is a forest of G.
Output : a stable set S of G so that $G[F \cup S]$ is a forest of G.

1: Determine $V_F^{\leq 1}$ and $V_F^{\geq 2}$.
2: Construct $H_X(F)$ and $H_Y(F)$.
3: Run Algorithm `GreedyStable` on $H_X(F)$ and $H_Y(F)$ to determine S_X and S_Y so that $G[F \cup S_X]$ and $G[F \cup S_Y]$ are forests of G.
4: Set $n_X \leftarrow |S_X| + |V_F^{\leq 1} \cap X|$ and $n_Y \leftarrow |S_Y| + |V_F^{\leq 1} \cap Y|$.
5: **if** $n_X \geq n_Y$ **then** Set $S \leftarrow S_X \cup (V_F^{\leq 1} \cap X)$ **else** Set $S \leftarrow S_Y \cup (V_F^{\leq 1} \cap Y)$.

Note that if we look for stable sets S with both $S \cap X$ and $S \cap Y$ possibly not empty and with $S \cap V_F^{\geq 2} = \emptyset$, then finding the largest one such that $G[F \cup S]$ is a forest of G is an easy problem. Indeed, $G[F \cup S]$ is a forest of G for all stable sets $S \subseteq V_F^{\leq 1}$, and the problem of finding the largest stable set in $G[V_F^{\leq 1}]$ can be solved in polynomial time since G is bipartite. Continuing on this path, note that instead of adding $V_F^{\leq 1} \cap X$ to S_X or $V_F^{\leq 1} \cap Y$ to S_Y at step 5 of Algorithm `ForestExtension`, one could do the following. Once S_X is obtained with Algorithm `GreedyStable`, define $F' = F \cup S_X$ and determine $V_{F'}^{\leq 1}$. We have seen that $G[F']$ is a forest of G, and by adding a maximum stable set S' of $V_{F'}^{\leq 1}$, one can get a forest $G[F \cup (S_X \cup S')]$ of G. Note however that $S = S_X \cup S'$ is not necessarily a stable set of G. Similarly, another forest $G[F \cup (S_Y \cup S')]$ can be obtained by determining a maximum stable set S' in $V_{F \cup S_Y}^{\leq 1}$ and adding it to $F \cup S_Y$. For the forest $G[F]$ of Figure 1, we have seen that $S_Y = \{l, q\}$, and instead of adding $V_F^{\leq 1} \cap Y = \{d\}$ to $F \cup S_Y$, one can add the maximum stable set $S' = \{c, e, k\}$ in $V_{F \cup S_Y}^{\leq 1} = \{c, d, e, k\}$. We thus get an induced forest with 13 vertices which is obtained from $G[F]$ by adding the non-stable set $\{c, e, k, l, q\}$.

The variations of Algorithm `ForestExtension` mentioned above necessitate the solution of maximum stable set problems in bipartite graphs. Even if this task can be done in polynomial time, it is too time consuming for the tabu search procedure described in the next subsection, where Algorithm `ForestExtension` has to be applied many times.

2.2 A tabu search

The proposed heuristic for producing a large induced forest in a given bipartite graph G is based on the tabu search metaheuristic, which is one of the most frequently used in combinatorial optimization [15]. Tabu search is a local search technique that visits a solution space $\mathcal{S}$ by moving step by step from a current solution $s \in \mathcal{S}$ to a neighbor solution $s' \in \mathcal{N}(s)$, where $\mathcal{N}(s)$ is a subset of $\mathcal{S}$ called the neighborhood of s. A tabu list forbids some moves which would bring the search back to a recently visited solution. The best move that does not belong to the tabu list is chosen at each iteration. Tabu search was introduced by Glover [14]. A description of the method and its concepts can be found in [15].

The proposed adaptation of the tabu search metaheuristic to our problem can be roughly described as follows. We first choose a vertex set F so that $G[F]$ is a forest of $G = (X, Y, E)$. Such a set F can be obtained, for example, by means of the following greedy algorithm: start with an empty set F, consider all vertices v of G in non-decreasing degree order, and sequentially add them to F when $G[F \cup \{v\}]$ is a forest of G. But this initial set F can be chosen in many other different ways. For example, it can be the vertex set of the best-known induced forest of G that

we try to improve. We can also impose $F \cap X = X$ so that F contains all the vertices of X or, on the contrary, we can impose $F \subseteq Y$ so that F does not contains any vertex of X. There are actually no restrictions on F except that $G[F]$ must be a forest of G.

As mentioned earlier, the proposed tabu search explores the space $\mathcal{S}$ that contains all the induced forests of G which have fixed numbers of vertices in X and in Y, and for each considered forest, we will try to extend it by adding a stable set of G. Once an initial induced forest F is chosen, we set $k_X = |F \cap X|$, $k_Y = |F \cap Y|$ and therefore define the solution space $\mathcal{S}$ as the set of all vertex sets F' such that $|F' \cap X| = k_X$, $|F' \cap Y| = k_Y$, and $G[F']$ is a forest of G.

The neighborhood $\mathcal{N}(F)$ of $F \in \mathcal{S}$ contains all vertex sets $F' \in \mathcal{S}$ that can be obtained from F by replacing a vertex $u \in F \cap X$ by a vertex $v \in X \setminus F$, or by replacing a vertex $u \in F \cap Y$ by a vertex $v \in Y \setminus F$. Such a move from F to F' is denoted $F' = F \oplus (u, v)$. The value $f(F)$ of a set $F \in \mathcal{S}$ is measured by using Algorithm `ForestExtension` that produces a stable set S such that $G[F \cup S]$ is a forest of G. More precisely, we fix $f(F)$ equal to $|F| + |S|$, where S is the output of Algorithm `ForestExtension`. Note that $F = F' \oplus (v, u)$, and there is therefore a risk that the algorithm removes v and adds u in the next iterations to return to F, which could cause the algorithm to cycle. To avoid this, when moving from a set F to $F \oplus (u, v)$, we put u and v in a tabu list, which means that we forbid the visit of sets $F' \in \mathcal{S}$ with $u \in F'$ or $v \notin F'$. These restrictions are only effective for a limited number of iterations. It is customary to keep them active for a number of iterations proportional to the square root of the number of other possible choices. Let $Z \in \{X, Y\}$ be so that u and v both belong to Z. Since $|Z| - k_Z$ vertices (including u) of Z do not belong to F' while k_Z of them are part of it (including v), we keep u and v in the tabu list for $\sqrt{|Z| - k_Z}$ and $\sqrt{k_Z}$ iterations, respectively. The stopping criterion is a time limit. Algorithm `LargeForest` describes this tabu search, where $G[F^*]$ is the best found forest of G, *BestValue* is the largest value $f(F')$ of a non-tabu neighbor $F' \in \mathcal{N}(F)$, and *ListBest* is the set of non-tabu neighbors F' of F with $f(F') =$*BestValue*. At each iteration, the algorithm moves from a set F to a neighbor F' chosen at random in *ListBest*. Note that if a neighbor F' of F is found at step 10 with value $f(F') > |F^*|$, then the possible tabu status of the move from F to F' is cancelled (i.e., the move from F to $F' = F \oplus (u, v)$ is accepted, even if u or v belong to the tabu list). Also, when $f(F') > |F^*|$, we stop exploring the neighborhood $\mathcal{N}(F)$ and start a new iteration from F'.

Algorithm `LargeForest` can be considered as an upper bounding procedure for the computation of the decycling number of arbitrary bipartite graphs. Indeed, its output F^* indicates that $G[F^*]$ is a forest of G, which is equivalent to say that $V \setminus F^*$ is a decycling set of G. Hence, $|V| - |F^*|$ is an upper bound on the decycling number $\phi(G)$ of G. In the following, this upper bound will be denoted $UB(G)$.

3 Computational experiments.

There are very few hard instances in the literature for the computation of the decycling number of bipartite graphs. We found several exceptions, where the gap between the best-known lower and upper bounds is not zero, namely, star graphs, bubble sort graphs, Fibonacci cubes, and hypercubes. These families of bipartite graphs are our test sets for the next subsections. We use a 3 GHz Intel Xeon X5675 machine with 8 GB of RAM.

3.1 Star graphs

The star graph S_n has a vertex for every permutation $\boldsymbol{v} = v_1, v_2, \ldots, v_n$ of the integers $1, 2 \ldots, n$, and two vertices $\boldsymbol{v} = v_1, v_2, \ldots, v_n$ and $\boldsymbol{u} = u_1, u_2, \ldots, u_n$ are adjacent if there is $i \in \{2, \ldots, n\}$

Algorithm `LargeForest`
Input : a bipartite graph $G = (X, Y, E)$
Output : a set F^* such that $G[F^*]$ is a forest of G.

Choose a set F such that $G[F]$ is a forest of G
Set $k_X \leftarrow |F \cap X|$, $k_Y \leftarrow |F \cap Y|$, $F^* \leftarrow F$ and $TabuList \leftarrow \emptyset$.
while the time limit is not reached **do**
  Set $BestValue \leftarrow 0$ and $ListBest \leftarrow \emptyset$.
  for all u, v with $u \in F \cap X$ and $v \in X \setminus F$ or $u \in F \cap Y$ and $v \in Y \setminus F$ **do**
    Set $F' = F \oplus (u, v)$.
    if $G[F']$ is a forest of G **then**
      Apply Algorithm `ForestExtension` to determine a stable set S such that $G[F' \cup S]$ is a forest of G.
      Set $f(F') = |F'| + |S|$.
      if $f(F') > |F^*|$ **then**
        Set $F^* \leftarrow F' \cup S$, $F \leftarrow F'$ and go to step 4.
      else
        if $f(F') \geq BestValue$ and $\{u, v\} \cap TabuList = \emptyset$ **then**
          if $f(F') > BestValue$ **then**
            Set $BestValue \leftarrow f(F')$ and $ListBest \leftarrow \{F'\}$.
          else
            Add F' to $ListBest$.
          end if
        end if
      end if
    end if
  end for
  Choose a set $F' \in ListBest$ an let u, v be such that $F' = F \oplus (u, v)$.
  if u and v both belong to X **then** set $Z \leftarrow X$ **else** set $Z \leftarrow Y$ **end if**
  Keep u and v in $TabuList$ for $\sqrt{|Z| - k_Z}$ and $\sqrt{k_Z}$ iterations, respectively.
  Set $F \leftarrow F'$.
end while

such that $v_1 = u_i$, $v_i = u_1$, and $v_j = u_j$ for all $j \neq 1, i$. For illustration, S_3 and S_4 are depicted in Figure 3. Clearly S_n is bipartite and has $n!$ vertices and $\frac{n!(n-1)}{2}$ edges. Star graphs have many nice topological properties that provide attractive interconnection schemes for massively parallel systems [1, 2]. Their decycling number is studied in [29] where it is proved that $\phi(S_1) = \phi(S_2) = 0$, $\phi(S_3) = 1$, $\phi(S_4) = 7$, and

$$L(S_n) = \frac{n!(n-3)}{2(n-2)} + 1 \leq \phi(S_n) \leq \frac{1}{2}(n! - \sum_{i=2}^{n-1} i!) - 1 = U(S_n) \quad \forall n \geq 5.$$

Algorithm `LargeForest` starts with the generation of a vertex set F such that $G[F]$ is a forest of G, and then sets $k_X = |F \cap X|$ and $k_Y = |F \cap Y|$, which means that every vertex set in the solution space $\mathcal{S}$ of the tabu search has k_X vertices in X and k_Y vertices in Y. As explained in Section 2.2, this initial set F can be obtained in various ways.

Our initial choice was to produce the initial forest with the help of a greedy algorithm. For example, for S_7, we got a forest of order $2,258$. Much larger forests are however easy to obtain by imposing that F must contain one side, say Y, of the bipartition (X, Y) of the vertex set of S_n. This can be done as follows : use Algorithm `GreedyStable` on $S_n = (X, Y, E)$ to obtain a stable set

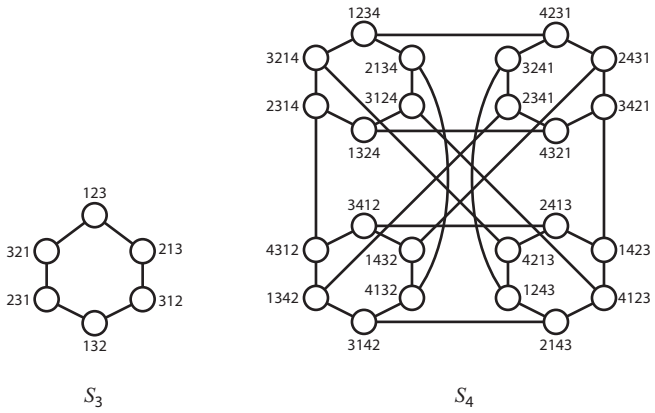

Figure 3: Star graphs S_3 and S_4.

$S \subseteq X$ so that $S_n[S \cup Y]$ is a forest of S_n, and then set $F = S \cup Y$. For S_7, this process has given a forest of order $2,931$, and we have therefore fixed $k_X = 2931 - \frac{1}{2}7! = 411$ and $k_Y = \frac{1}{2}7! = 2520$. For S_5 and S_6, the same initialization process resulted in forests of order 76 and 436, respectively. The results produced by Algorithm `LargeForest` with these initial forests are reported in Table 1. For every n, we indicate the order $|V|$ and the size $|E|$ of S_n, as well as the above-mentioned lower and upper bounds $L(S_n)$ and $U(S_n)$. For Algorithm `LargeForest`, we indicate the values of k_X and k_Y, and report the number of iterations (column 'Iter') and the CPU time (column 'time') needed to obtain the upper bound $UB(S_n)$ on $\phi(S_n)$. For S_7, we also indicate the time required to reach certain intermediate values. The complete evolution of the $UB(S_7)$ bound is shown in Figure 4.

			Bounds from [29]						
n	$\|V\|$	$\|E\|$	$L(S_n)$	$U(S_n)$					
1	1	0	0	0					
2	2	1	0	0					
3	6	6	1	1			`LargeForest`		
4	24	36	7	7	k_X	k_Y	$UB(S_6)$	Iter.	time
5	120	240	41	43	17	60	41	6	<1s.
6	720	1,800	271	283	76	360	271	209	3s.
7	5,040	15,120	2,017	2,083	411	2,520	2,100	9	2s.
							2,075	61	144s.
							2,050	557	1,850s.
							2,035	20,031	21h.49m.
							2,020	99,859	126h.48m.
							2,018	170,085	217h.27m.

Table 1: Result for star graphs S_n with $n \leq 7$.

We observe that for $n = 5$ and 6, our upper bound $UB(S_n)$ equals the lower bound $L(S_n)$, which means that we have proved that $\phi(S_5) = 41$ and $\phi(S_6) = 271$, and this proof is obtained in less than a second for S_5, and in 3 seconds for S_6. For S_7, the upper bound $U(S_7)$ is reached in 32 seconds. The algorithm was stopped after ten days of computation, and our upper bound $UB(S_7) = 2018$ is 65 units lower than that of Wang *et al.* [29], which represents a decrease of 98%

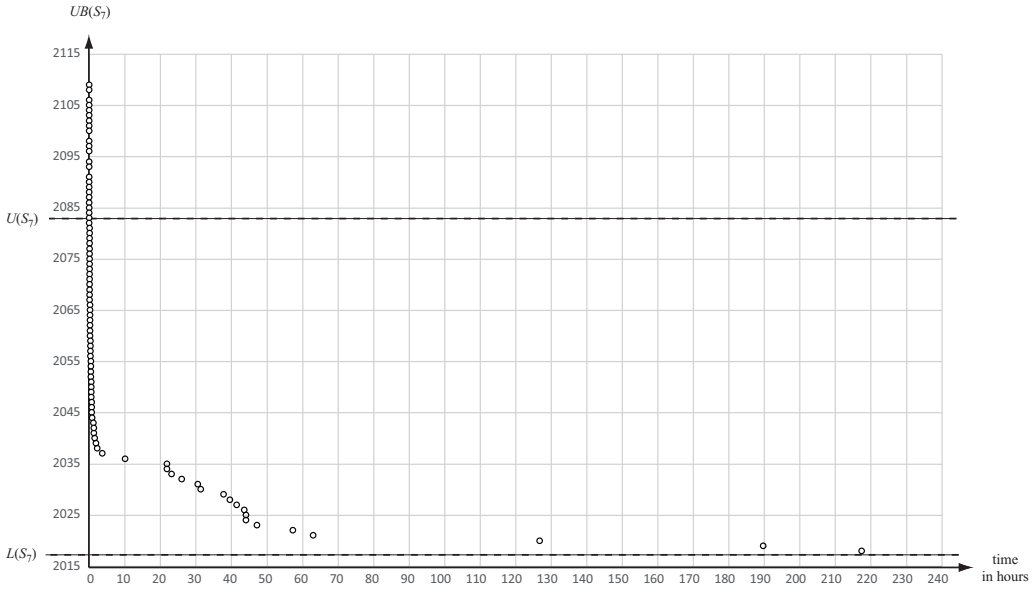

Figure 4: Results for S_7 with $k_X = 411$ and $k_Y = 2,520$.

of the gap. A smaller upper bound may have been reached if more time had been allocated. Note that $\phi(S_n) = L(S_n))$ for $n \leq 6$ which could mean that this is also true for larger values of n.

We then tried other values for k_X and k_Y. As explained in Section 2.2, knowledge of a large induced forest can be used to generate an initial set F at step 1 of Algorithm `LargeForest`. Such a large forest is known for S_n. Indeed, to demonstrate the validity of their upper bound $U(S_n)$ on $\phi(S_n)$, Wang *et al.* [29] have shown how to construct a decycling set D of $S_n = (X, Y, E)$ with $|D \cap X| = U(S_n)$ and $D \cap Y = \emptyset$. This means that $(X \setminus D) \cup Y$ induces a forest of order $n! - U(S_n)$ in S_n. Hence, by setting $F = X \setminus D$ at step 1 of Algorithm `LargeForest` and running `ForestExtension` with S_n and F as input, we immediately obtain a forest $S_n[F \cup Y]$ of S_n with $n! - U(S_n)$ vertices. In order to generate larger forests, we set $F = (X \setminus D) \cup R$, where R contains r vertices chosen at random in D, with $r \in \{1, \ldots, U(S_n)\}$. Hence, $k_X = \frac{1}{2}n! - U(S_n) + r$, $k_Y = 0$, and if Algorithm `LargeForest` succeeds in determining a forest $S_n[F^*]$ of S_n with $Y \subseteq F^*$, this means that we have found a forest of order $(\frac{1}{2}n! - U(S_n) + r) + \frac{1}{2}n! = n! - (U(S_n) - r)$, thus decreasing by r units the best-known upper bound on $\phi(S_n)$. We tested several values for r and the best results for S_7 were obtained with $r = 12$, which gives $k_X = 7! - 2083 + 12 = 449$ and $k_Y = 0$. With this setting, Algorithm `LargeForest` has reached the upper bound $U(S_7) = 2,083$ of Wang *et al.* [29] in 20 iterations that required one hour and 42 minutes of calculation. We were only able to improve it by 8 units in 2 days of calculation, thus producing a bound which is 57 units higher than that obtained with the other setting. These tests are therefore not reported in Table 1.

An analysis of the computing times shows that the first iterations take less than a second (because, at step 11 of `LargeForest`, we stop exploring $\mathcal{N}(F)$ when a neighbor F' of F is found with value $f(F') > |F^*|$), while subsequent iterations (which require $O(k_X(|X|-k_X)+k_Y(|Y|-k_Y))$ calls to Algorithm `ForestExtension`) can take a few seconds. Computing times are shorter with

$k_X = 411$ and $k_Y = 2520$ than with $k_X = 449$ and $k_Y = 0$, the reason being that with the first setting, Algorithm `ForestExtension` has to determine stable sets S having only a few tens of vertices, while with the second setting, S typically contains about 2500 vertices.

Still dealing with S_7, we have tried to initialize F by choosing 449 vertices at random in Y, without taking into account the decycling set described in [29]. This also gives $k_X = 449$ and $k_Y = 0$. With this other start, Algorithm `LargeForest` has produced an initial forest of $2,662$ vertices and therefore a decycling set of $2,378$ vertices. It would have taken days of computation to decrease this initial upper bound by 360 units (to get an upper bound equal to $2378 - 360 = 2018$). These tests are therefore not reported in Table 1.

3.2 Bubble sort graphs

The bubble sort graph B_n has the same vertex set as the star graph S_n, but not the same edge set. More precisely, there is a vertex in B_n for every permutation $\boldsymbol{v} = v_1, v_2, \ldots, v_n$ of the integers $1, 2 \ldots, n$, and two vertices $\boldsymbol{v} = v_1, v_2, \ldots, v_n$ and $\boldsymbol{u} = u_1, u_2, \ldots, u_n$ are adjacent if there is $i \in \{1, 2, \ldots, n-1\}$ such that $v_i = u_{i+1}$, $v_{i+1} = u_i$, and $v_j = u_j$ for all $j \neq i, i+1$. For illustration, B_3 and B_4 are depicted in Figure 5. Clearly B_n is bipartite and has $n!$ vertices and $\frac{n!(n-1)}{2}$ edges. Bubble sort graphs were first introduced by Akers and Krishnamurthy [2] as a new type of interconnection network. As mentioned in [30], bubble sort graphs are attractive because of their simple, symmetric and recursive structure. The decycling number of these graphs is studied in [30] and the authors have proved that $\phi(B_1) = \phi(B_2) = 0$, $\phi(B_3) = 1$, $\phi(B_4) = 7$, and

$$L(B_n) = \frac{n!(n-3)}{2(n-2)} + 1 \leq \phi(B_n) \leq \frac{n!(2n-3)}{4(n-1)} = U(B_n) \quad \forall n \geq 5.$$

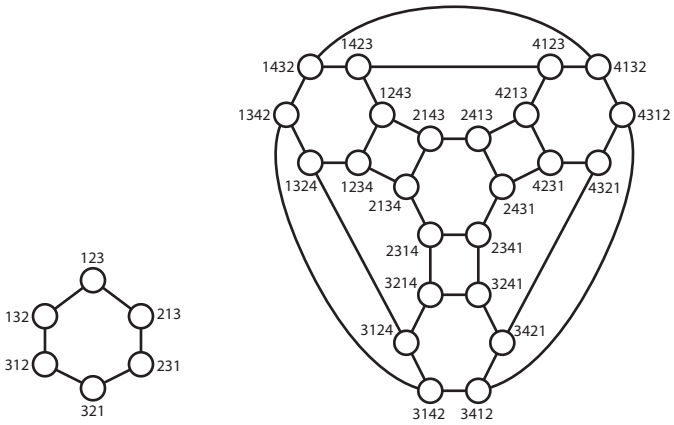

Figure 5: Bubble sort graphs B_3 and B_4.

Algorithm `LargeForest` is applied to B_5, B_6 and B_7, with the initialization process that produced the best results for star graphs. More precisely, Algorithm `GreedyStable` is run with $B_n = (X, Y, E)$ as input to obtain a stable set $S \subseteq X$ so that $B_n[S \cup Y]$ is a forest of B_n, and we then set $F = S \cup Y$. For B_7, this process has given a forest of order $2,932$, and we have therefore

fixed $k_X = 2932 - \frac{1}{2}7! = 412$ and $k_Y = \frac{1}{2}7! = 2520$. For B_5 and B_6, the same initialization process resulted in forests of order 76 and 437, respectively. The results produced by Algorithm `LargeForest` with these initial forests are reported in Table 2, with the same column labels as in Table 1.

			Bounds from [30]						
n	$\|V\|$	$\|E\|$	$L(B_n)$	$U(B_n)$					
1	1	0	0	0					
2	2	1	0	0					
3	6	6	1	1			`LargeForest`		
4	24	36	7	7	k_X	k_Y	$UB(B_n)$	Iter.	time
5	120	240	41	52	16	60	41	8	<1s.
6	720	1,800	271	324	77	360	272	796	14s.
							271	42,563	889s.
7	5,040	15,120	2,017	2,310	412	2,520	2,100	7	14s.
							2,075	55	474s.
							2,050	767	1h.31m.
							2,040	4,570	7h.46m.
							2,036	14,315	21h.10m.

Table 2: Results for bubble sort graphs B_n with $n \leq 7$.

We observe that for $n = 5$ and 6, our upper bound $UB(B_n)$ equals the lower bound $L(B_n)$, which means that we have proved that $\phi(B_5) = 41$ and $\phi(B_6) = 271$. For B_7, the process was stopped after one day of computation, and we could improve the best-known upper bound by 274 units, which represents a decrease of 94% of the gap. A smaller upper bound may have been reached if more time had been allocated. Note that $\phi(B_n) = L(B_n))$ for $n \leq 6$ which could mean that this is also true for larger values of n.

3.3 Fibonacci cubes

The Fibonacci numbers form a sequence of positive integers $\{f_n\}_{n=0}^{\infty}$ such that each number f_n is the sum of the two preceding ones, starting from $f_0 = 1$ and $f_1 = 2$. All non-negative integers i such that $i \leq f_n - 1$ can be uniquely represented as a sum of distinct non-consecutive Fibonacci numbers in the form $i = \sum_{j=0}^{n-1} b_j f_j$, where b_j is either 0 or 1, for $0 \leq j \leq n-1$, with the condition $b_j b_{j+1} = 0$ for $0 \leq j < n-1$ [32]. The sequence $(b_{n-1}, \ldots, b_1, b_0)$ is called the order-n Fibonacci code of i, and uniquely determines i. For example, $i = 19 \leq f_6 - 1 = 20$ has order-6 Fibonacci code $(1, 0, 1, 0, 0, 1)$. The Fibonacci cube of order n, denoted Γ_n, has vertex set $\{0, 1, \ldots, f_n - 1\}$, and two vertices i and j are adjacent if and only if their order-n Fibonacci codes differ in exactly one bit. The Fibonacci cubes for $n \leq 4$ are depicted in Figure 6.

Fibonacci cubes were first introduced by Hsu [16], and properties are described in [11, 21, 24, 31]. The decycling number of Γ_n is known for $n \leq 9$. For larger n, lower and upper bounds on $\phi(\Gamma_n)$ were obtained using various techniques. The best-known bounds $L(\Gamma_n)$ and $U(\Gamma_n)$ appear in [11] and are reported in Table 3 for $n \leq 14$.

Let (X, Y) be the bipartition of the vertex set of Γ_n, and assume without loss of generality that $|X| \leq |Y|$. It is easy to check that $|X| = |Y|$ or $|Y| - 1$.

Algorithm `LargeForest` was first tested with the initialization process of the previous sections, where the forests generated at step 1 contain all vertices of Y (i.e., $k_Y = |Y|$) and some of X. The results are reported in Table 3. We observe that the best-known upper bound $U(\Gamma_n)$ is easily

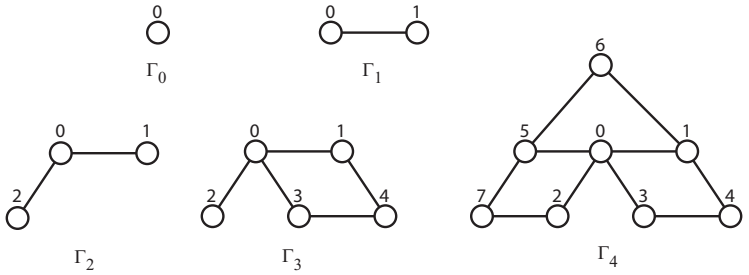

Figure 6: Fibonacci cubes Γ_n for $n \leq 4$.

			Bounds from [11]						
n	$\|V\|$	$\|E\|$	$L(\Gamma_n)$	$U(\Gamma_n)$					
1	2	1	0	0					
2	3	2	0	0					
3	5	5	1	1					
4	8	10	1	1					
5	13	20	3	3					
6	21	38	6	6					
7	34	71	11	11					
8	55	130	19	19	`LargeForest`				
9	89	235	33	33	k_X	k_Y	$UB(\Gamma_n)$	Iter.	time
10	144	420	53	55	13	72	55	58	<1s.
11	233	744	86	94	19	116	94	70	<1s.
12	377	1,308	139	158	26	189	157	51	1s.
13	610	2,285	225	264	37	305	259	134	7s.
14	987	3,970	364	439	62	494	439	5	<1s.
							427	1,494	31s.
							426	4,696	160s.
					54	0	439	80	213s.
					66	0	427	865	3,340s.
					67	0	426	5,217	5h16m.

Table 3: Result for Fibonacci cubes Γ_n with $n \leq 14$.

reached for all n, and we have improved it by one unit for $n = 12$, by 5 units for $n = 13$, and by 13 units for $n = 14$.

As done for star graphs, we have tested another initialization process by fixing $k_X = |X| - U(\Gamma_n) + r$ and $k_Y = 0$ with $r > 0$, but without taking into account the structure of the known forest of order $n! - U(\Gamma_n)$. In other words, our initial set F contains k_X vertices chosen at random in X. If Algorithm `LargeForest` succeeds in determining a forest $\Gamma_n[F^*]$ of Γ_n with $Y \subseteq F^*$, this means that we have found a forest of order $k_X + |Y| = (|X| - U(\Gamma_n) + r) + |Y| = f_n - (U(\Gamma_n) - r)$, thus decreasing by r units the best-known upper bound on $\phi(\Gamma_n)$. We can then stop the algorithm (since it cannot produce a forest with more that $k_X + \max\{|X|, |Y|\} = k_X + |Y|$ vertices) and rerun it with a larger value of r. The results produced with this setting are reported in Table 3 for some values of r. We observe that we reach the same bounds as with the previous setting, but in more time. As already indicated in Section 3.1, this can be explained by the fact that each iteration takes more time with $k_Y = 0$ than with $k_Y = |Y|$ since the stable sets S added to the forests F are larger with $k_Y = 0$.

3.4 Hypercubes

As final experiments, we illustrate how knowledge of the structure of the considered bipartite graph can help speed up the proposed upper bounding procedure. The n-dimensional hypercube Q_n is the graph with vertex set $\{0,1\}^n$, where two vertices $\boldsymbol{v}$ and $\boldsymbol{u}$ are linked by an edge if and only if there is $i \in \{1,2,\ldots,n\}$ such that $v_i = 1 - u_i$ and $v_j = u_j$ for all $j \neq i$. In what follows, we consider the bipartition (X,Y) of the vertex set of Q_n so that $\boldsymbol{v}$ belongs to X if and only if $\sum_{i=1}^{n} v_i$ is even.

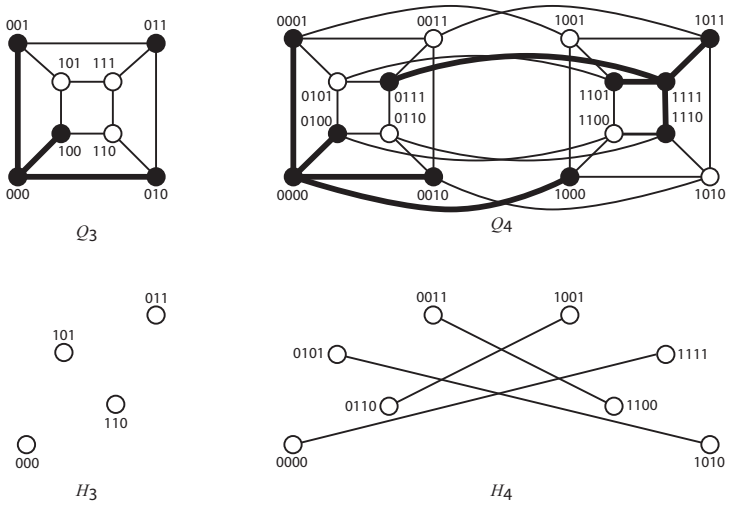

Figure 7: Hypercubes Q_3 and Q_4 with their associated graphs H_3 and H_4 and optimal induced forests represented with bold lines and black vertices.

The decycling number of Q_n is known for $n \leq 8$. As observed in [7], the best-known forests $G[F]$ of Q_n ($n \leq 8$) are unions of stars with n leaves plus some isolated vertices. For illustration optimal forests of Q_3 and Q_4 are depicted in Figure 7. A deeper analysis of these optimal forests $G[F]$ shows that all centers of stars belong to the same part, say X, of the partition (X,Y) of the vertex set of Q_n, and all vertices of Y are in $G[F]$. For example the largest induced forest of Q_3, depicted in Figure 7, contains all vertices of Y and 000 as only center of a star. For Q_4, the depicted largest induced forest contains all vertices of Y as well as 0000 and 1111 as centers of stars.

Assuming that optimal induced forests of Q_n have the above structure, we can speed up our upper bounding procedure by imposing that F must be a subset of X. Moreover, if $G[F \cup Y]$ is a union of stars, then all pairs of vertices in F must be at distance at least 4 in Q_n. So, let H_n be the graph with vertex set X, and where two vertices are linked by an edge if and only if their distance in Q_n is at least 4. For illustration, H_3 and H_4 are depicted in Figure 7. It follows that if F is a clique in H_n, then $G[F \cup Y]$ is a union of $|F|$ stars in Q_n. The following algorithm, called `CliqueSearch` is an adaptation of Algorithm `LargeForest`. It aims to determine a clique K of maximum cardinality in H_n, which will imply that $G[F^* = K \cup Y]$ is a union of stars (and therefore also a forest) of G. In this algorithm, $E(F)$ is the set of edges linking two vertices of F in H_n, which means that F is a clique in H_n if and only if $|E(F)| = \frac{|F|(|F|-1)}{2}$.

Algorithm `CliqueSearch` was tested on Q_n with $9 \leq n \leq 13$. The results are reported in Table 4, with the same column labels as in the previous tables. The best-known lower and upper bounds $L(Q_n)$ and $U(Q_n)$ for Q_n are taken from [25]. The upper bound $U(Q_n)$ is based on the theory of error-correcting codes, where

$$U(Q_n) = 2^{n-1} - 2^{n-r-1} \text{ with } n = 2^r - x \text{ and } 0 \leq x < 2^{r-1}.$$

We also report the upper bounds published one year earlier in [5] to demonstrate that Pike's work [25] has considerably reduced the value of these bounds.

Algorithm `CliqueSearch`
Input : graph H_n.
Output : a clique K in H_n so that $G[K \cup Y]$ is a forest of Q_n.

1: Generate a maximal clique F in H_n and set $K \leftarrow F$ and $k = |F| + 1$.
2: Choose a vertex $u \in X \setminus F$ and set $F \leftarrow F \cup \{u\}$, $F^* \leftarrow F$ and $TabuList \leftarrow \emptyset$.
3: **while** the time limit is not reached **do**
4: Set $BestValue \leftarrow 0$ and $ListBest \leftarrow \emptyset$.
5: **for** all $u \in F$ and $v \in X \setminus F$ **do**
6: Set $F' = F \oplus (u, v)$.
7: Set $f(F') = |E(F')|$.
8: **if** $f(F') > |F^*|$ **then**
9: Set $F^* \leftarrow F'$ and go to step 22.
10: **else**
11: **if** $f(F') \geq BestValue$ and $\{u, v\} \cap TabuList = \emptyset$ **then**
12: **if** $f(F') > BestValue$ **then**
13: Set $BestValue \leftarrow f(F')$ and $ListBest \leftarrow \{F'\}$.
14: **else**
15: Add F' to $ListBest$.
16: **end if**
17: **end if**
18: **end if**
19: **end for**
20: Choose a set $F' \in ListBest$ an let u, v be such that $F' = F \oplus (u, v)$.
21: Keep u and v in $TabuList$ for $\sqrt{|X| - k}$ and $\sqrt{k}$ iterations, respectively.
22: Set $F \leftarrow F'$.
23: **if** $|E(F)| = \frac{|F|(|F|-1)}{2}$ **then**
24: Set $K \leftarrow F$ and $k \leftarrow k + 1$.
25: Choose a vertex $u \in X \setminus F$ and set $F \leftarrow F \cup \{u\}$ and $F^* \leftarrow F$.
26: **end if**
27: **end while**

We observe that Algorithm `CliqueSearch` reaches the best-known upper bounds in at most 1 second. Note that Q_{13} has $8,192$ vertices and $53,248$ edges, which means that hours or days of computations would probably have been necessary to reach the same bounds with the original version of Algorithm `LargeForest`.

4 Concluding remarks

We have described a tabu search procedure that determines decycling sets of small size in arbitrary bipartite graphs G. The algorithm explores the set of induced forest of G of fixed order and tries

			Bounds from [25]					
n	$\|V\|$	$\|E\|$	$L(Q_n)$	$U(Q_n)$				
1	2	1	0	0				
2	4	4	1	1				
3	8	12	3	3				
4	16	32	6	6				
5	32	80	14	14				
6	64	192	28	28				
7	128	448	56	56	Upper bounds	`CliqueSearch`		
8	256	1,024	112	112	from [5]	$UB(Q_n)$	Iter.	time
9	512	2,304	225	236	237	236	89	<1s.
10	1,024	5,120	456	472	493	472	903	1s.
11	2,048	11,264	922	952	1,005	952	212	1s.
12	4,096	24,576	1,862	1,904	2,029	1,904	605	1s.
13	8,192	53,248	3,755	3,840	4,077	3,840	2,082	1s.

Table 4: Result for hypercubes Q_n with $n \leq 13$.

to extend these forests to larger ones by adding stable sets. We have shown that the procedure is effective on challenging families of bipartite graphs. Indeed, we have decreased most of the best-known upper bounds on $\phi(G)$, thus narrowing the gap to the best-known lower bounds. On four occasions we even managed to close this gap, which allows us to state that $\phi(B_5) = \phi(S_5) = 41$ and $\phi(B_6) = \phi(S_6) = 271$. The updated bounds for the tested graphs with a strictly positive gap are as follows, the old upper bounds being shown in parentheses:

$$\begin{array}{rcccl} 2017 & \leq & \phi(S_7) & \leq & 2018 \quad (2083) \\ 2017 & \leq & \phi(B_7) & \leq & 2036 \quad (2310) \\ 139 & \leq & \phi(\Gamma_{12}) & \leq & 157 \quad (158) \\ 225 & \leq & \phi(\Gamma_{13}) & \leq & 259 \quad (264) \\ 364 & \leq & \phi(\Gamma_{14}) & \leq & 426 \quad (439). \end{array}$$

Most of the largest forests produced by our algorithm contain all vertices of one part of the bipartition (X, Y) of the vertex set of G. Assuming $|X| \leq |Y|$, this has led us to fix $k_Y = 0$ or $k_Y = |Y|$, which helps speed up the algorithm since the moves from a solution F to a solution $F' = F \oplus (u, v)$ are restricted to pairs u, v of vertices that both belong to X. Such forests correspond to decycling sets fully contained in X. Notice however that it is easy to construct bipartite graphs $G = (X, Y, E)$ with no decycling set of minimum size fully contained in X or Y. In the example of Figure 8, the only optimal decycling set contains two adjacent vertices x and y which therefore belong to different parts of the bipartition of the vertex set. Algorithm `LargeForest` is flexible enough to allow the discovery of any induced forest of G (and therefore of any decycling set of G). Indeed, a forest $G[F^*]$ of G can be generated with $k_X = |F^* \cap X|$ and $k_Y \leq |F^* \cap Y|$: Algorithm `LargeForest` has to determine a set F so that $F \cap X = F^* \cap X$, $F \cap Y \subseteq F^* \cap Y$ and $G[F]$ is a forest of G, and Algorithm `GreedyStable` can then extend F to F^* by adding the stable set $F^* \setminus F \subseteq Y$.

We have observed in Section 3.4 that knowledge of the structure of the considered bipartite graph can help speed up the proposed upper bounding procedure. For example, several authors have realized that optimal forests in Q_n are unions of stars. With this assumption, we have been able to match the best-known upper bounds for $\phi(Q_n)$ ($n \leq 13$) in at most one second, while these graphs have up to $8,192$ vertices. Hence, an analysis of the structure of the best-known induced forests of a bipartite graph can facilitate the generation of large induced forests.

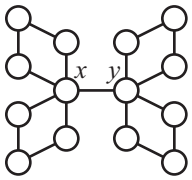

Figure 8: A bipartite graph with adjacent vertices in all optimal decycling sets.

Every iteration of Algorithm `LargeForest` requires $O(k_X(|X| - k_X) + k_Y(|Y| - k_Y))$ calls to Algorithm `ForestExtension`. In order to speed up the algorithm, we can fix an upper limit k_{max} on k_X and k_Y. If this limit has a too low value, it can prevent Algorithm `LargeForest` from generating an induced forest of G of maximum order. For example, if $k_{max} = 0$, then the output F^* of Algorithm `LargeForest` will be a stable set with $\max\{|X|, |Y|\}$ vertices. For the bubble sort graph B_7, this would lead to an upper bound of $2,520$ on $\phi(B_7)$ while it is very easy to generate decycling sets of B_7 with less than $2,100$ vertices. If an optimal forest has x vertices in X and y vertices in Y, then k_{max} should at least be equal to $\min\{x, y\}$. In summary, small values for k_{max} make it possible to speed up Algorithm `LargeForest`, but finding the smallest value of k_{max} which makes it possible to generate an optimal induced forest is a real challenge.

We want to mention that we did not seek to optimize the code by using structures making it possible to efficiently manage the fusion of connected components when adding a vertex to a forest, or their splitting when removing a vertex. This would certainly reduce the computing times somewhat. The aim of this article was to demonstrate that the proposed algorithm makes it possible to generate large induced forests in bipartite graphs and thus obtain small decycling sets.

Note finally that the smallest decycling sets generated by our algorithm for star graphs, bubble sort graphs, and Fibonacci cubes are archived online at `www.gerad.ca/~alainh/decycling.html`.

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