

# Minimum Activation Cost Edge-Disjoint Paths in Graphs with Bounded Tree-Width

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**Abstract.** In *activation network* problems we are given a directed or undirected graph  $G = (V, E)$  with a family  $\{f_{uv} : (u, v) \in E\}$  of monotone non-decreasing activation functions from  $D^2$  to  $\{0, 1\}$ , where  $D$  is a constant-size subset of the non-negative real numbers, and the goal is to find activation values  $x_v \in D$  for all  $v \in V$  of minimum total cost  $\sum_{v \in V} x_v$  such that the activated set of edges satisfies some connectivity requirements. We propose an algorithm that optimally solves the *minimum activation cost of  $k$  edge-disjoint st-paths* (*st-MAEDP*) problem in  $O(|V||D|^{tw+1}tw^3(k+1)^{(tw+3)^{2(tw+3)}}(tw+3)^{2(tw+3)+3})$  time for graphs with treewidth bounded by a constant  $tw$ .

## 1 Introduction

The *activation network* setting can be defined as follows. We are given a directed or undirected graph  $G = (V, E)$  together with a family  $\{f_{uv} : (u, v) \in E\}$  of monotone non-decreasing activation functions from  $D^2$  to  $\{0, 1\}$ , where  $D$  is a constant-size subset of the non-negative real numbers, such that the activation of an edge depends on the chosen values from the domain  $D$  at its endpoints. An edge  $(u, v) \in E$  is *activated* for chosen values  $x_u$  and  $x_v$  if  $f_{uv}(x_u, x_v) = 1$ , and the activation function  $f_{uv}$  is called *monotone non-decreasing* if  $f_{uv}(x_u, x_v) = 1$  implies  $f_{uv}(y_u, y_v) = 1$  for any  $y_u \geq x_u$ ,  $y_v \geq x_v$ . The objective of activation network problems is to find activation values  $x_v \in D$  for all  $v \in V$  such that the total activation cost  $\sum_{v \in V} x_v$  is minimized and the activated set of edges satisfies some connectivity requirements. Activation problems generalize several problems studied in the network literature such as power optimization, minimum broadcast tree and installation cost optimization. Several activation problems have been studied for arbitrary graphs in the recent research literature. Unfortunately, many of these problems are computationally hard. Obtaining a polynomial-time approximation algorithm is a typical approach to deal with an NP-hard problem. Another important approach is studying the problem on graph classes with nice decomposition properties such as bounded treewidth graphs to determine efficient algorithms. Panigrahi [8] shows that the fundamental problem of finding *minimum activation  $k$  edge-disjoint st-paths* (*st-MAEDP*) is NP-hard. It is an interesting area of research to investigate this problem for graphs with treewidth bounded by a constant  $tw$ . In this paper, we focus on developing a

polynomial-time algorithm that optimally solves the *st*-MAEDP for graphs of bounded treewidth. (Note that  $k$  is not a constant but part of the input.)

*Related Work.* Activation network problems were first introduced by Panigrahi in [8] and various of these problems have been investigated in [1, 2, 6–8]. The *minimum activation st-path* (MAP) problem is optimally solvable in polynomial-time [8]. However, the *minimum spanning activation tree* (MSPAT) problem is NP-hard to approximate within a factor of  $o(\log n)$  and there exists a  $O(\log n)$ -approximation algorithm for this problem [8]. Activation network problems generalize several problems studied in the network literature such as power optimization problems, which are modelled by a graph  $G = (V, E)$  where each edge  $(u, v)$  in  $G$  is assigned a threshold power requirement  $\theta_{uv}$ . For undirected graphs, an edge  $(u, v)$  is activated for chosen values  $x_u$  and  $x_v$  if each of these values is at least  $\theta_{uv}$ , and it is activated if  $x_u \geq \theta_{uv}$  for the directed case. The *minimum power st-path* (MPP) problem can be solved in polynomial time for both directed and undirected graphs [5]. For directed graphs, the *minimum power  $k$  node-disjoint st-paths* problem is also optimally solvable in polynomial time [4, 11]. However, the *minimum power  $k$  edge-disjoint st-paths* problem is unlikely to admit even a polylogarithmic approximation algorithm for both the directed and undirected variants [4].

There is a  $2k$ -approximation algorithm for the *st*-MAEDP problem and a 2-approximation algorithm for the *minimum activation node-disjoint st-paths* (*st*-MANDP) problem [6]. In [1], we have studied the *st*-MAEDP and *st*-MANDP problems when  $k = 2$ , denoted by *st*-MA2EDP and *st*-MA2NDP respectively. We proved that a  $\rho$ -approximation algorithm for the *st*-MA2NDP problem implies a  $\rho$ -approximation algorithm for the *st*-MA2EDP problem. In the same paper, we obtained a 1.5-approximation algorithm for the *st*-MA2NDP problem and hence for the *st*-MA2EDP problem. The problems *st*-MAEDP and *st*-MANDP for the restricted version of activation networks with  $|D| = 2$  and a single activation function for all edges have also been studied in [1]. Under this restriction, the *st*-MANDP problem is optimally solvable in polynomial time for arbitrary  $k$  (except for one case of the activation function, in which we require  $k = 2$ ). The *st*-MAEDP problem, however, remains NP-hard [1, 8]. So far these problems for an arbitrary constant-size  $D$  and fixed  $k \geq 2$  are not known to be NP-hard. Recently, in [2], we considered the *st*-MANDP and the problem of finding *minimum activation cost node-disjoint paths* (MANDP) between  $k$  disjoint terminal pairs,  $(s_1, t_1), \dots, (s_k, t_k)$ , for graphs of bounded treewidth. We proposed algorithms that optimally solve the *st*-MANDP problem in polynomial-time and the MANDP problem in linear-time for graphs with bounded treewidth [2]. There exists a polynomial-time algorithm that optimally solves the *st*-MA2EDP problem for graphs of bounded treewidth [1, 2]. Other relevant work, applications and motivations of activation network problems have been addressed in [1, 2, 6–8].

*Our Results.* We develop a polynomial-time algorithm for the *st*-MAEDP problem for graphs with treewidth bounded by  $tw$ . Our algorithm efficiently combines an edge-coloring scheme with dynamic programming over a *nice tree-decomposition* (see Sect. 2). The edge-coloring scheme was introduced in [12] to

develop a polynomial-time algorithm that solves the *minimum shared-edge  $kst$ -paths* (MSEP) problem (finding  $k$  paths between  $s$  and  $t$  with minimum number of shared edges among the paths) for graphs of bounded treewidth. The MSEP is a generalization of the problem of finding  $k$  edge disjoint  $st$ -paths.

The rest of the paper is organized as follows. In Sect. 2 we recall some definitions and results of the class of graphs with bounded treewidth. Then in Sect. 3, we present a polynomial-time algorithm that solves the  $st$ -MAEDP problem optimally for graphs with treewidth bounded by  $tw$ . We conclude the paper by stating some open problems in Sect. 4.

## 2 Preliminaries

In this paper we consider the class of graphs of bounded treewidth. A graph  $G = (V, E)$  has treewidth  $tw$  if it has a *tree-decomposition* of width  $tw$  [9]. The *tree-decomposition* concept is defined as follows.

**Definition 1.** Given a graph  $G = (V, E)$ , a tree  $\mathcal{T} = (I, F)$  and a family  $\mathcal{X} = \{X_i\}_{i \in I}$  of subsets of  $V$  (called *bags*). The pair  $(\mathcal{X}, \mathcal{T})$  is called a *tree-decomposition* of  $G$  if it satisfies the following conditions:

- $V = \bigcup_{i \in I} X_i$ .
- For every edge  $(v, w) \in E$ , there exists an  $i \in I$  with  $v \in X_i$  and  $w \in X_i$ .
- For every vertex  $v \in V$ , the nodes  $i \in I$  with  $v \in X_i$  form a subtree of  $\mathcal{T}$ .

The width of  $(\mathcal{X}, \mathcal{T})$  is the number  $\max_{i \in I} |X_i| - 1$ . The treewidth  $tw$  of the graph  $G$  is the minimum width among all possible tree-decompositions of the graph.

**Theorem 1** ([3]). For any fixed  $tw$ , there exists a linear-time algorithm that checks whether a given graph  $G = (V, E)$  has treewidth at most  $tw$ , and if so, outputs a tree-decomposition  $(\mathcal{X}, \mathcal{T})$  of  $G$  with width at most  $tw$ .

**Definition 2.** A tree-decomposition  $(\mathcal{X}, \mathcal{T})$  is called a *nice tree-decomposition*, if  $\mathcal{T}$  is a binary tree rooted at some  $r \in I$  that satisfies the following:

- Each node is either a leaf, or has exactly one or two children.
- Let  $i \in I$  be a leaf. Then  $X_i = \{u, v\}$  for some  $(u, v) \in E$ .
- For every edge  $(u, v) \in E$ , there is exactly one leaf  $i \in I$  such that  $u, v \in X_i$ . (We say that the edge  $(u, v)$  is associated with that leaf  $i \in I$ ).
- Let  $j \in I$  be the only child of  $i \in I$ , then either  $X_i = X_j \cup \{v\}$  or  $X_i = X_j \setminus \{v\}$ . The node  $i$  is called an *introduce node* or *forget node*, respectively.
- Let  $j, j' \in I$  be the two child nodes of a node  $i \in I$ , then  $X_j = X_{j'} = X_i$ . The node  $i$  is called a *join node* of  $\mathcal{T}$ .

Scheffler presented in [10] a special tree-decomposition that follows the structure of a nice tree-decomposition as defined above but with no restriction on the size of leaf bags (i.e., it does not require leaf bags to be of size 2). We call that type of tree-decomposition a Scheffler-type nice tree-decomposition. Any given tree-decomposition for a graph  $G = (V, E)$  with treewidth at most  $tw$  can be

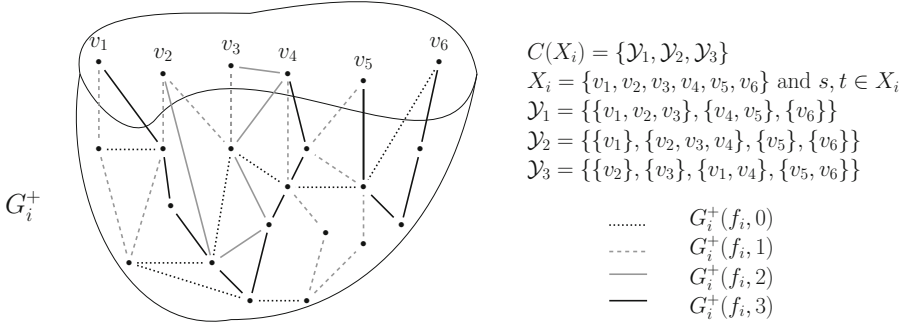
easily converted into a Scheffler-type nice tree-decomposition of width  $tw$  and size  $O(|V|)$  in linear time, if  $tw$  is a fixed constant [10]. One can also transform a Scheffler-type nice tree-decomposition to have leaves with bags of size 2 in linear time. Therefore, a nice tree-decomposition with leaf bags of size 2 and width  $tw$  can be constructed from any given tree-decomposition of the same treewidth. However, each leaf node of a Scheffler-type nice tree-decomposition may produce  $O(tw^3)$  new nodes in the construction of leaves with size 2. The resulting tree-decomposition, therefore, has  $O(|V|tw^3)$  nodes.

Throughout this paper, we consider simple undirected graphs  $G = (V, E)$  given as an input with a nice tree-decomposition of width at most  $tw$ . We define  $X_i^+$  to be the set of all vertices in  $X_j$  for all nodes  $j \in I$  such that  $j = i$  or  $j$  is a descendant of  $i$ . We denote by  $G_i^+$  a partial graph of  $G$ . For a leaf node  $i$ ,  $G_i^+$  is the subgraph of  $G$  with vertex set  $X_i$  and the edge of  $G$  that is associated with  $i$ . For a non-leaf node  $i$ ,  $G_i^+$  is the graph that is the union of  $G_j^+$  over all children  $j$  of  $i$ . Note that the graph  $G_r^+$  for the root  $r$  of the tree-decomposition is equal to  $G$ .

### 3 Minimum Activation Cost $k$ Edge-Disjoint $st$ -Paths

Given are an activation network  $G = (V, E)$  and a pair of source and destination vertices  $s, t \in V$ . In this section, we consider the  $st$ -MAEDP problem where the goal is to find activation values  $\{x_v : v \in V\}$  of minimum total cost  $\sum_{v \in V} x_v$  such that the activated set of edges contains  $k$  edge-disjoint  $st$ -paths  $\mathcal{P}^{st} = \mathcal{P}_1, \dots, \mathcal{P}_k$ . We present a polynomial-time algorithm that solves the  $st$ -MAEDP problem optimally in the case of graphs of bounded tree-width using dynamic programming techniques. The algorithm follows a bottom-up approach to compute a number of possible sub-solutions per nice tree-decomposition node  $i \in I$ . It is easy to compute the sub-solutions for a leaf node  $i \in I$  because the partial graph  $G_i^+$  consists of two vertices that are connected by an edge in  $G$  associated with  $i$ . For a non-leaf node, we use the information previously computed for its children. The algorithm also constructs a table  $tab_i$  to store the computed information for each node  $i \in I$ .

We use an edge-coloring scheme to compute the sub-solutions per tree node. Let  $C = \{0, 1, \dots, k\}$  be a set of colors. We consider a coloring  $f_i : E(G_i^+) \rightarrow C$  for the graph  $G_i^+$ . For each color  $c \in C$ , we define  $G_i^+(f_i, c)$  to be the subgraph of  $G_i^+$  induced by the edges with color  $c$ . Each color  $c \in C \setminus \{0\}$  represents the edges used by  $\mathcal{P}_c$ . Denote by  $P(X)$  the set of all possible partitions of the set  $X$ . We define  $C(X_i) = (\mathcal{Y}_1, \dots, \mathcal{Y}_k)$  to be a *color vector* on  $X_i$  such that  $\mathcal{Y}_c \in P(X_i \cup \{s, t\})$  for all  $c \in C \setminus \{0\}$ .  $P_{st}(X_i)$  denotes the set  $P(X_i \cup \{s, t\})$ . A color vector  $C(X_i) = (\mathcal{Y}_1, \dots, \mathcal{Y}_k)$  on  $X_i$  is called *active* if  $G_i^+$  has a coloring  $f_i$  such that every element of the partition  $\mathcal{Y}_c$ , for each  $c \in C \setminus \{0\}$ , is a set resulting from the intersection between  $X_i \cup \{s, t\}$  and the vertex set of a connected component of  $G_i^+(f_i, c)$  (see Fig. 1 for an example of an active color vector).  $\mathcal{Y}(X_i, f_i, c)$  denotes the partition  $\mathcal{Y}_c$  of  $X_i \cup \{s, t\}$ . This color vector concept was introduced in [12] for developing a polynomial-time algorithm that optimally



**Fig. 1.** An example of an active color vector

solves the MSEP problem for graphs with bounded treewidth. One simple way to compute the sub-solutions is by storing a color vector and an activation-value function per each row of table  $tab_i$ . However, the number of color vectors per node  $i$  can only be bounded by  $(tw + 3)^{k(tw+3)}$  because  $|X_i| \leq tw + 1$ ,  $|X_i \cup \{s, t\}| \leq tw + 3$  and  $|P_{st}(X_i)| \leq (tw + 3)^{tw+3}$ . Therefore, the number of color vectors is not always polynomially bounded. Hence the size of the table  $tab_i$  is also not always polynomially bounded. Therefore we define a mapping  $\gamma_i : P_{st}(X_i) \rightarrow \{0, 1, \dots, k\}$  to be a *count* on  $X_i$  to obtain a polynomial-time algorithm to optimally solve the  $st$ -MAEDP problem. We say that the count  $\gamma_i$  on  $X_i$  is an *active count* if  $G_i^+$  has a coloring  $f_i$  with a color vector  $C(X_i) = (\mathcal{Y}_1, \dots, \mathcal{Y}_k)$  such that  $\gamma_i(A) = |\{c \in C \setminus \{0\} : A = \mathcal{Y}_c\}|$  for each  $A \in P_{st}(X_i)$ . For any active count  $\gamma_i$ , it is clear that  $\sum_{A \in P_{st}(X_i)} \gamma_i(A) = k$ . In Fig. 1, the counts would be 1 for all partitions  $A \in \{\mathcal{Y}_1, \mathcal{Y}_2, \mathcal{Y}_3\} \subseteq P_{st}(X_i)$  and 0 otherwise. Consider a solution  $\mathcal{P} = \mathcal{P}_1, \dots, \mathcal{P}_k$  for the  $st$ -MAEDP problem. Let  $\mathcal{P}^i = \mathcal{P}[G_i^+]$  be the induced solution in a partial graph  $G_i^+$  (i.e., the set of vertices and edges that are both in  $\mathcal{P}$  and in  $G_i^+$ ). Since the interaction between  $\mathcal{P}^i$  in  $G_i^+$  and the rest of the graph happens only in vertices of  $X_i$ , we can consider an activation-value function  $\Lambda_i : X_i \rightarrow D$  and a count  $\gamma_i : P_{st}(X_i) \rightarrow \{0, 1, \dots, k\}$  to represent  $\mathcal{P}^i$  in  $G_i^+$ . The idea of using counts instead of color vectors is based on [12].

### 3.1 Processing the Tree Decomposition

For a tree node  $i \in I$ , the table  $tab_i$  has multiple rows and each row represents a solution of a unique combination of a count  $\gamma_i$  and a function of activation values  $\Lambda_i$ . Let  $val(\gamma_i, \Lambda_i)$  denote the minimum cost value of an assignment of activation values for  $G_i^+$  which satisfies the restrictions  $\Lambda_i$  and activates an edge-colored subgraph of  $G_i^+$  that satisfies the count  $\gamma_i$ . The value  $val(\gamma_i, \Lambda_i)$  is also stored in  $tab_i$ . We compute the sub-solution tables starting at the leaves towards the root.

*Leaf.* Let  $i \in I$  be a leaf,  $X_i = \{u, v\}$ . Let  $(\gamma_i, \Lambda_i)$  be any row of  $tab_i$ . We distinguish the following cases and define the value  $val(\gamma_i, \Lambda_i)$  for each case.

Each case corresponds to a possible sub-solution in  $G_i^+$ . Recall that  $G_i^+$  is a single edge. The sub-solution's cost  $val(\gamma_i, \Lambda_i)$  is set to  $\sum_{v \in X_i} \Lambda_i(v)$  if one of the following cases applies:

- $\gamma_i(A') = k - 1$  for  $A' = \{\{u\}, \{v\}, \{s\}, \{t\}\}$  and  $\gamma_i(A'') = 1$  for  $A'' = \{\{u, v\}, \{s\}, \{t\}\}$  and  $\gamma_i(A) = 0$  for all  $A \in P_{st}(X_i) \setminus \{A', A''\}$ , where  $f_{uv}(\Lambda_i(u), \Lambda_i(v)) = 1$  and  $s, t \notin X_i$ . Intuitively, this means that the sub-solution is a path with one edge not containing  $s$  or  $t$ .
- $\gamma_i(A') = k - 1$  for  $A' = \{\{u\}, \{v\}\}$  and  $\gamma_i(A'') = 1$  for  $A'' = \{\{u, v\}\}$  and  $\gamma_i(A) = 0$  for all  $A \in P_{st}(X_i) \setminus \{A', A''\}$ , where  $f_{uv}(\Lambda_i(u), \Lambda_i(v)) = 1$  and  $s, t \in X_i$ . Intuitively, this means that the sub-solution is a path with one edge containing  $s$  and  $t$ .
- $\gamma_i(A') = k - 1$  for  $A' = \{\{u\}, \{v\}, \{s\}\}$  and  $\gamma_i(A'') = 1$  for  $A'' = \{\{u, v\}, \{s\}\}$  and  $\gamma_i(A) = 0$  for all  $A \in P_{st}(X_i) \setminus \{A', A''\}$ , where  $f_{uv}(\Lambda_i(u), \Lambda_i(v)) = 1$  and  $s \notin X_i$  and  $t \in X_i$ . Intuitively, the sub-solution is a path with one edge and one endpoint equal to  $t$ . (The roles of  $s$  and  $t$  can be exchanged.)
- $\gamma_i(A') = k$  for  $A' = \{\{u\}, \{v\}, \{s\}, \{t\}\}$  and  $\gamma_i(A) = 0$  for all  $A \in P_{st}(X_i) \setminus \{A'\}$ , where  $s, t \notin X_i$ . Intuitively, this means that the sub-solution has no edges.
- $\gamma_i(A') = k$  for  $A' = \{\{u\}, \{v\}\}$  and  $\gamma_i(A) = 0$  for all  $A \in P_{st}(X_i) \setminus \{A'\}$ , where  $s, t \in X_i$ . Intuitively, this means that the sub-solution has no edges.
- $\gamma_i(A') = k$  for  $A' = \{\{u\}, \{v\}, \{s\}\}$  and  $\gamma_i(A) = 0$  for all  $A \in P_{st}(X_i) \setminus \{A'\}$ , where  $s \notin X_i$  and  $t \in X_i$ . Intuitively, this means that the sub-solution has no edges. (The roles of  $s$  and  $t$  can be exchanged.)

In these cases we construct an edge-colored subgraph of  $G_i^+$  plus activation-values that may be part of a global solution. If none of the above cases applies,  $val(\gamma_i, \Lambda_i) = +\infty$ .

*Introduce.* Let  $i \in I$  be an introduce node, and  $j \in I$  its only child. We have  $X_j \subset X_i$ ,  $|X_i| = |X_j| + 1$  and let  $v$  be the additional vertex in  $X_i$ . The vertex  $v$  is isolated since  $i$  does not introduce edges in  $G_i^+$ . For every row  $(\gamma_j, \Lambda_j)$  in  $tab_j$ , there are  $|D|$  rows in  $tab_i$  such that for all  $A \in P_{st}(X_j)$  and all  $u \in X_i \setminus \{v\}$ ,  $\gamma_i(A \cup \{\{v\}\}) = \gamma_j(A)$  if  $v \notin \{s, t\}$  and  $\gamma_i(A) = \gamma_j(A)$  if  $v \in \{s, t\}$  and  $\Lambda_i(u) = \Lambda_j(u)$ . The sub-solution's cost  $val(\gamma_i, \Lambda_i)$  for these rows is set to  $val(\gamma_j, \Lambda_j) + \Lambda_i(v)$ .

*Forget.* Let  $i \in I$  be a forget node, and  $j \in I$  its only child. We have  $X_i \subset X_j$ ,  $|X_j| = |X_i| + 1$  and let  $v$  be the discarded vertex. Let  $A^{-v}$  be the partition  $A \in P_{st}(X_j)$  after removing the vertex  $v$  if  $v \notin \{s, t\}$  and  $A^{-v}$  equals the partition  $A$  if  $v \in \{s, t\}$ . Note that  $A^{-v} \in P_{st}(X_i)$  for all  $A \in P_{st}(X_j)$ . For  $A \in P_{st}(X_i)$ , we say that  $W(A)$  is the set of all partitions  $B \in P_{st}(X_j)$  such that  $B^{-v} = A$  (i.e.,  $W(A) = \{B \in P_{st}(X_j) : B^{-v} = A\}$ ). For each row  $(\gamma_i, \Lambda_i)$ , we consider all  $(\gamma_j, \Lambda_j)$  such that for all  $u \in X_i$  and  $A \in P_{st}(X_i)$ ,  $\Lambda_i(u) = \Lambda_j(u)$  and  $\gamma_i(A) = \sum_{B \in W(A)} \gamma_j(B)$ . The sub-solution's cost  $val(\gamma_i, \Lambda_i)$  is the minimum of  $val(\gamma_j, \Lambda_j)$  over all these  $(\gamma_j, \Lambda_j)$ .

*Join.* Let  $i \in I$  be a join node, and  $j, j' \in I$  its two children. We have  $X_i = X_j = X_{j'}$ . We call a mapping  $\beta_i : P_{st}(X_i) \times P_{st}(X_i) \rightarrow \{0, 1, \dots, k\}$  a pair-count on  $X_i$ . We define  $\beta_i$  to be an active pair-count if  $G_i^+$  has a coloring  $f_i$  such that  $\beta_i(A_j, A_{j'}) = |\{c \in C : A_j = \mathcal{V}(X_j, f_j, c) \text{ and } A_{j'} = \mathcal{V}(X_{j'}, f_{j'}, c)\}|$  for each  $A_j, A_{j'} \in P_{st}(X_i)$  where  $f_j$  and  $f_{j'}$  are the restriction of  $f_i$  to  $E(G_j^+)$  and  $E(G_{j'}^+)$ , respectively. Let  $val(\beta_i, A_i)$  denote the minimum cost value of an assignment of activation values for  $G_i^+$  which satisfies the restrictions  $A_i$  and activates an edge-colored subgraph of  $G_i^+$  that satisfies the pair-count  $\beta_i$ . The algorithm computes all  $val(\beta_i, A_i)$  of all  $\beta_i$  on  $X_i$  from all pairs of sub-solutions  $(\gamma_j, A_j)$  and  $(\gamma_{j'}, A_{j'})$  such that both have the same activation-value function ( $A_j(u) = A_{j'}(u)$  for all  $u \in X_i$ ) and the pair of active counts  $\gamma_j$  and  $\gamma_{j'}$  satisfy the following conditions:

- C1.  $\gamma_j(A_j) = \sum_{A \in P_{st}(X_i)} \beta_i(A_j, A)$  for all  $A_j \in P_{st}(X_i)$ .
- C2.  $\gamma_{j'}(A_{j'}) = \sum_{A \in P_{st}(X_i)} \beta_i(A, A_{j'})$  for all  $A_{j'} \in P_{st}(X_i)$ .

The value  $val(\beta_i, A_i)$  is set to be the summation value of the pair of sub-solutions  $(\gamma_j, A_j)$  and  $(\gamma_{j'}, A_{j'})$  that satisfy the above conditions minus the activation cost of  $X_i$ . To determine the pair  $(\gamma_i, A_i)$  that corresponds to the pair  $(\beta_i, A_i)$  of pair-count and an activation-value function, we construct a bipartite graph for each pair of partitions with a pair-count greater than 0 as follows. For each pair  $A_j, A_{j'} \in P_{st}(X_i)$  where  $\beta_i(A_j, A_{j'}) \geq 1$ , we construct a bipartite graph  $H_{\beta_i}^{(A_j, A_{j'})} = (A_j \cup A_{j'}, E_{\beta_i}^{(A_j, A_{j'})})$  with partite sets  $A_j$  and  $A_{j'}$ , where the vertices  $a_j \in A_j$  and  $a_{j'} \in A_{j'}$  are joined by an edge in  $E_{\beta_i}^{(A_j, A_{j'})}$  iff  $a_j \cap a_{j'} \neq \emptyset$ . Assume that  $D_1, D_2, \dots, D_b$  are the connected components of the graph  $H_{\beta_i}^{(A_j, A_{j'})}$ . We define  $\mathcal{U}(A_j, A_{j'})$  to be the family of vertex sets  $\{\bigcup_{v \in D_l} v : 1 \leq l \leq b\}$ . We set the value  $val(\gamma_i, A_i)$  to be the minimum  $val(\beta_i, A_i)$  over all  $(\beta_i, A_i)$  such that for each  $A_i \in P_{st}(X_i)$ :

$$\gamma_i(A_i) = \sum \beta_i(A_j, A_{j'})$$

where the summation above is taken over all pairs  $A_j, A_{j'} \in P_{st}(X_i)$  such that  $A_i = \mathcal{U}(A_j, A_{j'})$ .

*Extracting the Solution at the Root.* The algorithm checks all the pairs  $(\gamma_r, A_r)$  of the root bag  $X_r$  such that for all  $A \in P_{st}(X_r)$  where  $\gamma_r(A) \geq 1$ , there is a set in  $A$  containing both  $s$  and  $t$ . In this case  $(\gamma_r, A_r)$  corresponds to a feasible solution. The output of the algorithm is the minimum cost value among all the feasible solutions obtained at the root. For each row  $(\gamma_i, A_i)$  of bag  $X_i$ , we store the rows of  $X_i$ 's children that were used in the calculation of  $val(\gamma_i, A_i)$ . Computing the optimum solution is possible by traversing top-down in the decomposition tree to the leaves (traceback) to get the activation values, and then running a maximum flow algorithm on the unit-capacity graph of the activated edges to get the  $k$  edge-disjoint  $st$ -paths.

### 3.2 Analysis

The following lemmas analyse the running time of the algorithm and show that we can efficiently compute an optimal solution for the *st*-MAEDP problem for graphs with bounded treewidth. Let an instance of the problem be given by an activation network  $G = (V, E)$  with treewidth bounded by  $tw$  and terminals  $s, t \in V$ . Let  $\mathcal{P}^{OPT}$  represent an optimal solution for this instance. We use  $\mathcal{C}(\mathcal{P}^i)$  to denote the activation cost of a sub-solution  $\mathcal{P}^i$  in a partial graph  $G_i^+$ .

**Lemma 1.** *The *st*-MAEDP algorithm requires  $O(|V||D|^{tw+1}(k+1)^{(tw+3)^{2(tw+3)}}(tw+3)^{2(tw+3)+3}tw^3)$  time.*

*Proof.* The running time of the algorithm depends on the size of the tables and the combination of tables during the bottom-up traversal. For each set of vertices  $X$ , there are at most  $|X|^{|X|}$  possible partitions. Therefore, for each node  $i$ ,  $|P_{st}(X_i)| \leq (|X_i| + 2)^{(|X_i|+2)} \leq (tw+3)^{(tw+3)}$ . That means there are at most  $(k+1)^{(tw+3)^{(tw+3)}}$  possible active counts  $\gamma$  on  $X_i$ . The table  $tab_i$  in a processed bag  $X_i$  contains no more than  $(k+1)^{(tw+3)^{(tw+3)}}|D|^{tw+1}$  rows corresponding to the possible active counts  $\gamma$  and the  $|D|$  possible activation values for each vertex of  $X_i$ . Consider all possible row combinations with equal activation functions for two tables for a join node. Since there are at most  $(tw+3)^{2(tw+3)}$  possible pairs of partitions, there are at most  $(k+1)^{(tw+3)^{2(tw+3)}}$  possible active pair-counts on  $X_i$ . For each pair-count  $\beta_i$ , there is a pair of active counts  $\gamma_j$  and  $\gamma_{j'}$  such that the conditions  $C_1$  and  $C_2$  are satisfied. Computing  $\gamma_j$  and  $\gamma_{j'}$  from  $\beta_i$  takes  $O((tw+3)^{2(tw+3)})$  time. Therefore, the computation of the value  $val(\beta, \Lambda)$  for each combination of pair-count  $\beta$  and activation function  $\Lambda$  needs  $O((tw+3)^{2(tw+3)})$  time. Thus we see that all pairs  $(\beta, \Lambda)$  and  $val(\beta, \Lambda)$  can be computed in  $O((k+1)^{(tw+3)^{2(tw+3)}}(tw+3)^{2(tw+3)}|D|^{tw+1})$  time. Since  $|A| \leq tw+3$  for all  $A \in P_{st}(X_i)$ , the bipartite graph  $H_{\beta_i}^{(A_j, A_{j'})}$  defined in the join node contains no more than  $(tw+3)^2$  edges and can be constructed in  $O((tw+3)^3)$  time. Therefore, for each  $\beta_i$ , one can compute the active count  $\gamma_i$  that satisfies  $\gamma_i(A_i) = \sum \beta_i(A_j, A_{j'})$  where the summation is taken over all  $A_j, A_{j'} \in P_{st}(X_i)$  such that  $A_i = \mathcal{U}(A_j, A_{j'})$  and update the value  $val(\gamma_i, A_i)$  to be equal to  $val(\beta_i, A_i)$  if  $val(\beta_i, A_i) \leq val(\gamma_i, A_i)$  in  $O((tw+3)^{2(tw+3)}(tw+3)^3)$  time. Thus, we can compute all  $val(\gamma, \Lambda)$  of all combinations of active count  $\gamma$  and activation-value function  $\Lambda$  on  $X_i$  in time  $O((k+1)^{(tw+3)^{2(tw+3)}}(tw+3)^{2(tw+3)+3}|D|^{tw+1})$ . Since the tree-decomposition  $T$  has  $O(|V|tw^3)$  nodes, one can compute all the tables for all nodes in  $O(|V||D|^{tw+1}tw^3(k+1)^{(tw+3)^{2(tw+3)}}(tw+3)^{2(tw+3)+3})$  time.  $\square$

**Lemma 2.** *For any processed bag  $X_i$ , let  $\mathcal{P}_i^{OPT}$  be the induced solution of  $\mathcal{P}^{OPT}$  in  $G_i^+$  and  $(\gamma_i^{OPT}, \Lambda_i^{OPT})$  be the corresponding count and activation values, then*

$$val(\gamma_i^{OPT}, \Lambda_i^{OPT}) \leq \mathcal{C}(\mathcal{P}_i^{OPT}) \quad (1)$$

*Proof.* We use induction over the tree decomposition to prove that the value of  $(\gamma_i^{OPT}, \Lambda_i^{OPT})$  is at most the activation cost of  $\mathcal{P}_i^{OPT}$ . The base case are the



leaf nodes where the hypothesis clearly holds. Let us assume that the induction hypothesis holds for all descendants of bag  $X_i$ . We want to prove that the induction hypothesis holds also for  $X_i$  and that can be proven by showing that the hypothesis holds for all different types of the bag  $X_i$ .

*Introduce.* Let us assume that  $i$  is an introduce node and  $j$  its only child and let  $v$  be the additional vertex in  $X_i$ . The statement (1) holds for  $X_j$  because it is a child of the bag  $X_i$ . Let  $(\gamma_j^{OPT}, \Lambda_j^{OPT})$  be the corresponding active count and activation values of  $\mathcal{P}_j^{OPT}$ . The following statement holds:

$$val(\gamma_j^{OPT}, \Lambda_j^{OPT}) \leq \mathcal{C}(\mathcal{P}_j^{OPT})$$

$(\gamma_i^{OPT}, \Lambda_i^{OPT})$  and  $(\gamma_j^{OPT}, \Lambda_j^{OPT})$  both agree on the activation values for all vertices in  $X_i \setminus \{v\}$ . The additional vertex  $v$  is an isolated vertex in the induced graph  $G_i^+$  and the one vertex set  $\{v\}$  is in  $A$  for all  $A \in P_{st}(X_i)$  such that  $\gamma_i^{OPT}(A) \geq 1$ . Since  $\Lambda_i^{OPT}(v) \in D$  is the activation value of the vertex  $v$ , then:

$$\begin{aligned} val(\gamma_i^{OPT}, \Lambda_i^{OPT}) &= val(\gamma_j^{OPT}, \Lambda_j^{OPT}) + \Lambda_i^{OPT}(v) \\ &\leq \mathcal{C}(\mathcal{P}_j^{OPT}) + \Lambda_i^{OPT}(v) \\ &= \mathcal{C}(\mathcal{P}_i^{OPT}) \end{aligned}$$

The statement (1) holds for an introduce node.

*Forget.* Assume that  $i$  is a forget node and  $j$  its only child and let  $v$  be the discarded vertex in  $X_i$ . The statement (1) holds for  $X_j$  because it is a child of the bag  $X_i$ . Let  $(\gamma_j^{OPT}, \Lambda_j^{OPT})$  be the corresponding active count and activation values of  $\mathcal{P}_j^{OPT}$ . The following statement holds:

$$val(\gamma_j^{OPT}, \Lambda_j^{OPT}) \leq \mathcal{C}(\mathcal{P}_j^{OPT})$$

$(\gamma_i^{OPT}, \Lambda_i^{OPT})$  and  $(\gamma_j^{OPT}, \Lambda_j^{OPT})$  both agree on activation values for all vertices in  $X_j \setminus \{v\}$ . The discarded vertex  $v$  is either an isolated vertex or part of a connected component in the induced solution  $\mathcal{P}_j^{OPT}$  and that means the value  $val(\gamma_j^{OPT}, \Lambda_j^{OPT})$  is one of the values that has been considered when calculating  $val(\gamma_i^{OPT}, \Lambda_i^{OPT})$ . Then:

$$val(\gamma_i^{OPT}, \Lambda_i^{OPT}) \leq val(\gamma_j^{OPT}, \Lambda_j^{OPT}) \leq \mathcal{C}(\mathcal{P}_j^{OPT}) = \mathcal{C}(\mathcal{P}_i^{OPT})$$

Then the statement (1) holds for a forget node.

*Join.* Assume that  $i$  is a join node and  $j$  and  $j'$  its children. Let  $(\gamma_i^{OPT}, \Lambda_i^{OPT})$  be the corresponding active count and activation values of the induced solution  $\mathcal{P}_i^{OPT}$  in  $G_i^+$ .  $\mathcal{P}_i^{OPT}$  is the union of its children sub-solutions  $\mathcal{P}_j^{OPT}$  and  $\mathcal{P}_{j'}^{OPT}$ . Let  $(\gamma_j^{OPT}, \Lambda_j^{OPT})$  and  $(\gamma_{j'}^{OPT}, \Lambda_{j'}^{OPT})$  be the corresponding active count and activation values of  $\mathcal{P}_j^{OPT}$  and  $\mathcal{P}_{j'}^{OPT}$ , respectively. The statement (1) holds for  $X_j$  and  $X_{j'}$  because they are children of the bag  $X_i$ . For all  $l \in \{j, j'\}$  the following statement holds:

$$val(\gamma_l^{OPT}, \Lambda_l^{OPT}) \leq \mathcal{C}(\mathcal{P}_l^{OPT})$$

Let  $(\beta_i^{OPT}, \Lambda_i^{OPT})$  be the corresponding active pair-count and activation values of the induced solution  $\mathcal{P}_i^{OPT}$  in  $G_i^+$  which satisfies conditions  $C_1$  and  $C_2$  for  $(\gamma_j^{OPT}, \Lambda_j^{OPT})$  and  $(\gamma_{j'}^{OPT}, \Lambda_{j'}^{OPT})$ . We know that  $\mathcal{P}_i^{OPT}$  is a sub-solution in  $G_i^+$ , therefore,  $val(\beta_i^{OPT}, \Lambda_i^{OPT})$  is one of the values that has been considered when computing  $val(\gamma_i^{OPT}, \Lambda_i^{OPT})$ . Then:

$$\begin{aligned} val(\gamma_i^{OPT}, \Lambda_i^{OPT}) &\leq val(\beta_i^{OPT}, \Lambda_i^{OPT}) \\ &= val(\gamma_j^{OPT}, \Lambda_j^{OPT}) + val(\gamma_{j'}^{OPT}, \Lambda_{j'}^{OPT}) - \sum_{v \in X_i} \Lambda_i^{OPT}(v) \\ &\leq \mathcal{C}(\mathcal{P}_j^{OPT}) + \mathcal{C}(\mathcal{P}_{j'}^{OPT}) - \sum_{v \in X_i} \Lambda_i^{OPT}(v) \\ &= \mathcal{C}(\mathcal{P}_i^{OPT}) \end{aligned}$$

Then the induction hypothesis holds for a join node.  $\square$

**Lemma 3.** *For any processed bag  $X_i$ , any pair  $(\gamma_i, \Lambda_i)$  where  $val(\gamma_i, \Lambda_i) = c_i < \infty$  corresponds to an edge-coloring plus activation-values  $\mathcal{P}^i = (f_i, \Lambda_i^+)$  where  $f_i : E(G_i^+) \rightarrow \{0, \dots, k\}$  and  $\Lambda_i^+ : X_i^+ \rightarrow D$  with the following properties:*

- The active count of  $\mathcal{P}^i$  in  $X_i$  is  $\gamma_i$ .
- The activation values of  $\mathcal{P}^i$  in  $X_i$  are  $\Lambda_i$ .
- The total activation cost in  $X_i^+$  is  $c_i$ .

*Proof.* We prove by induction that for any bag  $X_i$  there exists an edge-coloring and activation-values  $\mathcal{P}^i$  with the above properties. The base case are the leaf nodes where the hypothesis clearly holds. Let us assume that the induction hypothesis holds for all the descendants of bag  $X_i$ . We want to prove that the induction hypothesis holds also for  $X_i$  and that can be proved by showing that the hypothesis holds for all different types of the bag  $X_i$ .

*Introduce.* Assume that  $i$  is an introduce node and  $j$  its only child and  $v$  the additional vertex in  $X_i$ . Let  $(\gamma_i, \Lambda_i)$  be some entry with  $val(\gamma_i, \Lambda_i) = c_i < \infty$ . The induction hypothesis holds for  $X_j$ . Let  $(\gamma_j, \Lambda_j)$  be the corresponding active count and activation function that have been used for the calculation of  $val(\gamma_i, \Lambda_i)$ . Then  $(\gamma_j, \Lambda_j)$  corresponds to an edge-coloring and activation-values  $\mathcal{P}^j$  that satisfies the above properties. From the algorithm,  $(\gamma_i, \Lambda_i)$  and  $(\gamma_j, \Lambda_j)$  both agree on the activation values for all vertices in  $X_i \setminus \{v\}$ . Moreover, the additional vertex  $v$  is an isolated vertex in the induced graph  $G_i^+$  and the one vertex set  $\{v\}$  is in  $A$  for all  $A \in P_{st}(X_i)$  such that  $\gamma_i(A) \geq 1$ . The value of  $(\gamma_i, \Lambda_i)$  is equal to the value of  $(\gamma_j, \Lambda_j)$  added to the activation value of the vertex  $v$ . The union of the isolated vertex  $v$  and  $\mathcal{P}^j$  is an edge-coloring with activation-values that satisfies all properties of the induction hypothesis. Thus, the induction hypothesis holds for an introduce node.

*Forget.* Assume that  $i$  is a forget node and  $j$  its only child and  $v$  the discarded vertex. Let  $(\gamma_i, \Lambda_i)$  be some entry with  $val(\gamma_i, \Lambda_i) = c_i < \infty$ . The induction hypothesis holds for  $X_j$ . Let  $(\gamma_j, \Lambda_j)$  be the corresponding active count and

activation function that have been used for the calculation of  $val(\gamma_i, A_i)$ . Then  $(\gamma_j, A_j)$  corresponds to an edge-coloring and activation-values  $\mathcal{P}^j$  that satisfies the above properties. From the algorithm,  $(\gamma_j, A_j)$  has the minimum value cost among all possible rows of  $X_j$  that can produce  $(\gamma_i, A_i)$  and both agree on activation values for all vertices in  $X_j \setminus \{v\}$ . The discarded vertex  $v$  is part of the edge-coloring  $\mathcal{P}^j$ . Therefore  $(\gamma_i, A_i)$  corresponds to the edge-coloring and activation-values  $\mathcal{P}^j$  and we can set  $\mathcal{P}^i = \mathcal{P}^j$ . The induction hypothesis holds for a forger node.

*Join.* Assume that  $i$  is a join node and  $j$  and  $j'$  its children. Let  $(\gamma_i, A_i)$  be some entry with  $val(\gamma_i, A_i) = c_i < \infty$ . Since  $val(\gamma_i, A_i) = c_i < \infty$ , then there is a pair of active pair-count and activation function  $(\beta_i, A_i)$  such that  $val(\gamma_i, A_i) = val(\beta_i, A_i)$  and for each  $A_i \in P_{st}(X_i)$ ,  $\gamma_i(A_i) = \sum \beta_i(A_j, A_{j'})$  where the summation is taken over all pairs  $A_j, A_{j'} \in P_{st}(X_i)$  satisfying  $A_i = \mathcal{U}(A_j, A_{j'})$ . Let  $(\gamma_j, A_j)$  and  $(\gamma_{j'}, A_{j'})$  be the pairs that have been used for the calculation of  $val(\beta_i, A_i)$  which satisfy the conditions  $C_1$  and  $C_2$ . We know that  $val(\gamma_j, A_j) < \infty$  and  $val(\gamma_{j'}, A_{j'}) < \infty$  and the induction hypothesis holds for  $X_j$  and  $X_{j'}$ . Therefore,  $(\gamma_j, A_j)$  and  $(\gamma_{j'}, A_{j'})$  correspond to edge-colorings with activation-values  $\mathcal{P}^j = (f_j, A_j^+)$  and  $\mathcal{P}^{j'} = (f_{j'}, A_{j'}^+)$ , respectively. For each pair of partitions  $A_j$  and  $A_{j'}$  where  $\gamma_j(A_j) = r_j$ ,  $\gamma_{j'}(A_{j'}) = r_{j'}$  and  $\beta_i(A_j, A_{j'}) = r > 0$ , there are  $r_l$  colors for partition  $A_l$  in  $\mathcal{P}^l = (f_l, A_l)$  for  $l \in \{j, j'\}$ . Therefore, we choose  $r$  unused colors  $z_1^j, z_2^j, \dots, z_r^j$  for  $A_j$  in  $\mathcal{P}^j$  and  $r$  unused colors  $z_1^{j'}, z_2^{j'}, \dots, z_r^{j'}$  for  $A_{j'}$  in  $\mathcal{P}^{j'}$  and then recolor both  $z_w^j$  in  $G_j^+(f_j, z_w^j)$  and  $z_w^{j'}$  in  $G_{j'}^+(f_{j'}, z_w^{j'})$  with a new color  $z_w^{jj'}$  for all  $w \in \{1, \dots, r\}$ . The colors  $z_1^j, z_2^j, \dots, z_r^j$  in  $\mathcal{P}^j$  and colors  $z_1^{j'}, z_2^{j'}, \dots, z_r^{j'}$  in  $\mathcal{P}^{j'}$  are now marked as used. Note that there are always enough unused colors because the conditions  $C_1$  and  $C_2$  are satisfied. After recoloring  $\mathcal{P}^j$  and  $\mathcal{P}^{j'}$  and since for each  $A_i \in P_{st}(X_i)$ ,  $\gamma_i(A_i) = \sum \beta_i(A_j, A_{j'})$  where the summation is taken over all pairs  $A_j, A_{j'} \in P_{st}(X_i)$  satisfying  $A_i = \mathcal{U}(A_j, A_{j'})$ , it follows that the active count of the union  $\mathcal{P}^j \cup \mathcal{P}^{j'}$  in  $X_i$  is  $\gamma_i$ . Moreover,  $(\gamma_i, A_i)$ ,  $(\gamma_j, A_j)$  and  $(\gamma_{j'}, A_{j'})$  all have the same activation-value function ( $A_i(u) = A_j(u) = A_{j'}(u)$  for all  $u \in X_i$ ). That means the activation values of the union  $\mathcal{P}^j \cup \mathcal{P}^{j'}$  in  $X_i$  are  $A_i$ . The algorithm also computes the cost value of  $(\gamma_i, A_i)$  as follows:

$$val(\gamma_i, A_i) = val(\beta_i, A_i) = val(\gamma_j, A_j) + val(\gamma_{j'}, A_{j'}) - \sum_{v \in X_i} A_i(v)$$

Since  $val(\gamma_j, A_j)$  and  $val(\gamma_{j'}, A_{j'})$  are the total activation costs in  $X_j^+$  and  $X_{j'}^+$ , respectively, then the summation of these activation costs minus the activation values for all  $u \in X_i$  is the total activation cost in  $X_i^+$ . We can set  $\mathcal{P}^i$  to be the edge-coloring and activation values  $\mathcal{P}^j \cup \mathcal{P}^{j'}$  that satisfies the properties of the lemma. The induction hypothesis holds for a join node.  $\square$

We obtain the following theorem by combining the above lemmas.

**Theorem 2.** *The st-MAEDP problem for graphs with treewidth bounded by  $tw$  can be solved optimally in  $O(|V||D|^{tw+1}tw^3(k+1)^{(tw+3)2^{(tw+3)}}(tw+3)^{2(tw+3)+3})$  time.*

**Corollary 1.** *For any fixed  $k$ , the  $st$ -MAEDP problem for graphs of bounded treewidth can be solved optimally in linear-time FPT parameterized by the treewidth  $tw$ .*

## 4 Conclusion

To the best of our knowledge, the  $st$ -MAEDP problem for graphs with treewidth bounded by  $tw$  considered here has not been addressed before. We established an algorithm that solves the problem optimally in  $O(|V||D|^{tw+1}tw^3(k+1)^{(tw+3)^{2(tw+3)}}(tw+3)^{2(tw+3)+3})$  time. Our algorithm also solves the  $st$ -MAEDP problem when  $k = 2$  in linear-time and this is an improvement over the cubic algorithm obtained in [2]. It would be interesting if one can obtain a faster or even linear-time algorithm for the  $st$ -MAEDP problem in graphs with treewidth bounded by a constant.

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Combinatorial Algorithms

26th International Workshop, IWOCA 2015, Verona,  
Italy, October 5-7, 2015, Revised Selected Papers

Lipták, Z.; Smyth, W.F. (Eds.)

2016, XIX, 360 p. 62 illus. in color., Softcover

ISBN: 978-3-319-29515-2