

# Chromatic Correlation Clustering

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## ABSTRACT

We study a novel clustering problem in which the pairwise relations between objects are *categorical* rather than *numerical*, as usual. This problem can be naturally viewed as clustering the vertices of a graph whose edges are of different types (*colors*). We tackle this problem by introducing an objective function which aims to partition the graph such that the edges within each cluster have, as much as possible, the same color. We show that the problem is NP-hard and propose a randomized algorithm with approximation guarantee proportional to the maximum degree of the input graph. The algorithm iteratively picks a random edge as pivot, builds a cluster around it, and removes the cluster from the graph. Although being fast, easy-to-implement, and parameter free, this algorithm tends to produce a relatively large number of clusters. To overcome this issue, we introduce a variant based on two heuristics concerning the way the pivot is chosen and the way the cluster is built around it. Finally, to address the case where a fixed number of output clusters is required, we devise a third algorithm that directly optimizes the objective function via a strategy based on the *alternating minimization* paradigm.

We test our algorithms both on synthetic as well as real data coming from domains such as protein-protein interactions, social media, and bibliometrics. Experimental evidence show that our algorithms outperform a baseline algorithm both in the task of reconstructing a ground-truth clustering and in terms of objective function value.

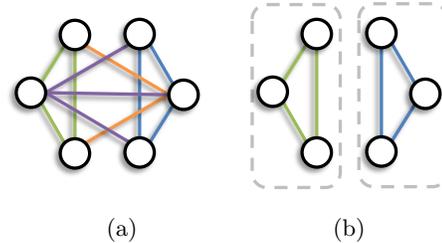
## 1. INTRODUCTION

Clustering is one of the most well-studied problems in data mining. The goal of clustering is to partition a set of objects in different clusters, so that objects in the same cluster are more similar to each other than to objects in other clusters. A common trait underlying most clustering paradigms is the existence of a function  $\text{sim}(x, y)$  representing the similarity between pairs of objects  $x$  and  $y$ . The similarity function is either provided explicitly as input, or it can be computed implicitly from the representation of the objects.

In this paper, we consider a different clustering setting

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**Figure 1: An example of chromatic clustering: (a) input graph, (b) the optimal solution for chromatic-correlation-clustering (Problem 2).**

where the relationship among objects is represented by a relation type, such as a label  $\ell(x, y)$  from a finite set of possible labels  $L$ . In other words, the *range* of the similarity function  $\text{sim}(x, y)$  can be viewed as being *categorical*, instead of *numerical*. Moreover, we model the case where two objects  $x$  and  $y$  do not have any relation at all with a special label  $l_0 \notin L$ .

Our framework has a natural graph interpretation: the input to our problem can be viewed as an edge-labeled graph  $G = (V, E, L, \ell)$ , where the set of vertices  $V$  is the set of objects to be clustered, the set of edges  $E \subseteq V \times V$  is implicitly defined as  $E = \{(x, y) \in V \times V \mid \ell(x, y) \neq l_0\}$ , and each edge has a *label* in  $L$  or, as we like to think about it, a *color*.

The key objective in our framework is to find a partition of the vertices of the graph such that the edges in each cluster have, as much as possible, the same color. An example of clustering objects with chromatic relations is shown in Figure 1. Intuitively, a **red** edge  $(x, y)$  provides positive evidence that the vertices  $x$  and  $y$  should be clustered in such a way that the edges in the subgraph induced by that cluster are mostly **red**. Furthermore, in the case that most edges of a cluster are **red**, it is reasonable to label the whole cluster with the **red** color. Note that a clustering algorithm for this problem should also deal with inconsistent evidence, as a **red** edge  $(x, y)$  provides evidence for the vertex  $x$  to participate in a cluster with **red** edges, while a **green** edge  $(x, z)$  provides contradicting evidence for the vertex  $x$  to participate in a cluster with **green** edges. Aggregating such inconsistent information is resolved by optimizing a properly-defined objective function.

**Applications.** The study of edge-labeled graphs is motivated by many real-world applications and is receiving increasing attention in the data-mining literature [8, 10, 16]. As an example, biologists study protein-protein interaction networks, where vertices represent proteins and edges represent interactions occurring when two or more proteins bind

together to carry out their biological function. Those interactions can be of different types, e.g., physical association, direct interaction, co-localization, etc. In these networks, for instance, a cluster containing mainly edges labeled as co-localization, might represent a *protein complex*, i.e., a group of proteins that interact with each other at the same time and place, forming a single multi-molecular machine [11].

As a further example, social networks are commonly represented as graphs, where the vertices represent individuals and the edges capture relationships among these individuals. Again, these relationships can be of various types, e.g., colleagues, neighbors, schoolmates, football-mates.

In bibliographic data, co-authorship networks represent collaborations among authors: in this case the topic of the collaboration can be seen as an edge label, and a cluster of vertices represents a topic-coherent community of researchers. In our experiments in Section 5 we show how our framework can be applied in all the above domains.

**Contributions.** In this paper we address the problem of clustering data with categorical similarity, achieving the following contributions:

- We define CHROMATIC-CORRELATION-CLUSTERING, a novel clustering problem for objects with categorical similarity, by revisiting the well-studied *correlation clustering* framework [3]. We show that our problem is a generalization of the traditional CORRELATION-CLUSTERING problem, implying that it is **NP**-hard.
- We introduce a randomized algorithm, named **Chromatic Balls**, that provides approximation guarantee proportional to the maximum degree of the graph.
- Though of theoretical interest, **Chromatic Balls** has some limits when it comes to practice. Trying to overcome these limits, we introduce two alternative algorithms: a more practical *lazy* version of **Chromatic Balls**, and an algorithm that directly optimizes the proposed objective function via an iterative process based on the *alternating minimization* paradigm.
- We empirically assess our algorithms both on synthetic and real datasets. Experiments on synthetic data show that our algorithms outperform a baseline algorithm in the task of reconstructing a ground-truth clustering. Experiments on real-world data confirm that CHROMATIC-CORRELATION-CLUSTERING provides meaningful clusters in terms of objective clustering value.

The rest of the paper is organized as follows. In the next section we recall the traditional correlation clustering problem and introduce our new formulation. In Section 3 we introduce the **Chromatic Balls** algorithm and we prove its approximation guarantees. In Section 4 we present the two more practical algorithms, namely **Lazy Chromatic Balls** and **Alternating Minimization**. In Section 5 we report our experimental analysis. In Section 6 we discuss related work, while Section 7 concludes the paper.

## 2. PROBLEM DEFINITION

Given a set of objects  $V$ , a clustering problem asks to partition the set  $V$  into clusters of similar objects. Assuming that cluster identifiers are represented by natural numbers, a clustering  $\mathcal{C}$  can be seen as a function  $\mathcal{C} : V \rightarrow \mathbb{N}$ . Typically, the goal is to find a clustering  $\mathcal{C}$  that optimizes an objective function that measures the quality of the clustering

Numerous formulations and objective functions have been considered in the literature. One of these, considered both in the area of theoretical computer science and data mining, is that at the basis of the CORRELATION-CLUSTERING problem [3].

### Problem 1 (CORRELATION-CLUSTERING)

Given a set of objects  $V$  and a pairwise similarity function  $\text{sim} : V \times V \rightarrow [0, 1]$ , find a clustering  $\mathcal{C} : V \rightarrow \mathbb{N}$  that minimizes the cost

$$\text{cost}(\mathcal{C}) = \sum_{\substack{(x,y) \in V \times V \\ \mathcal{C}(x) = \mathcal{C}(y)}} (1 - \text{sim}(x,y)) + \sum_{\substack{(x,y) \in V \times V \\ \mathcal{C}(x) \neq \mathcal{C}(y)}} \text{sim}(x,y). \quad (1)$$

The intuition underlying the above problem is that the cost of assigning two objects  $x$  and  $y$  to the same cluster should be equal to the dissimilarity  $1 - \text{sim}(x, y)$ , while the cost of assigning the objects in different clusters should correspond to their similarity  $\text{sim}(x, y)$ . A common case is when the similarity is binary, that is,  $\text{sim} : V \times V \rightarrow \{0, 1\}$ . In this case, Equation (1) reduces to counting the number of pairs of objects that have similarity 0 and are put in the same cluster plus the number of pairs of objects that have similarity 1 and belong to different clusters. Or equivalently, in a graph-based terminology, the objective function counts the number of “positive” edges that are cut plus the number of “negative” (i.e., non-existing) edges that are not cut.

In CHROMATIC-CORRELATION-CLUSTERING, which we formally define below, we still have negative edges (i.e.,  $l_0$ -edges), but the positive edges may have different colors, representing different kinds of relations among the objects.

### Problem 2 (CHROMATIC-CORRELATION-CLUSTERING)

Given a set  $V$  of objects, a set  $L$  of labels, a special label  $l_0$ , and a pairwise labeling function  $\ell : V \times V \rightarrow L \cup \{l_0\}$ , find a clustering  $\mathcal{C} : V \rightarrow \mathbb{N}$  and a cluster labeling function  $\mathcal{c} \ell : \mathcal{C}[V] \rightarrow L$  so to minimize the cost

$$\text{cost}(\mathcal{C}, \mathcal{c} \ell) = \sum_{\substack{(x,y) \in V \times V, \\ \mathcal{C}(x) = \mathcal{C}(y)}} (1 - I[\ell(x,y) = \mathcal{c} \ell(\mathcal{C}(x))]) + \sum_{\substack{(x,y) \in V \times V, \\ \mathcal{C}(x) \neq \mathcal{C}(y)}} I[\ell(x,y) \neq l_0]. \quad (2)$$

Equation (2) is composed by two terms, representing intra- and inter-cluster costs, respectively. In particular, according to the intra-cluster cost term, any pair of objects  $(x, y)$  assigned to the same cluster should pay a cost if and only if their relation type  $\ell(x, y)$  is other than the predominant relation type of the cluster indicated by the function  $\mathcal{c} \ell$ . For the inter-cluster cost, the objective function does not penalize a pair of objects  $(x, y)$  only if they do not have any relation, i.e.,  $\ell(x, y) = l_0$ . If  $\ell(x, y) \neq l_0$ , the objective function incurs a cost, regardless of the label  $\ell(x, y)$ .

**Example 1** For the problem instance in Figure 1(a), the solution in Figure 1(b) has a cost of 5: there is no intra-cluster cost, because the two clusters are cliques and their edges are monochromatic, while we have an inter-cluster cost of 5 as equal to the number of edges that are cut.

It is trivial to observe that, when  $|L| = 1$ , the CHROMATIC-CORRELATION-CLUSTERING problem corresponds to the binary version of CORRELATION-CLUSTERING. Thus, our problem is a generalization of the standard problem. Since

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**Algorithm 1** Chromatic Balls

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**Input:** Edge-labeled graph  $G = (V, E, L, \ell)$ **Output:** Clustering  $\mathcal{C} : V \rightarrow \mathbb{N}$ ; cluster labeling function  $cl : \mathcal{C}[V] \rightarrow L$ 

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1:  $V' \leftarrow V$ ;  $i \leftarrow 1$ 
2: while there exist  $u, v \in V'$  such that  $(u, v) \in E$  do
3:   randomly pick  $u, v \in V'$  such that  $(u, v) \in E$ 
4:    $C \leftarrow \{u, v\} \cup \{x \in V' \mid \ell(u, x) = \ell(v, x) = \ell(u, v)\}$ 
5:   for all  $x \in C$  do  $\mathcal{C}(x) \leftarrow i$ 
6:    $cl(i) = \ell(u, v)$ 
7:    $V' \leftarrow V' \setminus C$ ;  $i \leftarrow i + 1$ 
8: for all  $x \in V'$  do
9:    $\mathcal{C}(x) \leftarrow i$ 
10:   $cl(i) \leftarrow$  a random label from  $L$ 
11:   $i \leftarrow i + 1$ 
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CORRELATION-CLUSTERING is **NP**-hard, we can easily conclude that CHROMATIC-CORRELATION-CLUSTERING is **NP**-hard too.

The previous observation motivates us to consider whether applying standard CORRELATION-CLUSTERING algorithms, just ignoring the different colors, is a good solution to the problem. As we show in the following example, such an approach does not guarantee to produce good solutions.

**Example 2** For the problem instance in Figure 1(a), the optimal solution for the standard CORRELATION-CLUSTERING which does not consider the different colors, would be composed by a single cluster containing all the six vertices, as, according to Equation (1), this solution has a (minimum) cost of 4 corresponding to the number of missing edges within the cluster. Conversely, this solution has a non-optimal cost 12 when evaluated according to the CHROMATIC-CORRELATION-CLUSTERING formulation, i.e., according to Equation (2). Instead, the optimum in this case would correspond to the cost 5 solution depicted in Figure 1(b).

Although the example shows that for the chromatic version of the problem we cannot directly apply algorithms developed for the CORRELATION-CLUSTERING problem, we can use such algorithms at least as a starting point, as shown in the next section.

### 3. THE Chromatic Balls ALGORITHM

We present next a randomized approximation algorithm for the CHROMATIC-CORRELATION-CLUSTERING problem. This algorithm, called **Chromatic Balls**, is motivated by the **Balls** algorithm [1], which is an approximation algorithm for standard CORRELATION-CLUSTERING.

For completeness, we briefly review the **Balls** algorithm. The algorithm works in iterations. Initially all objects are considered *uncovered*. In each iteration the algorithm produces a cluster, and the objects participating in the cluster are considered *covered*. In particular, the algorithm picks as pivot a random object currently uncovered, and forms a cluster consisting of the pivot itself along with all currently uncovered objects that are connected to the pivot.

The outline of our **Chromatic Balls** is summarized in Algorithm 1. The main difference with the **Balls** algorithm is that the edge labels are taken into account in order to build clusters around the pivots. To this end, the pivot chosen at each iteration of **Chromatic Balls** is an edge rather than a single object. The **Chromatic Balls** algorithm employs a set  $V'$  to keep all the objects that have not been assigned to any cluster yet; hence, initially,  $V' = V$ . At each iteration, a random edge  $(u, v)$  such that both objects  $u$  and  $v$  are

currently in the set  $V'$  is selected as pivot (line 3). Given the pivot  $(u, v)$ , a cluster  $C$  is formed around it. Beyond the objects  $u$  and  $v$ , the cluster  $C$  additionally contains all other objects  $x \in V'$  for which the *triangle*  $(u, v, x)$  is monochromatic, that is,  $\ell(u, x) = \ell(v, x) = \ell(u, v)$  (lines 4 and 5). Since the label  $\ell(u, v)$  forms the basis for creating the cluster  $C$ , the cluster is labeled with this label (line 6). All objects added in  $C$  are removed from  $V'$  (line 7), and the algorithm terminates when  $V'$  does not contain any pair of objects that share an edge, i.e., that is labeled with a label other than  $l_0$  (line 2). All objects remaining in the set  $V'$ , if any, are eventually made singleton clusters (lines 8-11).

**Computational complexity.** The complexity of the **Chromatic Balls** algorithm is determined by two steps: (i) picking the pivots (line 3), and (ii) building the clusters (line 4). Choosing the pivots requires  $\mathcal{O}(m \log n)$  time, where  $n = |V|$  and  $m = |E|$ , as selecting random edges can be implemented by building a priority queue of edges with random priorities, and subsequently removing edges; each edge is removed once from the priority queue, whether it is selected as pivot or not. Building a single cluster  $C$ , instead, requires to access all neighbors of the pivot edge  $(u, v)$ . As the current cluster is removed from the set of uncovered objects at the end of each iteration, the neighbors of any pivot are not considered again in the remainder of the algorithm. Thus, the step of selecting the objects to be included into the current clusters requires visiting each edge at most once; therefore, the process takes  $\mathcal{O}(m)$  time. In conclusion, we can state that the computational complexity of the **Chromatic Balls** algorithm is  $\mathcal{O}(m \log n)$ .

#### 3.1 Theoretical analysis

We analyze next the quality of the solutions obtained by **Chromatic Balls**. Our main result, given in Theorem 1, shows that the approximation guarantee of the algorithm depends on the number of *bad triplets* incident to a pair of objects in the input dataset. The notion of bad triplet is defined below; however, here we note that this result gives a constant-factor guarantee for bounded-degree graphs.

Even though the **Chromatic Balls** algorithm is similar to the **Balls** algorithm, which can be shown to provide a constant-factor approximation guarantee for general graphs too, the theoretical analysis of **Chromatic Balls** is much more complicated and requires several additional and nontrivial arguments. Due to the limited space of this paper, we report next only an outline of our analysis. Further details, including complete proofs, can be found in an extended technical report.<sup>1</sup>

We begin our analysis by defining special types of triplets and quadruples among the vertices of the graph.

**Definition 1 (SC-triplet)** We say that  $\{x, y, z\}$  is a same-color triplet (*SC-triplet*) if the induced triangle is monochromatic, i.e.,  $\ell(x, y) = \ell(x, z) = \ell(y, z) \neq l_0$ .

**Definition 2 (B-triplet)** We say that  $\{x, y, z\}$  is a bad triplet (*B-triplet*) if the induced triangle is non-monochromatic and it has at most one pair labeled with  $l_0$ .

**Definition 3 (B-quadruple)** A Bad-quadruple is a set  $\{x, y, z, w\} \subseteq V$  that contains at least one SC-triplet and at least one B-triplet.

Note that, according to the cost function of our problem as defined in Equation (2), there is no way to partition a B-triplet without paying any cost. Next we define the notions of *hitting* and *d-hitting*.

<sup>1</sup><http://francescobonchi.com/CCC.pdf>

**Definition 4 (hitting)** Consider a pair of objects  $(x, y)$  and a triplet  $t$ , which can be either SC-triplet or B-triplet. We say that  $t$  hits  $(x, y)$  if  $x \in t$  and  $y \in t$ . Additionally, if  $q$  is a B-quadruple, we say that  $q$  hits  $(x, y)$  if  $x \in q$ ,  $y \in q$ , and there exists  $z \in q$  such that  $\{x, y, z\}$  is a B-triplet.

**Definition 5 (d-hitting)** Given any pair of objects  $(x, y)$  and any B-quadruple  $q = \{x, y, z, w\}$ , we say that  $q$  deeply hits (d-hits)  $(x, y)$  if  $q$  hits  $(x, y)$  and either  $\{x, z, w\}$  or  $\{y, z, w\}$  is an SC-triplet.

In reference to the above notions, we hereinafter denote by  $\mathcal{S}$ ,  $\mathcal{T}$ , and  $\mathcal{Q}$  the sets of all SC-triplets, B-triplets, and B-quadruples for an instance of our problem. Moreover, given a pair  $(x, y) \in V \times V$  we define the following sets:  $\mathcal{T}_{xy} \subseteq \mathcal{T}$  denotes the set of all B-triplets in  $\mathcal{T}$  that hit  $(x, y)$ ;  $\mathcal{Q}_{xy} \subseteq \mathcal{Q}$  denotes the set of all B-quadruples in  $\mathcal{Q}$  that hit  $(x, y)$ ;  $\mathcal{Q}_{xy}^d \subseteq \mathcal{Q}_{xy} \subseteq \mathcal{Q}$  denotes the set of all B-quadruples in  $\mathcal{Q}$  that d-hit  $(x, y)$ .

Let us now consider some events that may arise during the execution of the Chromatic Balls algorithm. Given an object  $x \in V$ ,  $P_x^{(i)}$  denotes the event “ $x$  is chosen as pivot in the  $i$ -th iteration.” Given a set  $\{x_1, \dots, x_n\} \subseteq V$ , with  $n \geq 2$ ,  $A_{x_1 \dots x_n}^{(i)}$  denotes the event “all objects  $x_1, \dots, x_n$  enter the  $i$ -th iteration of the algorithm, while two of them are chosen as pivot in the same iteration.”

Additionally, the events  $T_{z|xy}^{(i)}$  and  $Q_{zw|xy}^{(i)}$  are defined in reference to a pair  $(x, y)$ . Given a B-triplet  $\{x, y, z\} \in \mathcal{T}_{xy}$ ,  $T_{z|xy}^{(i)}$  denotes the event “ $A_{xyz}^{(i)}$  occurs while  $x$  and  $y$  are not chosen both as pivots in the  $i$ -th iteration.” Given a B-quadruple  $\{x, y, z, w\} \in \mathcal{Q}_{xy}^d$ ,  $Q_{zw|xy}^{(i)}$  denotes the event “ $A_{xyzw}^{(i)}$  occurs while neither  $x$  nor  $y$  are chosen as pivots in  $i$ -th iteration.”

For the events  $A_{x_1 \dots x_n}^{(i)}$ ,  $T_{z|xy}^{(i)}$ , and  $Q_{zw|xy}^{(i)}$ , defined above, we also consider their counterparts that assert that the events occur at some iteration  $i$ . For instance,  $A_{x_1 \dots x_n}$  denotes the event “ $A_{x_1 \dots x_n}^{(i)}$  happens at some iteration  $i$ ,” while  $T_{z|xy}$  and  $Q_{zw|xy}$  are defined analogously. Formally:

$$A_{x_1 \dots x_n} \Leftrightarrow \bigvee_i A_{x_1 \dots x_n}^{(i)}, \quad (3)$$

$$T_{z|xy} \Leftrightarrow \bigvee_i T_{z|xy}^{(i)} \Leftrightarrow \bigvee_i \left( A_{xyz}^{(i)} \wedge \neg \left( P_x^{(i)} \wedge P_y^{(i)} \right) \right), \quad (4)$$

$$Q_{zw|xy} \Leftrightarrow \bigvee_i Q_{zw|xy}^{(i)} \Leftrightarrow \bigvee_i \left( A_{xyzw}^{(i)} \wedge \neg P_x^{(i)} \wedge \neg P_y^{(i)} \right). \quad (5)$$

As reported in the next two lemmas, the probabilities of the events  $T_{z|xy}$  and  $Q_{zw|xy}$  can be expressed in terms of the probabilities of the events  $A_{xyz}$  and  $A_{xyzw}$ .

**Lemma 1** Given a pair  $(x, y) \in V \times V$  and a B-triplet  $\{x, y, z\} \in \mathcal{T}_{xy}$ , it holds that  $\frac{1}{2} \Pr[A_{xyz}] \leq \Pr[T_{z|xy}] \leq \Pr[A_{xyz}]$ .

**Lemma 2** Given a pair  $(x, y) \in V \times V$  and a B-quadruple  $\{x, y, z, w\} \in \mathcal{Q}_{xy}^d$ , it holds that  $\frac{1}{6} \Pr[A_{xyzw}] \leq \Pr[Q_{zw|xy}] \leq \frac{1}{4} \Pr[A_{xyzw}]$ .

Analyzing carefully the probabilities of events  $T_{z|xy}$  and  $Q_{zw|xy}$  is crucial for deriving the desired approximation factor, as shown next.

We consider an instance  $G = (V, E, L, \ell)$  of our problem and rewrite the cost function in Equation (2) as sum of the costs paid by any single pair  $(x, y)$ . To this end, in order

to simplify the notation, we hereinafter write the cost by omitting  $\mathcal{C}$  and  $\ell$  while keeping  $G$  only:

$$c(G) = \sum_{(x,y) \in V \times V} c_{xy}(G), \quad (6)$$

where  $c_{xy}(G)$  denotes the aforementioned contribution of the pair  $(x, y)$  to the total cost. Moreover, let  $E[c(G)]$  denote the expected cost of Chromatic Balls over the random choices made by the algorithm. By the linearity of expectation, the expected cost  $E[c(G)]$  can be expressed as

$$E[c(G)] = \sum_{(x,y) \in V \times V} E[c_{xy}(G)]. \quad (7)$$

Finally, let  $c^*(G)$  be the cost of the optimal solution on  $G$ .

To derive an approximation factor  $r(G)$  on the performance of the Chromatic Balls algorithm, we look for an upper bound  $Ub(G)$  on the expected cost  $E[c(G)]$  of the algorithm, and a lower bound  $Lb(G)$  on the cost  $c^*(G)$  of the optimal solution, so that

$$\frac{E[c(G)]}{c^*(G)} \leq \frac{Ub(G)}{Lb(G)} = r(G). \quad (8)$$

We next show how to derive such upper and lower bounds.

**Deriving the upper bound  $Ub(G)$ .** For a pair  $(x, y)$  we define the collection of events  $\Omega_{xy} = \{T_{z|xy} \mid \{x, y, z\} \in \mathcal{T}_{xy}\} \cup \{Q_{zw|xy} \mid \{x, y, z, w\} \in \mathcal{Q}_{xy}^d\}$ . As the following two lemmas show, if pair  $(x, y)$  contributes to the cost paid by the algorithm, then exactly one of the events in  $\Omega_{xy}$  occurs.

**Lemma 3** If  $c_{xy}(\mathcal{C}, \ell, G) > 0$  then at least one of the events in  $\Omega_{xy}$  occurs.

**Lemma 4** The events within the collection  $\Omega_{xy}$  are disjoint.

Combining Lemmas 3 and 4 with the expressions of the probabilities of the events  $T_{z|xy}$  (Lemma 1) and  $Q_{zw|xy}$  (Lemma 2) we can derive an upper bound on the expected contribution  $E[c_{xy}(G)]$  of a pair  $(x, y)$  to the total cost.

**Lemma 5** For a pair  $(x, y) \in V \times V$  the following bound holds.

$$E[c_{xy}(G)] \leq \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \Pr[A_{xyz}] + \sum_{\{x,y,z,w\} \in \mathcal{Q}_{xy}^d} \frac{1}{4} \Pr[A_{xyzw}].$$

The bound in Lemma 5 together with Equation (7) can be used to give the desired (upper) bound on the overall expected cost  $E[c(G)]$ .

**Lemma 6** The expected cost  $E[c(G)]$  of the Chromatic Balls algorithm can be bounded as follows

$$E[c(G)] \leq Ub(G) = \sum_{\{x,y,z\} \in \mathcal{T}} \left( 3 \Pr[A_{xyz}] + \frac{3}{4} X_{xyz} + \frac{1}{2} Y_{xyz} \right),$$

where:

$$X_{xyz} = \sum_{w \in W_{xyz}} \frac{\Pr[A_{xyzw}]}{\mathcal{T}_{xyzw}}, \quad Y_{xyz} = Y_{xyz}^{xy} + Y_{xyz}^{xz} + Y_{xyz}^{yz},$$

$$Y_{xyz}^{xy} = \sum_{w \in W_{xyz}^{xy}} \frac{\Pr[A_{xyzw}]}{\mathcal{T}_{xyzw}}, \quad Y_{xyz}^{xz} = \sum_{w \in W_{xyz}^{xz}} \frac{\Pr[A_{xyzw}]}{\mathcal{T}_{xyzw}},$$

$$\text{and } Y_{xyz} = \sum_{w \in W_{xyz}^{yz}} \frac{\Pr[A_{xyzw}]}{\tau_{xyzw}}.$$

Finally,  $\tau_{xyzw}$  denotes the number of B-triplets contained in any B-quadruple  $\{x, y, z, w\}$ .

**Deriving the lower bound  $Lb(G)$ .** Recalling that a B-triplet incurs a non-zero cost in any solution, a lower bound on the cost of the optimal solution  $c^*(G)$  can be obtained by counting the number of disjoint B-triplets in the input. Considering the set  $\mathcal{T}$  of B-triplets we can restate the following result of Ailon et al. [1] that provides a lower bound on the optimal by “fractionally assigning” all pairs of objects in  $V \times V$  to the triplets in  $\mathcal{T}$ .

**Lemma 7 (Ailon et al. [1])** *Let  $\{\alpha_{xyz} \mid \{x, y, z\} \in \mathcal{T}\}$  be any assignment of nonnegative weights to the B-triplets in  $\mathcal{T}$  satisfying  $\sum_{\{x, y, z\} \in \mathcal{T}_{x'y'}} \alpha_{xyz} \leq 1$  for all  $(x', y') \in V \times V$ . It holds that  $c^*(G) \geq \sum_{\{x, y, z\} \in \mathcal{T}} \alpha_{xyz}$ .*

We can then obtain a lower bound on the optimal solution by finding a suitable set of weights  $\alpha_{xyz}$  that satisfies the conditions of the previous lemma. We derive such a set of weights in the following further lemma.

**Lemma 8** *For any pair  $(x, y) \in V \times V$  the following condition holds.*

$$\sum_{\{x, y, z\} \in \mathcal{T}_{xy}} \frac{1}{1 + |\mathcal{T}_{xy}|} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6} X_{xyz} + \frac{1}{6} Y_{xyz} \right) \leq 1.$$

Thus, combining Lemmas 7 and 8, we can obtain the desired lower bound  $Lb(G)$  as follows.

**Lemma 9** *The cost  $c^*(G)$  of the optimal solution on any input instance  $G$  is lower bounded as follows*

$$\begin{aligned} c^*(G) &\geq Lb(G) = \\ &= \sum_{\{x, y, z\} \in \mathcal{T}} \frac{1}{1 + t_{max}} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6} X_{xyz} + \frac{1}{6} Y_{xyz} \right), \end{aligned}$$

where  $t_{max} = \max_{(x, y) \in V \times V} |\mathcal{T}_{xy}|$  is the maximum number of B-triplets that hit a pair of objects.

**The approximation ratio  $r(G)$ .** The upper and lower bounds obtained in Lemmas 6 and 9 are at the basis if the final form of the approximation ratio of Chromatic Balls.

**Theorem 1** *The approximation ratio of the Chromatic Balls algorithm on any input instance  $G$  is*

$$r(G) = \frac{\mathbb{E}[c(G)]}{c^*(G)} \leq 6(1 + t_{max}),$$

where  $t_{max} = \max_{(x, y) \in V \times V} |\mathcal{T}_{xy}|$  is the maximum number of B-triplets that hit a pair of objects.

Theorem 1 shows that the approximation factor of the Chromatic Balls algorithm is bounded by the maximum number  $t_{max}$  of B-triplets that hit a pair of objects. The result is meaningful as it quantifies the quality of the performance of the algorithm as a property of the input graph. For example, as the following corollary shows, the algorithm provides a constant-factor approximation for bounded-degree graphs.

**Corollary 1** *The approximation ratio of the Chromatic Balls algorithm on any input instance  $G$  is*

$$r(G) \leq 6(2D_{max} - 1),$$

where  $D_{max} = \max_{x \in V} |\{y \mid y \in V \wedge \ell(x, y) \neq l_0\}|$  is the maximum degree in the problem instance.

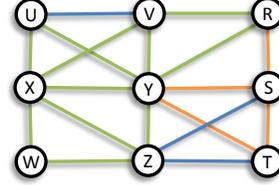
## 4. OTHER ALGORITHMS

In this section we present two additional algorithms for the CHROMATIC-CORRELATION-CLUSTERING problem. The first one is a variant of the Chromatic Balls algorithm that attempts to overcome some weaknesses of Chromatic Balls by employing two heuristics, one for pivot selection and one for cluster selection. The second one is an alternating minimization method that is designed to optimize directly the objective function.

### 4.1 Lazy Chromatic Balls

The algorithm we present next is motivated by the following example, in which we discuss what may go wrong during the execution of the Chromatic Balls algorithm.

**Example 3** *Consider the graph in Figure 2: it has a fairly evident green cluster formed by vertices  $\{U, V, R, X, Y, W, Z\}$ .*



**Figure 2: An example of an edge-labeled graph.**

However, as all the edges have the same probability of being selected as pivots, Chromatic Balls might miss this green cluster, depending on which edge is selected first. For instance, suppose that the first pivot chosen is  $(Y, S)$ . Chromatic Balls forms the red cluster  $\{Y, S, T\}$  and removes it from the graph. Removing vertex  $Y$  makes the edge  $(X, Y)$  missing, which would have been a good pivot to build a green cluster. At this point, even if the second selected pivot edge is a green one, say  $(X, Z)$ , Chromatic Balls would form only a small green cluster  $\{X, W, Z\}$ .

Motivated by the previous example we introduce the Lazy Chromatic Balls heuristic, which tries to minimize the risk of bad choices. Given a vertex  $x \in V$ , and a label  $l \in L$ , let  $d(x, l)$  be the number of edges incident to  $x$  having label  $l$ . Also, we denote by  $\Delta(x) = \max_{l \in L} d(x, l)$ , and  $\lambda(x) = \arg \max_{l \in L} d(x, l)$ . Lazy Chromatic Balls differs from Chromatic Balls in two ways:

**Pivot random selection.** At each iteration Lazy Chromatic Balls selects a pivot edge in two steps. First, a vertex  $u$  is picked up with probability directly proportional to  $\Delta(u)$ . Then, a second vertex  $v$  is selected among the neighbors of  $u$  with probability proportional to  $d(v, \lambda(u))$ .

**Ball formation.** Given the pivot  $(u, v)$ , Chromatic Balls forms a cluster by adding all vertices  $x$  such that  $\langle u, v, x \rangle$  is a monochromatic triangle. Lazy Chromatic Balls instead, iteratively adds vertices  $x$  in the cluster as long as they form a triangle  $\langle X, Z, w \rangle$  of color  $\ell(u, v)$ , where  $X$  is either  $u$  or  $v$ , and  $Z$  can be any other vertex already belonging to the current cluster.

**Example 4** Consider again the example in Figure 2. Vertices  $X$  and  $Y$  have the maximum number of edges of one color: they both have 5 green edges. Hence, one of them is chosen as first pivot vertex  $u$  by **Lazy Chromatic Balls** with higher probability than the remaining vertices. Suppose that  $X$  is picked up, i.e.,  $u = X$ . Given this choice, the second pivot  $v$  is chosen among the neighbors of  $X$  with probability proportional to  $d(v, \lambda(u))$ , i.e., the higher the number of green edges of the neighbor, the higher the probability for it to be chosen. In this case, hence, **Lazy Chromatic Balls** would likely choose  $Y$  as a second pivot vertex  $v$ , thus making  $(X, Y)$  the selected pivot edge. Afterwards, **Lazy Chromatic Balls** adds to the being formed cluster the vertices  $\{U, V, Z\}$  because each of them forms a green triangle with the pivot edge. Then,  $R$  enters the cluster too, because it forms a green triangle with  $Y$  and  $V$ , which is already in the cluster. Similarly,  $W$  enters the cluster thanks to  $Z$ .

**Computational complexity.** Like Chromatic Balls, the running time of the Lazy Chromatic Balls algorithm is determined by picking the pivots and building the various clusters. Picking the first pivot  $u$  can be implemented with a priority queue with priorities  $\Delta \times \text{rnd}$ , where  $\text{rnd}$  is a random number. This requires computing  $\Delta$  for all objects, which takes  $\mathcal{O}(nh + m)$  (where  $h = |L|$ ). Managing the priority queue itself requires instead  $\mathcal{O}(n \log n)$ , as each object is put into/removed from the queue only once during the execution of the algorithm. Given  $u$ , the second pivot  $v$  is selected by probing all (non-chosen) neighbors of  $u$ . This takes  $\mathcal{O}(m)$  time, as for each pivot  $u$ , its neighbors are accessed only once throughout the execution of the algorithm. Finally, building the current cluster takes  $\mathcal{O}(m)$  time, as it requires a visit of the graph, where each edge is accessed  $\mathcal{O}(1)$  times. In conclusion, the computational complexity of Lazy Chromatic Balls is  $\mathcal{O}(n(\log n + h) + m)$ , which, for small  $h$ , is better than the complexity of Chromatic Balls.

## 4.2 An alternating-minimization approach

A nice feature of the previous algorithms is that they are parameter-free: they produce clusterings by using information that is local to the pivot edges, without forcing the number of output clusters in any way. However, in some cases, it could be desired having an output clustering composed by a pre-specified number  $K$  of clusters. To this purpose, we present here an algorithm based on the *alternating minimization* paradigm [7], that receives in input the number  $K$  of output clusters and attempts to minimize Equation (2) directly. The pseudocode of the proposed algorithm, called **Alternating Minimization**, is given in Algorithm 2.

In a nutshell, **AM** tries to produce a solution by alternating between two optimization steps. In the first step the algorithm finds the best cluster assignment for every  $x \in V$  given the assignments of every other  $y \in V$  and the current cluster labels. In the second step, it finds the best label for every cluster given the current assignment of objects to clusters. Below we show that both steps can be solved optimally. As a consequence the value of Equation (2) is guaranteed to decrease in every step, until convergence. Finding the global minimum is obviously hard, but the algorithm is guaranteed to converge to a local optimum.

**Definitions.** For presentation sake, we adopt matrix notation. We denote matrices by uppercase boldface romans and vectors by lowercase boldface romans. We write  $\mathbf{X}_{ij}$

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### Algorithm 2 Alternating Minimization (AM)

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**Input:** Edge-labeled graph  $G = (V, E, L, \ell)$ ;  
number  $K$  of output clusters  
**Output:** Clustering  $\mathcal{C} : V \rightarrow \mathbb{N}$ ; cluster labeling function  $cl : \mathcal{C}[V] \rightarrow L$   
1: initialize  $\mathbf{A} = [\mathbf{a}_1, \dots, \mathbf{a}_N]$  and  $\mathbf{C} = [\mathbf{c}_1, \dots, \mathbf{c}_K]$  at random  
2: **repeat**  
3: for all  $x \in V$  compute optimal  $\mathbf{a}_x$  according to Proposition 1  
4: for all  $k \in [1..K]$  compute optimal  $\mathbf{c}_k$  according to Proposition 2  
5: **until** neither  $\mathbf{A}$  nor  $\mathbf{C}$  changed

---

for the  $(i, j)$  coordinate of matrix  $\mathbf{X}$ , and  $\mathbf{x}(i)$  for the  $i$ -th coordinate of vector  $\mathbf{x}$ .

The parameter space of Problem 2 consists of a *cluster assignment* for every object  $x \in V$ , given by the binary matrix  $\mathbf{A}$ , and a *label assignment* for every cluster  $k \in \{1, \dots, K\}$ , given by the binary matrix  $\mathbf{C}$ . We have  $\mathbf{A}_{kx} = 1$  when object  $x$  is assigned to cluster  $k$ , and  $\mathbf{A}_{kx} = 0$  otherwise. Similarly, we set  $\mathbf{C}_{lk} = 1$  when label  $l$  is assigned to cluster  $k$ , and  $\mathbf{C}_{lk} = 0$  otherwise. Since every object must belong to one and only one cluster, and every cluster must have one and only one label assigned, both  $\mathbf{A}$  and  $\mathbf{C}$  are constrained to consist of all zeros with a single 1 on every column. Denote by  $\mathbf{a}_x$  the column of  $\mathbf{A}$  that corresponds to object  $x$ .

The input is represented by a set of binary matrices, with a matrix  $\mathbf{Z}_x$  for every  $x \in V$ . These matrices encode the labeling function  $\ell$  as follows. Let  $\mathbf{z}_{xy}$  denote the column of  $\mathbf{Z}_x$  that corresponds to the object  $y \in V$ . We have  $\mathbf{z}_{xy}(l) = 1$  if and only if  $\ell(x, y) = l$ , otherwise  $\mathbf{z}_{xy}(l) = 0$ . Every  $\mathbf{Z}_x$  consists thus of zeros, with exactly one 1 on every column. Finally, denote by  $\mathbf{b}$  a special binary vector where  $\mathbf{b}(l) = 1$  when  $l = l_0$  and  $\mathbf{b}(l) = 0$  otherwise. We have then  $\mathbf{z}_{xy}^T \mathbf{b} = 1$  if and only if  $\ell(x, y) = l_0$ .

The above formulation of the problem assumes that the input is represented by many large matrices. Note however that this representation is only conceptual. In the actual implementation we do not have to materialize these matrices and we can represent the input with the minimal amount of space required, as shown next. The benefit of our formulation is that it allows to write our objective function and our optimization process using linear-algebra operations, and argue about the optimality of the local optimization steps.

**Optimal cluster assignment.** Denote by  $N_{xk}^-$  the number of objects  $y \in V$  in cluster  $k$  that have  $\ell(x, y) = l_0$ . Since  $\ell(x, y) = l_0 \Leftrightarrow \mathbf{z}_{xy} \mathbf{b} = 1$ , we have  $N_{xk}^- = (\mathbf{A} \mathbf{Z}_x \mathbf{b})(k)$ . Similarly, let  $N_{xk}^+$  denote the number of objects  $y \in V$  in cluster  $k$  that have  $\ell(x, y) = cl(k)$ . Since  $y \in k$ , we have  $\ell(x, y) = cl(k) \Leftrightarrow \mathbf{z}_{xy} \mathbf{C} \mathbf{a}_y = 1$  and can write  $N_{xk}^+ = (\mathbf{A} \mathbf{w}_x)(k)$ , where  $\mathbf{w}_x = [\mathbf{z}_{x1}^T \mathbf{C} \mathbf{a}_1 \dots \mathbf{z}_{xn}^T \mathbf{C} \mathbf{a}_n]$ .

**Proposition 1** *The optimal cluster assignment for  $x \in V$  given  $\mathbf{A}$  and  $\mathbf{C}$  is  $k^* = \arg \min_k N_{xk}^- - N_{xk}^+$ .*

**PROOF.** We can rewrite Equation (2) as follows:

$$\sum_{x,y} \mathbf{a}_x^T \mathbf{a}_y (1 - \mathbf{z}_{xy}^T \mathbf{C} \mathbf{a}_y) + (1 - \mathbf{a}_x^T \mathbf{a}_y) (1 - \mathbf{z}_{xy}^T \mathbf{b}) = \quad (9)$$

$$= \sum_x \mathbf{a}_x^T \mathbf{A} (\mathbf{1} - \mathbf{w}_x) + (\mathbf{1}^T - \mathbf{a}_x^T \mathbf{A}) (\mathbf{1} - \mathbf{Z}_x^T \mathbf{b}),$$

where  $\mathbf{w}_x$  is defined as above, and  $\mathbf{1}$  denotes the  $|V|$ -dimensional vector of all 1s. Terms that correspond to a

fixed  $x \in V$  further simplify to

$$\mathbf{a}_x^T \mathbf{A}^T \mathbf{Z}_x \mathbf{b} - \mathbf{a}_x^T \mathbf{A} \mathbf{w}_x + d_x,$$

where the constant  $d_x = \mathbf{1}^T \mathbf{1} - \mathbf{1}^T \mathbf{Z}_x^T \mathbf{b}$  is the ‘‘degree’’ of object  $x$ , the number of objects  $y \in V$  where  $\ell(x, y) \neq l_0$ . Since we must assign exactly one cluster for  $x$ , the above expression is minimized simply by assigning  $x$  to the cluster  $k$  that minimizes  $(\mathbf{A} \mathbf{Z}_x^T \mathbf{b})(k) - (\mathbf{A} \mathbf{w}_x)(k) = N_{xk}^- - N_{xk}^+$ .

The result is quite intuitive. The best cluster for  $x$  is the one having the least ‘‘push’’ in terms of  $l_0$  connections, and the most ‘‘pull’’ given by connections having the appropriate label. However, evaluating  $N_{xk}^-$  in practice is very slow, as it involves checking all  $l_0$  connections of  $x$ . Ideally the update rule should only require access to edges having some label other than  $l_0$ . This is easy to achieve, however.

Let  $N_{xk}^0$  denote the remaining objects in cluster  $k$ , that is, those with  $\ell(x, y) \neq \ell(k) \neq l_0$ . Also, let  $S_k$  denote the size of cluster  $k$ . Clearly we have  $S_k = N_{xk}^+ + N_{xk}^0 + N_{xk}^-$  for every  $x \in V$ . Using this we obtain  $N_{xk}^- - N_{xk}^+ = S_k - 2N_{xk}^+ - N_{xk}^0$ , which is much faster to evaluate.

**Optimal label assignment.** The update rule for the cluster label is intuitive as well. Denote by  $E_k$  the number of ordered  $(x, y)$  pairs so that both  $x$  and  $y$  belong to cluster  $k$ , and  $\ell(x, y) = \ell(k)$ .

**Proposition 2** *The optimal label assignment for cluster  $k$  given  $\mathbf{A}$  is  $l^* = \arg \min_l S_k^2 - E_k$ .*

PROOF. We can partition the cost in Equation (9) as a sum over clusters. That is, for a fixed cluster  $k$  we sum only over those  $x$  and  $y$  that belong both to  $k$ . Also, the second term in Equation (9) does not depend on  $\mathbf{C}$  and can therefore be omitted. This leaves us with the sum

$$\sum_{x \in k, y \in k} (1 - \mathbf{z}_{xy}^T \mathbf{C} \mathbf{a}_y),$$

where we can replace  $\mathbf{C} \mathbf{a}_y$  with the binary vector  $\mathbf{c}_k$  that indicates the label assigned to cluster  $k$ . Clearly we have  $\sum_{x \in k, y \in k} 1 = S_k^2$ , and it is easy to see that  $\sum_{x \in k, y \in k} \mathbf{z}_{xy} \mathbf{c}_k$  counts all  $(x, y)$  pairs having the same label that is currently assigned to  $k$ , which is by definition equal to  $E_k$ .

This means that the optimal label for cluster  $k$  is simply the label shared by the majority of the pairs in  $k$ .

**Computational complexity.** The running time of Alternating Minimization depends on the (optimal) cluster and label assignment steps. Cluster assignment requires two sub-steps: evaluating  $S_k - 2N_{xk}^+ - N_{xk}^0$  for each vertex and cluster, which can be performed in  $\mathcal{O}(m)$  by a simple visit of the input graph, and looking at all clusters to choose the best one for each vertex, which clearly takes  $\mathcal{O}(Kn)$ . Label assignment requires to compute the number of intra-cluster edge labels for each cluster  $k$  and label  $l$ . This takes  $\mathcal{O}(m)$ , as it can be performed, again, by visiting the input graph. Then, the assignment of labels to clusters by evaluating  $S_k^2 - E_k$  can be performed in  $\mathcal{O}(Kh)$ . In conclusion, as usually  $h = |L| \ll n$ , the computational complexity of Alternating Minimization is  $\mathcal{O}(s(Kn + m))$ , where  $s$  is the number of iterations to convergence.

## 5. EXPERIMENTAL EVALUATION

In this section, we report our empirical assessments. We experiment with all three proposed algorithms, **Chromatic**

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### Algorithm 3 Synthetic data generator

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**Input:** number of vertices  $n$ , number of clusters  $K$ , number of labels  $h$ , probability  $p$  of intra-cluster edges, probability  $q$  of inter-cluster edges, probability  $w$  that an edge inside a cluster has a color different from the cluster

**Output:** edge labeled graph  $G = (V, E, L, \ell)$

- 1:  $V \leftarrow [1, n]$ ,  $E \leftarrow \emptyset$ ,  $L \leftarrow \{l_1, \dots, l_h\}$
- 2: assign each vertex  $x \in V$  to a randomly selected cluster
- 3: assign to each cluster a randomly selected label from  $L$
- 4: **for all** pairs  $(x, y) \in V \times V$  **do**
- 5:   pick 3 random numbers  $r_1, r_2, r_3$  ranging within  $[0, 1]$
- 6:   **if**  $\mathcal{C}(x) = \mathcal{C}(y)$  **then**
- 7:     **if**  $r_1 < p$  **then**
- 8:       **if**  $r_2 < w$  **then**
- 9:          $E \leftarrow E \cup (x, y)$
- 10:         $\ell(x, y) \leftarrow$  a random label from  $L \setminus \{\mathcal{C}(x)\}$
- 11:     **else**
- 12:         $E \leftarrow E \cup (x, y)$
- 13:         $\ell(x, y) \leftarrow \mathcal{C}(x)$ ;
- 14:     **else if**  $r_3 < q$  **then**
- 15:         $E \leftarrow E \cup (x, y)$
- 16:         $\ell(x, y) \leftarrow$  a random label from  $L$

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**Balls**, **Lazy Chromatic Balls**, and **Alternating Minimization**, to which we refer by **CB**, **LCB**, and **AM**, respectively. We also evaluate the performance of the baseline described in the Introduction, namely the ‘‘standard’’ **Balls** algorithm [1] that ignores colors. We refer to this baseline as **B**. All measurements reported are averaged over 50 runs.

### 5.1 Experiments on synthetic data

We evaluate our algorithms on synthetic datasets generated by the process outlined in Algorithm 3. In a nutshell, the generator initially assigns vertices and labels to clusters uniformly at random, and then adds noise according to the probability parameters  $p$ ,  $q$ , and  $w$ . Given the assignment of vertices to clusters, intra-cluster edges are sampled with probability  $p$ , and they are given the correct label (the label of the cluster they are assigned to) with probability  $1 - w$ , while, inter-cluster edges are sampled with probability  $q$ .

The initial assignment of objects and labels to clusters can be interpreted as a ground truth underlying the corresponding synthetic dataset. We compare the resulting clusterings with the ground-truth clustering using the well-known  $F$ -measure external cluster-validity criterion. Given a ground-truth clustering  $\hat{\mathcal{C}}$  and a clustering solution  $\mathcal{C}$  having  $\hat{K}$  and  $K$  clusters, respectively,  $F$ -measure is defined in terms of *precision* and *recall* as follows:

$$F(\mathcal{C}, \hat{\mathcal{C}}) = \frac{1}{n} \sum_{\hat{k}=1}^{\hat{K}} S_{\hat{k}} \max_{k \in [1..K]} F_{\hat{k}k},$$

where  $F_{\hat{k}k} = (2P_{\hat{k}k}R_{\hat{k}k})/(P_{\hat{k}k} + R_{\hat{k}k})$  such that  $P_{\hat{k}k} = S_{\hat{k} \cap k} / S_k$  and  $R_{\hat{k}k} = S_{\hat{k} \cap k} / S_{\hat{k}}$ , while  $S_{\hat{k} \cap k}$  denotes the number of common objects between the  $\hat{k}$ -th cluster of  $\hat{\mathcal{C}}$  and the  $k$ -th cluster of  $\mathcal{C}$ , and  $S_{\hat{k}}$  and  $S_k$  are the sizes of clusters  $\hat{k}$  and  $k$ , respectively. It is easy to see that  $F \in [0, 1]$ .

We generate datasets with a fixed number of objects ( $n = 1000$ ), and we vary (i) the noise level (controlled by playing with  $p$ ,  $q$ , and  $w$ ); (ii) the number of labels  $h$ ; and (iii) the number of clusters  $K$  in the ground truth. Even though we perform tests by varying all parameters  $p$ ,  $q$ , and  $w$ , due to space limitations we only report results obtained for varying  $q$  and keeping  $p$  and  $w$  equal to 0.5.

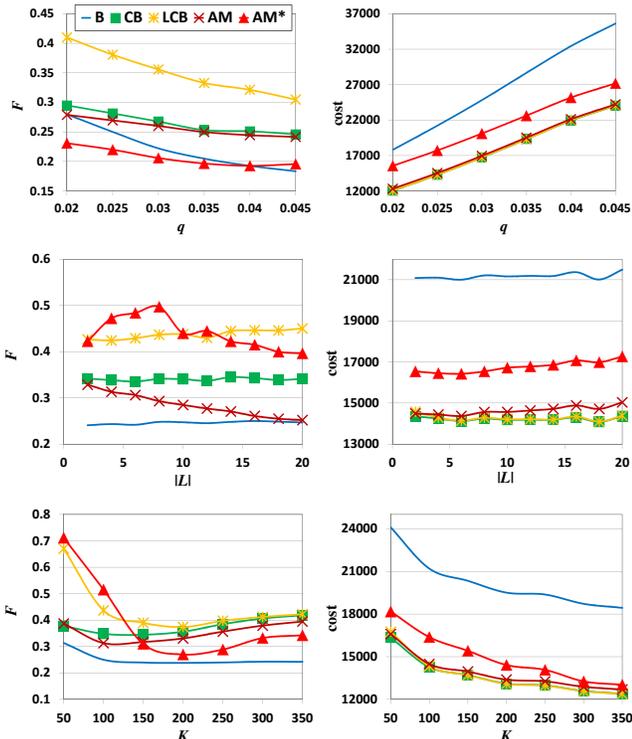


Figure 3: Accuracy on synthetic datasets in terms of  $F$ -measure (left) and solution cost (right), by varying level of noise (1st row), number of labels (2nd row), and number of ground-truth clusters (3rd row).

For the number of clusters required as input for the AM algorithm, we consider two options: the average number of clusters produced by the CB algorithm, and the number of clusters in the ground truth. We refer to these two settings by AM and AM\*, respectively. In Figure 3 we report the performance of our algorithms in terms of  $F$ -measure, as well as solution cost (Equation (2)).

All trends observed by varying the parameters  $q$ ,  $h$ , and  $K$  are intuitive. Indeed, for all methods, the performance decreases as the noise level  $q$  increases (Figure 3, 1st row). On the other hand, all methods give better solutions, in terms of cost, as the number of ground truth clusters  $K$  increases (Figure 3, 3rd row right). The reason is that since CB and LCB tend to produce a large number of clusters, by setting a larger  $K$  the difference tends to disappear.

We also note that, in terms of  $F$ -measure, CB, LCB, and AM perform better than AM\*. The poorer performance of AM\* can again be explained by the input number of clusters, which is likely far away from the “ideal” number of clusters according to the objective function cost. In terms of  $F$ -measure scores, LCB outperforms all other methods, followed by CB. The performance of AM and AM\* is variable, while remaining generally better than the baseline.

## 5.2 Experiments on real data

We experiment with three real datasets (Table 1).

**String.** A protein-protein interaction (PPI) network obtained from `string-db.org`, i.e., a database of known protein interactions for a large number of organisms. The dataset is an undirected graph where vertices represent proteins and edges represent protein interactions. Edges are labeled with 4 types of interactions. The PPI datasets are usually very sparse, therefore, we keep only the 30-core of the entire network, i.e., we recursively remove the vertices with degree less than 30 until a fix-point has been reached.

**Youtube.** This dataset represents a network of associations in the `youtube` site. The vertices of the network represent

Table 1: Characteristics of real data.  $n$ : number of vertices;  $m$ : number of edges;  $d$ : average degree;  $|L|$ : number of labels;  $c$ : clustering coefficient.

dataset	$n$	$m$	$d$	$ L $	$c$
String	18 152	401 582	44.25	4	0.731
Youtube	15 088	19 923 067	2 640.92	5	0.495
DBLP	312 416	2 110 470	13.51	100	0.204

users, videos, and channels. Entities in the network have five different types of associations: *contact*, *co-contact*, *co-subscription*, *co-subscribed*, and *favorite*; these are the edge labels considered by our algorithms. For edges with multiple labels we picked one label at random from the available ones. The dataset has been compiled by Tang et al. [14] and it is available at <http://www.public.asu.edu/~ltang9/>.

**DBLP.** We obtain a recent snapshot of the DBLP co-authorship network (<http://dblp.uni-trier.de/xml/>). For each co-authorship edge, we consider the bag of words obtained by merging the titles of all papers coauthored by the two authors. Words are stemmed and stop-words are removed. We then apply *Latent Dirichlet Allocation* (LDA) [5] to automatically identify 100 topics on each edge. After LDA topic-modeling, for each edge, we assign its most prominent topic discovered as edge label.

**Results.** Table 2 summarizes the results obtained on real data. Like in synthetic data, all proposed algorithms clearly outperform the baseline B. CB is the best method on Youtube and DBLP, achieving up to 27.74% of improvement with respect to the baseline in terms of solution cost. Instead, CB is slightly outperformed by LCB and AM on String, while LCB outperforms AM on String and DBLP.

As far as the runtime, we observe that the baseline is faster than the proposed methods, as expected. This is mainly due to a smaller complexity in choosing vertex pivots compared to choosing edge pivots. However, all proposed methods remain very efficient, as they take a few seconds (CB and LCB) or minutes (AM) on large and dense graphs like Youtube and DBLP. All runtimes comply with the computational complexity analysis reported previously. Indeed, AM is the slowest method, mostly due to the typically high number of iterations needed to convergence, while LCB is faster than CB, especially on dense datasets like Youtube.

Finally, Figure 4 shows an example cluster from the DBLP co-authorship network recognized by the LCB algorithm, containing 23 authors (vertices). Among the 71 intra-cluster edges, 58 have the same label, i.e., Topic 18, whose most representative (stemmed) keywords are: *queri*, *effici*, *spatial*, *tempor*, *search*, *index*, *similar*, *data*, *dimension*, *aggreg*. Other topics (edge colors) that appears are “sensor networks”, “frequent pattern mining”, “algorithms on graphs and trees”, “support vector machines”, “classifiers and Bayesian learning”.

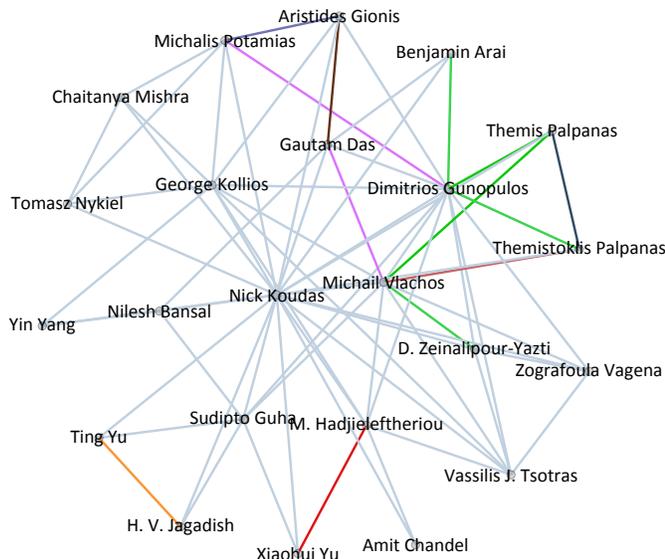
## 6. RELATED WORK

**Edge-labeled graphs and multidimensional networks.** Graphs in which edges are labeled with a type of relation occurring among the connected vertices are receiving increasingly attention. To the best of our knowledge no previous work has investigated the problem of clustering in such graphs. The problems studied so far on this kind of graphs are mainly on label-constrained reachability queries [10, 12, 16, 8], whose main goal is to answer whether a vertex  $u$  can reach vertex  $v$  through a path whose edge labels belong to a given set.

Clustering has been studied, instead, in so called *multidimensional networks*, i.e., networks defined as a collection of multiple networks over the same set of actors. In our jargon these are simply graphs where each edge can have more than

**Table 2: Results on real datasets: average cost, runtime (s), and average number of output clusters**

dataset	cost				runtime (s)				#clusters			
	B	CB	LCB	AM	B	CB	LCB	AM	B	CB	LCB	AM
String	163 305	160 060	155 881	156 976	0.5	1.4	1.3	21.0	1086	1451	784	1451
Youtube	23 550 213	18 956 000	22 644 858	19 670 899	22.4	117.8	40.5	1 038.9	568	1 078	672	1 078
DBLP	2 260 065	1 633 149	1 678 714	2 018 952	4.3	10.2	5.5	2 116.1	66 276	123 197	99 948	123 197



**Figure 4: An example cluster from DBLP.**

one color [15, 4, 13]. Although the input of that problem might seem close to ours, the objective is semantically far away. In clustering multidimensional networks, the objective is to find a partitioning of vertices which is meaningful and relevant in *all dimensions at the same time*. Taking again the colors metaphor, in that setting is a clustering is considered as good if it makes sense in the **green** network *and* as well as the **red** network, and so on. In our work, we are rather interested in finding groups of objects that induce color-coherent clusters while looking at all the colors together.

**Correlation Clustering.** The problem of CORRELATION-CLUSTERING was first defined by Bansal et al. [3] in its binary version. Ailon et al. [1] proposed the Balls algorithm that achieves expected approximation factor 5 if the weights obey the probability condition. If the weights  $X_{ij}$  obey also the triangle inequality, then the algorithm achieves expected approximation factor 2. Giotis and Guruswami [9] consider correlation clustering when the number of clusters is given, while Ailon and Liberty [2] study a variant of correlation clustering where the goal is to minimize the number of disagreements between the produced clustering and a given ground truth clustering. We recently extended correlation clustering to allow overlaps, i.e., objects belonging to more than one cluster [6].

## 7. CONCLUSIONS

In this paper, we introduce a variant of the correlation-clustering problem, in which the pairwise relations between objects are categorical. The problem has interesting applications, such as clustering social networks where individuals are connected with different types of relations, or clustering protein networks, where proteins are associated with different types of interactions. We propose three algorithms that

we evaluate on synthetic and real datasets.

Our problem is a novel clustering formulation well-suited for mining multi-labeled and heterogeneous datasets that are becoming increasingly common. We believe that there are many interesting extensions and fruitful future research directions. For example, we would like to extend the problem formulation in order to capture overlapping clusters as well as multiple-labeled edges.

**Acknowledgements.** This research was partially supported by the Torres Quevedo Program of the Spanish Ministry of Science and Innovation, co-funded by the European Social Fund, and by the Spanish Centre for the Development of Industrial Technology under the CENIT program, project CEN-20101037, “Social Media” (<http://www.centisocialmedia.es/>).

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## APPENDIX

### A. DETAILS ABOUT THE THEORETICAL ANALYSIS OF THE Chromatic Balls ALGORITHM

#### A.1 Proofs of Lemmas 1–2

To prove Lemmas 1 and 2, we need to introduce the following additional Lemma 10 and Corollaries 2-3 to show the disjointness of the events  $A_{x_1 \dots x_n}^{(i)}$ ,  $T_{z|xy}^{(i)}$ , and  $Q_{zw|xy}^{(i)}$  with respect to the iterations of the algorithm.

**Lemma 10** *Given any  $\{x_1, \dots, x_n\} \subseteq V$ ,  $n \geq 2$ , it holds that  $A_{x_1 \dots x_n}^{(i)}$  and  $A_{x_1 \dots x_n}^{(j)}$  are disjoint, for any two iterations  $i$  and  $j$  of Chromatic Balls such that  $i \neq j$ .*

PROOF. As soon as  $A_{x_1 \dots x_n}^{(i)}$  happens at any iteration  $i$ , exactly two objects in  $\{x_1, \dots, x_n\}$  are chosen as pivots and therefore removed from the set of non-chosen objects that inputs the next iteration(s). Thus, the set  $\{x_1, \dots, x_n\}$  is no longer available as a whole after  $i$ ; this implies that no  $A_{x_1 \dots x_n}^{(j)}$  may occur for any  $j \neq i$ , because this would require that all objects within  $\{x_1, \dots, x_n\}$  input iteration  $j$ .  $\square$

**Corollary 2** *Given any B-triplet  $\{x, y, z\}$ , it holds that  $T_{z|xy}^{(i)}$  and  $T_{z|xy}^{(j)}$  are disjoint, for any two iterations  $i$  and  $j$  such that  $i \neq j$ .*

**Corollary 3** *Given any B-quadruple  $\{x, y, z, w\}$ , it holds that  $Q_{zw|xy}^{(i)}$  and  $Q_{zw|xy}^{(j)}$  are disjoint, for any two iterations  $i$  and  $j$  such that  $i \neq j$ .*

**Lemma 1** *Given a pair  $(x, y) \in V \times V$  and a B-triplet  $\{x, y, z\} \in \mathcal{T}_{xy}$ , it holds that  $\frac{1}{2} \Pr[A_{xyz}] \leq \Pr[T_{z|xy}] \leq \Pr[A_{xyz}]$ .*

PROOF. Given the disjointness conditions proved in Lemma 10 and Corollary 2, it can be noted that:

$$\Pr[A_{xyz}] = \Pr\left[\bigvee_i A_{xyz}^{(i)}\right] = \sum_i \Pr[A_{xyz}^{(i)}]$$

and

$$\Pr[T_{z|xy}] = \Pr\left[\bigvee_i T_{z|xy}^{(i)}\right] = \sum_i \Pr[T_{z|xy}^{(i)}]$$

Thus, it holds that:

$$\begin{aligned} \Pr[T_{z|xy}] &= \Pr\left[\bigvee_i T_{z|xy}^{(i)}\right] = \sum_i \Pr[T_{z|xy}^{(i)}] = \\ &= \sum_i \Pr\left[A_{xyz}^{(i)} \wedge \neg\left(P_x^{(i)} \wedge P_y^{(i)}\right)\right] = \\ &= \sum_i \underbrace{\Pr\left[\neg\left(P_x^{(i)} \wedge P_y^{(i)}\right) \mid A_{xyz}^{(i)}\right]}_p \Pr[A_{xyz}^{(i)}] \quad (10) \end{aligned}$$

The probability  $p = \Pr\left[\neg\left(P_x^{(i)} \wedge P_y^{(i)}\right) \mid A_{xyz}^{(i)}\right]$  depends on how the B-triplet  $\{x, y, z\}$  is composed. Three cases may arise:

1. No pair of objects within  $\{x, y, z\}$  is labeled with  $l_0$ . In this case, all possible outcomes concerning the choice of the pivots are three:  $(x, y)$ ,  $(x, z)$ , and  $(y, z)$ . Among

these, only the last two make the event  $\neg\left(P_x^{(i)} \wedge P_y^{(i)}\right)$  true. All outcomes have equal probability as the random choice of the pivot is uniform in Chromatic Balls. This gives  $p = \frac{2}{3}$  probability.

2.  $(x, y)$  is an  $l_0$ -labeled pair. In this case, there are only two possible outcomes for the pivot choice, because an  $l_0$ -labeled pair cannot be picked up. These choices correspond to the pairs  $(x, z)$  and  $(y, z)$  which both make the event  $\neg\left(P_x^{(i)} \wedge P_y^{(i)}\right)$  true. Thus,  $p = 1$  in this case.
3. Either  $(x, z)$  or  $(y, z)$  is an  $l_0$ -labeled pair. Suppose  $\ell(x, z) = l_0$  (an analogous reasoning holds if  $\ell(y, z) = l_0$ ). Again, there are only two possible choices for the pivot, that are in this case  $(x, y)$  and  $(y, z)$ . Among these, only the latter make the event  $\neg\left(P_x^{(i)} \wedge P_y^{(i)}\right)$  true. This gives  $p = \frac{1}{2}$ .

The above reasoning implies  $\frac{1}{2} \leq p \leq 1$ ; hence, Equation (10) can be rewritten as follows:

$$\begin{aligned} \Pr[T_{z|xy}] &= \sum_i \Pr\left[\neg\left(P_x^{(i)} \wedge P_y^{(i)}\right) \mid A_{xyz}^{(i)}\right] \Pr[A_{xyz}^{(i)}] \Rightarrow \\ &\Rightarrow \frac{1}{2} \sum_i \Pr[A_{xyz}^{(i)}] \leq \Pr[T_{z|xy}] \leq \sum_i \Pr[A_{xyz}^{(i)}] \Leftrightarrow \\ &\Leftrightarrow \frac{1}{2} \Pr\left[\bigvee_i A_{xyz}^{(i)}\right] \leq \Pr[T_{z|xy}] \leq \left[\bigvee_i A_{xyz}^{(i)}\right] \Leftrightarrow \\ &\Leftrightarrow \frac{1}{2} \Pr[A_{xyz}] \leq \Pr[T_{z|xy}] \leq \Pr[A_{xyz}] \quad \square \end{aligned}$$

**Lemma 2** *Given a pair  $(x, y) \in V \times V$  and a B-quadruple  $\{x, y, z, w\} \in \mathcal{Q}_{xy}^d$ , it holds that  $\frac{1}{6} \Pr[A_{xyzw}] \leq \Pr[Q_{zw|xy}] \leq \frac{1}{4} \Pr[A_{xyzw}]$ .*

PROOF. The proof is similar to Lemma 1. Indeed, we exploit the disjointness results shown in Lemma 10 and Corollary 3, and note that:

$$\begin{aligned} \Pr[Q_{zw|xy}] &= \Pr\left[\bigvee_i Q_{zw|xy}^{(i)}\right] = \sum_i \Pr[Q_{zw|xy}^{(i)}] = \\ &= \sum_i \Pr\left[A_{xyzw}^{(i)} \wedge (\neg P_x^{(i)} \wedge \neg P_y^{(i)})\right] = \\ &= \sum_i \Pr\left[(\neg P_x^{(i)} \wedge \neg P_y^{(i)}) \mid A_{xyzw}^{(i)}\right] \Pr[A_{xyzw}^{(i)}] \Rightarrow \\ &\Rightarrow \frac{1}{6} \sum_i \Pr[A_{xyzw}^{(i)}] \leq \Pr[Q_{zw|xy}] \leq \frac{1}{4} \sum_i \Pr[A_{xyzw}^{(i)}] \Leftrightarrow \quad (11) \\ &\Rightarrow \frac{1}{6} \Pr\left[\bigvee_i A_{xyzw}^{(i)}\right] \leq \Pr[Q_{zw|xy}] \leq \frac{1}{4} \Pr\left[\bigvee_i A_{xyzw}^{(i)}\right] \Leftrightarrow \\ &\Leftrightarrow \frac{1}{6} \Pr[A_{xyzw}] \end{aligned}$$

where (11) is derived in a way similar to Lemma 1. Indeed, given  $A_{xyzw}^{(i)}$ , all possible choices of pivots may vary from four of the six pairs of objects within  $\{x, y, z, w\}$  (if two of them are labeled with  $l_0$ ), up to all six pairs (if none of them is labeled with  $l_0$ ). Among these choices, only one (i.e.,  $(z, w)$ ) guarantees  $\neg P_x^{(i)} \wedge \neg P_y^{(i)}$  true. This gives probability values ranging from  $\frac{1}{6}$  to  $\frac{1}{4}$ .  $\square$

## A.2 Proof of Lemma 3

**Lemma 3** *If  $c_{xy}(C, \mathcal{cl}, G) > 0$  then at least one of the events in  $\Omega_{xy}$  occurs.*

PROOF. According to the cost function defined in Equation (2),  $c_{xy}(G) > 0$  if and only if either 1)  $x$  and  $y$  are put in different clusters while  $\ell(x, y) \neq l_0$ , or 2)  $x$  and  $y$  belong to the same cluster  $C$  while  $\ell(x, y)$  is not equal to the label of  $C$ . Let us analyze both cases next.

1) According to the outline of Chromatic Balls,  $(x, y)$  is split when, at some iteration  $i$ , it happens that  $x$  is put into the being formed cluster  $C$ , while  $y$  does not or vice versa. Assuming that the object chosen to belong to  $C$  is  $x$  (an analogous reasoning holds considering  $y$  as belonging to  $C$ ), we have two further cases:

(a)  $x$  is chosen as pivot at the iteration  $i$ , along with any other object  $z \neq y$ . Thus, both the events  $A_{xyz}^{(i)}$  and  $\neg(P_x^{(i)} \wedge P_y^{(i)})$  are true. Also, as  $(x, y)$  is split, either  $\ell(x, y) \neq \ell(x, z)$  or  $\ell(x, z) \neq \ell(x, y)$  (cf. Line 5 in Algorithm 1); hence,  $\{x, y, z\}$  must be a B-triplet hitting  $(x, y)$ . Combining these results and resorting to Equation (4), it results that:

$$\begin{aligned} & \left( A_{xyz}^{(i)} \wedge \neg \left( P_x^{(i)} \wedge P_y^{(i)} \right) \right) \Rightarrow \\ & \Rightarrow \bigvee_i \left( A_{xyz}^{(i)} \wedge \neg \left( P_x^{(i)} \wedge P_y^{(i)} \right) \right) \Leftrightarrow \\ & \Leftrightarrow \bigvee_i T_{z|xy}^{(i)} \Leftrightarrow T_{z|xy}. \end{aligned}$$

(b) The pivots chosen at the iteration  $i$  are  $z$  and  $w$ , with  $z \neq x, z \neq y, w \neq x, w \neq y$ . In this case, it is easy to see that  $\{x, z, w\}$  is a SC-triplet and both  $A_{xyzw}^{(i)}$  and  $\neg P_x^{(i)} \wedge \neg P_y^{(i)}$  are true. Also,  $\{y, z, w\}$  must be a B-triplet, because  $y$  is not chosen as belonging to the current cluster and, therefore, either  $\ell(y, z) \neq \ell(z, w)$  or  $\ell(y, w) \neq \ell(z, w)$ . As a result,  $\{x, y, z, w\}$  must be a B-quadruple  $d$ -hitting  $(x, z)$ , which implies that (cf. Equation (5)):

$$\begin{aligned} & A_{xyzw}^{(i)} \wedge \neg P_x^{(i)} \wedge \neg P_y^{(i)} \Rightarrow \\ & \Rightarrow \bigvee_i \left( A_{xyzw}^{(i)} \wedge \neg P_x^{(i)} \wedge \neg P_y^{(i)} \right) \Leftrightarrow \\ & \Leftrightarrow \bigvee_i Q_{zw|xy}^{(i)} \Leftrightarrow Q_{zw|xy}. \end{aligned}$$

2) Two further cases may arise in this case too.

(a) Either  $x$  or  $y$  is chosen as pivot at the iteration  $i$ , along with any other object  $z$ . This situation is analogous to case 1)-(a); therefore, it is easy to see that the event  $T_{z|xy}^{(i)}$  is true in this case too.

(b) The pivots chosen at the iteration  $i$  are  $z$  and  $w$ , with  $z \neq x, z \neq y, w \neq x, w \neq y$ . As both  $x$  and  $y$  are chosen as being part of the current cluster  $C$ , then both  $\{x, z, w\}$  and  $\{x, y, w\}$  are SC-triplets. Moreover, denoting by  $l_C$  the label of  $C$ , by hypothesis it holds that  $\ell(x, y) \neq l_C = \ell(z, w)$ , which implies that both  $\{x, y, z\}$  and  $\{x, y, w\}$  are B-triplets. This is sufficient to recognize  $\{x, y, z, w\}$  as a B-quadruple  $d$ -hitting  $(x, y)$  and have a situation analogous to case 1)-(b). Thus, the event  $Q_{zw|xy}$  is true in this case too.

In conclusion, we can state that  $c_{xy}(G) > 0$  only if either  $T_{z|xy}$  occurs for any  $z$  (cases 1)-(a) and 2)-(a)) or  $Q_{zw|xy}$  occurs for any  $z, w$  (cases 1)-(b) and 2)-(b)). This proves the lemma.  $\square$

## A.3 Proof of Lemma 4

To prove Lemma 4 we first need to show:

- Some straightforward implications arising from the probability events  $T_{z|xy}^{(i)}$  and  $Q_{zw|xy}^{(i)}$ , i.e., (i) if  $T_{z|xy}^{(i)}$  occurs, then the pivots chosen at iteration  $i$  must be  $z$  along with either  $x$  or  $y$  (Lemma 11), and (ii) when  $Q_{zw|xy}^{(i)}$  happens, the pivots chosen at iteration  $i$  are  $z$  and  $w$  (Lemma 12).
- The disjointness of the events  $T_{z|xy}^{(i)}$  with respect to each other in reference to both the same iteration (Lemma 13) and different iterations (Lemma 14).
- The disjointness of the events  $Q_{zw|xy}^{(i)}$  with respect to each other in reference to both the same iteration (Lemma 15) and different iterations (Lemma 16).
- The disjointness of the events  $T_{z|xy}^{(i)}$  and  $Q_{zw|xy}^{(i)}$  with respect to one another in the same iteration (Lemma 17) as well as among different iterations (Lemma 18)

**Lemma 11** *It holds that  $T_{z|xy}^{(i)} \Rightarrow \left( P_x^{(i)} \oplus P_y^{(i)} \right) \wedge P_z^{(i)}$ .*

PROOF. By definition,  $T_{z|xy}^{(i)} \Leftrightarrow A_{xyz}^{(i)} \wedge \neg \left( P_x^{(i)} \wedge P_y^{(i)} \right)$  and  $A_{xyz}^{(i)} \Rightarrow \left( P_x^{(i)} \wedge P_y^{(i)} \right) \oplus \left( P_x^{(i)} \wedge P_z^{(i)} \right) \oplus \left( P_y^{(i)} \wedge P_z^{(i)} \right)$ . Thus,  $T_{z|xy}^{(i)} \Rightarrow \left( \left( P_x^{(i)} \wedge P_y^{(i)} \right) \oplus \left( P_x^{(i)} \wedge P_z^{(i)} \right) \oplus \left( P_y^{(i)} \wedge P_z^{(i)} \right) \right) \wedge \neg \left( P_x^{(i)} \wedge P_y^{(i)} \right) \Leftrightarrow \left( P_x^{(i)} \wedge P_z^{(i)} \right) \oplus \left( P_y^{(i)} \wedge P_z^{(i)} \right) \Leftrightarrow \left( P_x^{(i)} \oplus P_y^{(i)} \right) \wedge P_z^{(i)}$ .  $\square$

**Lemma 12** *It holds that  $Q_{zw|xy}^{(i)} \Rightarrow P_z^{(i)} \wedge P_w^{(i)}$ .*

PROOF. By definition,  $A_{xyzw}^{(i)}$  implies that the pivots chosen at iteration  $i$  correspond to one of the (six) unordered pairs that may be defined over the set  $\{x, y, z, w\}$ . However, as  $Q_{zw|xy}^{(i)} \Leftrightarrow A_{xyzw}^{(i)} \wedge \neg P_x^{(i)} \wedge \neg P_y^{(i)}$ , the pairs containing either  $x$  or  $y$  clearly make  $Q_{zw|xy}^{(i)}$  false. Thus, the only remaining choice is the pair  $(z, w)$ , which leads to the event  $P_z^{(i)} \wedge P_w^{(i)}$ .  $\square$

**Lemma 13** *It holds that  $T_{z|xy}^{(i)} \Rightarrow \neg T_{z'|xy}^{(i)}$ , for all  $z' \neq z$ .*

PROOF. According to Lemma 11,  $T_{z|xy}^{(i)}$  implies that the pair of pivots chosen in  $i$  is either  $(x, z)$  or  $(y, z)$ , whereas  $T_{z'|xy}^{(i)}$  implies one among  $(x, z')$  and  $(y, z')$ . These two situations are clearly conflicting as  $z' \neq z$ .  $\square$

**Lemma 14** *It holds that  $T_{z|xy}^{(i)} \Rightarrow \neg T_{z'|xy}^{(j)}$ , for all  $j \neq i$ , and for all  $z'$ .*

PROOF. The Chromatic Balls algorithm always removes the pivots from the set of objects available in the next iterations. Thus, if  $T_{z|xy}^{(i)}$  occurs, either  $x$  or  $y$  are no longer available for any next iteration  $j$ , as Lemma 11 states that one of them must be chosen as pivot; this clearly implies that  $T_{z'|xy}^{(j)}$  cannot be true in any iteration  $j \neq i$ , even for  $z' \neq z$ .  $\square$

**Lemma 15** *It holds that  $Q_{zw|xy}^{(i)} \Rightarrow \neg Q_{z'w'|xy}^{(i)}$ , for all  $(z', w') \neq (z, w)$ .<sup>2</sup>*

PROOF. According to Lemma 12, it holds that  $Q_{zw|xy}^{(i)} \Rightarrow P_z^{(i)} \wedge P_w^{(i)}$  and  $Q_{z'w'|xy}^{(i)} \Rightarrow P_{z'}^{(i)} \wedge P_{w'}^{(i)}$ ; but,  $P_z^{(i)} \wedge P_w^{(i)}$  and  $P_{z'}^{(i)} \wedge P_{w'}^{(i)}$  are mutually exclusive as  $(z', w') \neq (z, w)$  by hypothesis.  $\square$

**Lemma 16** *It holds that  $Q_{zw|xy}^{(i)} \Rightarrow \neg Q_{z'w'|xy}^{(j)}$ , for all  $j \neq i$ , and for all  $z', w'$ .*

PROOF. The proof is similar to Lemma 14. By definition, any event  $Q_{zw|xy}^{(i)}$  involves a B-quadruple  $\{x, y, z, w\} \in \mathcal{Q}_{xy}^d$ , thus implying that either  $\{x, z, w\}$  or  $\{y, z, w\}$  is a SC-triplet. This, along with the fact that, given  $Q_{zw|xy}^{(i)}$ , the pivots chosen are necessarily  $z$  and  $w$  (Lemma 12), is sufficient for the Chromatic Balls algorithm to put either  $x$  or  $y$  in the cluster being formed at iteration  $i$  and, therefore, make it/them not available in any next iteration  $j$ .  $\square$

**Lemma 17** *It holds that  $T_{z|xy}^{(i)} \Rightarrow \neg Q_{z'w'|xy}^{(i)}$ , for all  $z', w'$  and  $Q_{zw|xy}^{(i)} \Rightarrow \neg T_{z'|xy}^{(i)}$ , for all  $z'$ .*

PROOF. According to Lemma 11, it holds that  $T_{z|xy}^{(i)} \Rightarrow P_x^{(i)} \oplus P_y^{(i)}$ ; this contradicts  $Q_{z'w'|xy}^{(i)} \Rightarrow \neg P_x^{(i)} \wedge \neg P_y^{(i)} \Leftrightarrow \neg (P_x^{(i)} \vee P_y^{(i)})$ , which holds by definition. Thus,  $T_{z|xy}^{(i)}$  and  $Q_{z'w'|xy}^{(i)}$  are mutually exclusive.  $\square$

**Lemma 18** *It holds that  $T_{z|xy}^{(i)} \Rightarrow \neg Q_{z'w'|xy}^{(j)}$ , for all  $j \neq i$ , for all  $z', w'$  and  $Q_{zw|xy}^{(i)} \Rightarrow \neg T_{z'|xy}^{(j)}$ , for all  $j \neq i$ , and for all  $z'$ .*

PROOF. If  $T_{z|xy}^{(i)}$  happens, either  $x$  or  $y$  is chosen as pivot (Lemma 11) and, therefore, no longer available for making any  $Q_{z'w'|xy}^{(j)}$  true in any next iteration  $j$ . On the other hand, if  $Q_{zw|xy}^{(i)}$  happens, according to the same reasoning explained in Lemma 16, either  $x$  or  $y$  is put in the cluster being formed at iteration  $i$  and, therefore, not available for  $T_{z'|xy}^{(j)}$  to be true in any next  $j$ .  $\square$

Given the results shown in Lemma 11-18, we can now prove Lemma 4.

**Lemma 4** *The events within the collection  $\Omega_{xy}$  are disjoint.*

PROOF. As  $\Omega_{xy} = \{T_{z|xy} \mid \{x, y, z\} \in \mathcal{T}_{xy}\} \cup \{Q_{zw|xy} \mid \{x, y, z, w\} \in \mathcal{Q}_{xy}^d\}$ , to prove the theorem, we need to demonstrate that 1) the events  $T_{z|xy}$  are each other disjoint, 2) the events  $Q_{zw|xy}$  are each other disjoint, and 3) the events in  $\mathcal{T}_{xy}$  are disjoint from the events  $Q_{zw|xy}$  and vice versa. We account for these three cases separately.

1) We need to prove that  $T_{z|xy} \Rightarrow \neg T_{z'|xy}$ , for all  $z' \neq z$ . Denoting by  $i$  and  $j$  two generic iterations of the

<sup>2</sup> $(z', w') \neq (z, w) \Leftrightarrow (z' \neq z \wedge z' \neq w) \vee (w' \neq z \wedge w' \neq w)$

Chromatic Balls algorithm, we note that:

$$\begin{aligned} (T_{z|xy} \Rightarrow \neg T_{z'|xy}, \forall z' \neq z) &\Leftrightarrow \\ &\Leftrightarrow \left( \bigvee_i T_{z|xy}^{(i)} \Rightarrow \neg \left( \bigvee_j T_{z'|xy}^{(j)} \right), \forall z' \neq z \right) \Leftrightarrow \\ &\Leftrightarrow \left( T_{z|xy}^{(i)} \Rightarrow \bigwedge_j \neg T_{z'|xy}^{(j)}, \forall i, \forall z' \neq z \right) \Leftrightarrow \\ &\Leftrightarrow \left( T_{z|xy}^{(i)} \Rightarrow \neg T_{z'|xy}^{(i)} \wedge \bigwedge_{j \neq i} \neg T_{z'|xy}^{(j)}, \forall i, \forall z' \neq z \right). \end{aligned}$$

The latter is true as, given any iteration  $i$ ,  $T_{z|xy}^{(i)} \Rightarrow \neg T_{z'|xy}^{(i)}$ , for all  $z' \neq z$  according to Lemma 13, while  $T_{z|xy}^{(i)} \Rightarrow \neg T_{z'|xy}^{(j)}$ , for all  $j \neq i$ , and for all  $z'$  according to Lemma 14.

2) It should be demonstrated that  $Q_{zw|xy} \Rightarrow \neg Q_{z'w'|xy}$ , for all  $(z', w') \neq (z, w)$ , which is equivalent to  $Q_{zw|xy}^{(i)} \Rightarrow \neg Q_{z'w'|xy}^{(i)} \wedge \bigwedge_{j \neq i} \neg Q_{z'w'|xy}^{(j)}$ ,  $\forall i, \forall (z', w') \neq (z, w)$ , according to a similar reasoning to the previous case. Again, the latter is true given the results derived previously, i.e., in Lemma 15 ( $Q_{zw|xy}^{(i)} \Rightarrow \neg Q_{z'w'|xy}^{(i)}$ , for all  $(z', w') \neq (z, w)$ ) and Lemma 16 ( $Q_{zw|xy}^{(i)} \Rightarrow \neg Q_{z'w'|xy}^{(j)}$ , for all  $j \neq i$ , and for all  $z', w'$ ).

3) Here, we need to derive  $T_{z|xy} \Rightarrow \neg Q_{z'w'|xy}$ , for all  $z', w'$  and  $Q_{zw|xy} \Rightarrow \neg T_{z'|xy}$ , and for all  $z'$ . The former is equivalent to  $T_{z|xy}^{(i)} \Rightarrow \neg Q_{z'w'|xy}^{(i)} \wedge \bigwedge_{j \neq i} \neg Q_{z'w'|xy}^{(j)}$ ,  $\forall i, z', w'$ , which holds according to Lemma 17 and 18. Analogously, the other statement can be rewritten as  $Q_{zw|xy}^{(i)} \Rightarrow \neg T_{z'|xy}^{(i)} \wedge \bigwedge_{j \neq i} \neg T_{z'|xy}^{(j)}$ ,  $\forall i, z'$  and proved to be true by resorting again to the same lemmas (i.e., Lemma 17 and 18).  $\square$

## A.4 Proofs of Lemmas 5–6

**Lemma 5** *For a pair  $(x, y) \in V \times V$  the following bound holds.*

$$E[c_{xy}(G)] \leq \sum_{\{x, y, z\} \in \mathcal{T}_{xy}} \Pr[A_{xyz}] + \sum_{\{x, y, z, w\} \in \mathcal{Q}_{xy}^d} \frac{1}{4} \Pr[A_{xyzw}].$$

PROOF. According to Lemma 3, any pair  $(x, y)$  pays a cost only if an event in  $\Omega_{xy}$  occurs, while Lemma 4 shows that all events in  $\Omega_{xy}$  are each other disjoint. This is sufficient for stating that  $E[c_{xy}(G)] = \sum_{\omega \in \Omega_{xy}} \Pr[\omega] c_{xy|\omega}$ , where  $c_{xy|\omega}$  denotes the cost paid by  $(x, y)$  if the event  $\omega$  happens. Clearly,  $c_{xy|\omega} \leq 1$ , as Lemma 3 gives only a necessary condition. Hence, it holds that:

$$\begin{aligned} E[c_{xy}(G)] &= \sum_{\omega \in \Omega_{xy}} \Pr[\omega] c_{xy|\omega} \leq \\ &\leq \sum_{\{x, y, z\} \in \mathcal{T}_{xy}} \Pr[T_{z|xy}] + \sum_{\{x, y, z, w\} \in \mathcal{Q}_{xy}^d} \Pr[Q_{zw|xy}] \leq \\ &\leq \sum_{\{x, y, z\} \in \mathcal{T}_{xy}} \Pr[A_{xyz}] + \sum_{\{x, y, z, w\} \in \mathcal{Q}_{xy}^d} \frac{1}{4} \Pr[A_{xyzw}], \end{aligned}$$

where the latter comes from the results shown in Lemma 1 ( $\Pr[T_z|xy] \leq \Pr[A_{xyz}]$ ) and Lemma 2 ( $\Pr[Q_{zw}|xy] \leq \frac{1}{4} \Pr[A_{xyzw}]$ ).  $\square$

To prove Lemma 6, we need to first introduce the following, additional Lemmas 19 and 20, and Corollary 4

**Lemma 19** *It holds that:*

$$\begin{aligned} & \sum_{(x,y) \in V \times V} \sum_{\{x,y,z,w\} \in \mathcal{Q}_{xy}^d} \Pr[A_{xyzw}] = \\ & = \sum_{\{x,y,z\} \in \mathcal{T}} \sum_{(x',y') \in \{x,y,z\}} \sum_{\substack{w \in V \setminus \{x,y,z\}, \\ \{x,y,z,w\} \in \mathcal{Q}_{x'y'}^d}} \frac{\Pr[A_{xyzw}]}{\tau_{xyzw}}, \end{aligned}$$

where  $\tau_{xyzw}$  denotes the number of B-triplets contained in any B-quadruple  $\{x, y, z, w\}$ .

PROOF. By definition, any B-quadruple that d-hits a pair  $(x, y)$  must contain a B-triplet that hits in turn  $(x, y)$ . Therefore, for any  $(x, y)$ , any sum over all B-quadruples within  $\mathcal{Q}_{xy}^d$  can be split into two sums, one over all B-triplets that hit  $(x, y)$  and one over all objects  $w$  that make these B-triplets B-quadruples too:

$$\begin{aligned} & \sum_{(x,y) \in V \times V} \sum_{\{x,y,z,w\} \in \mathcal{Q}_{xy}^d} \Pr[A_{xyzw}] = \\ & = \sum_{(x,y) \in V \times V} \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \sum_{\substack{w \in V \setminus \{x,y,z\}, \\ \{x,y,z,w\} \in \mathcal{Q}_{xy}^d}} \frac{\Pr[A_{xyzw}]}{\tau_{xyzw}} = \\ & = \sum_{\{x,y,z\} \in \mathcal{T}} \sum_{(x',y') \in \{x,y,z\}} \sum_{\substack{w \in V \setminus \{x,y,z\}, \\ \{x,y,z,w\} \in \mathcal{Q}_{x'y'}^d}} \frac{\Pr[A_{xyzw}]}{\tau_{xyzw}}, \end{aligned}$$

where the scaling factor  $\tau_{xyzw}$  is introduced because, in the original sum, any B-quadruple within  $\mathcal{Q}_{xy}^d$  is taken into account only once, while in the split sum it is considered as many times as the number of its B-triplets.  $\square$

**Lemma 20** *Given any B-triplet  $\{x, y, z\} \in \mathcal{T}$ , it holds that:*

$$\sum_{(x',y') \in \{x,y,z\}} \sum_{\substack{w \in V \setminus \{x,y,z\}, \\ \{x,y,z,w\} \in \mathcal{Q}_{x'y'}^d}} \frac{\Pr[A_{xyzw}]}{\tau_{xyzw}} = 3 X_{xyz} + 2 Y_{xyz},$$

where

$$X_{xyz} = \sum_{w \in W_{xyz}} \frac{\Pr[A_{xyzw}]}{\tau_{xyzw}}, \quad (12)$$

$$Y_{xyz} = Y_{xyz}^{xy} + Y_{xyz}^{xz} + Y_{xyz}^{yz}, \quad (13)$$

$$Y_{xyz}^{xy} = \sum_{w \in W_{xyz}^{xy}} \frac{\Pr[A_{xyzw}]}{\tau_{xyzw}},$$

$$Y_{xyz}^{xz} = \sum_{w \in W_{xyz}^{xz}} \frac{\Pr[A_{xyzw}]}{\tau_{xyzw}},$$

$$Y_{xyz}^{yz} = \sum_{w \in W_{xyz}^{yz}} \frac{\Pr[A_{xyzw}]}{\tau_{xyzw}}.$$

PROOF. For each  $(x, y) \in V \times V$ , the internal sum in the statement of the lemma is over all  $w$  that make  $\{x, y, z, w\}$  a B-quadruple d-hitting  $(x, y)$ . We split this sum as described next. In order to satisfy the general B-quadruple conditions, at least one SC-triplet may be contained in  $\{x, y, z, w\}$ . In principle,  $\{x, y, z, w\}$  may contain up to three SC-triplets as at most four distinct triplets can be defined over any set of four objects and at least one of them is not an SC-triplet (i.e.,  $\{x, y, z\}$ , which is a B-triplet by definition); however, the case where exactly three SC-triplets are contained in  $\{x, y, z, w\}$  cannot arise, as it is easy to verify that this would imply  $\{x, y, z\}$  to be an SC-triplet too. Thus, we can define four possible sets of objects which  $w$  should belong to in order to make  $\{x, y, z, w\}$  a B-quadruple:

1.  $W_{xyz} = \{w \in V \setminus \{x, y, z\} \text{ s.t. two among } \{x, y, w\}, \{x, z, w\}, \{y, z, w\} \text{ are SC-triplets}\}.$
2.  $W_{xyz}^{xy} = \{w \in V \setminus \{x, y, z\} \text{ s.t. } \{x, y, w\} \text{ is an SC-triplet, while } \{x, z, w\} \text{ and } \{y, z, w\} \text{ are not}\}.$
3.  $W_{xyz}^{xz} = \{w \in V \setminus \{x, y, z\} \text{ s.t. } \{x, z, w\} \text{ is an SC-triplet, while } \{x, y, w\} \text{ and } \{y, z, w\} \text{ are not}\}.$
4.  $W_{xyz}^{yz} = \{w \in V \setminus \{x, y, z\} \text{ s.t. } \{y, z, w\} \text{ is an SC-triplet, while } \{x, y, w\} \text{ and } \{x, z, w\} \text{ are not}\}.$

Now, it can easily be verified that:

- $w \in W_{xyz}$  implies that  $\{x, y, z, w\}$  d-hits all  $(x, y)$ ,  $(x, z)$ ,  $(y, z)$ .
- $w \in W_{xyz}^{xy}$  implies that  $\{x, y, z, w\}$  d-hits  $(x, z)$  and  $(y, z)$ , but not  $(x, y)$ .
- $w \in W_{xyz}^{xz}$  implies that  $\{x, y, z, w\}$  d-hits  $(x, y)$  and  $(y, z)$ , but not  $(x, z)$ .
- $w \in W_{xyz}^{yz}$  implies that  $\{x, y, z, w\}$  d-hits  $(x, y)$  and  $(x, z)$ , but not  $(y, z)$ .

Within this view, it holds that:

$$\begin{aligned} & \sum_{(x',y') \in \{x,y,z\}} \sum_{\substack{w \in V \setminus \{x,y,z\}, \\ \{x,y,z,w\} \in \mathcal{Q}_{x'y'}^d}} \frac{\Pr[A_{xyzw}]}{\tau_{xyzw}} = \\ & = 3 X_{xyz} + 2 Y_{xyz}^{xy} + 2 Y_{xyz}^{xz} + 2 Y_{xyz}^{yz} = 3 X_{xyz} + 2 Y_{xyz}. \end{aligned}$$

$\square$

Combining the results in Lemmas 19 and 6 leads to the following straightforward corollary.

**Corollary 4** *It holds that:*

$$\sum_{(x,y) \in V \times V} \sum_{\{x,y,z,w\} \in \mathcal{Q}_{xy}^d} \Pr[A_{xyzw}] = \sum_{\{x,y,z\} \in \mathcal{T}} (3 X_{xyz} + 2 Y_{xyz}).$$

**Lemma 6** *The expected cost  $E[c(G)]$  of the Chromatic Balls algorithm can be bounded as follows*

$$E[c(G)] \leq Ub(G) = \sum_{\{x,y,z\} \in \mathcal{T}} \left( 3 \Pr[A_{xyz}] + \frac{3}{4} X_{xyz} + \frac{1}{2} Y_{xyz} \right).$$

PROOF.

$$\begin{aligned}
E[c(G)] &= \sum_{(x,y) \in V \times V} E[c_{xy}(G)] \leq \\
&\leq \sum_{(x,y) \in V \times V} \left( \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \Pr[A_{xyz}] + \right. \\
&\quad \left. + \sum_{\{x,y,z,w\} \in \mathcal{Q}_{xy}^d} \frac{1}{4} \Pr[A_{xyzw}] \right) = \quad (14) \\
&= \sum_{(x,y) \in V \times V} \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \Pr[A_{xyz}] + \\
&\quad + \sum_{(x,y) \in V \times V} \sum_{\{x,y,z,w\} \in \mathcal{Q}_{xy}^d} \frac{1}{4} \Pr[A_{xyzw}] = \\
&= \sum_{\{x,y,z\} \in \mathcal{T}} \Pr[A_{xyz}] \sum_{(x',y') \in \{x,y,z\}} 1 + \\
&\quad + \frac{1}{4} \sum_{\{x,y,z\} \in \mathcal{T}} (3 X_{xyz} + 2 Y_{xyz}) = \quad (15) \\
&= \sum_{\{x,y,z\} \in \mathcal{T}} \left( 3\Pr[A_{xyz}] + \frac{3}{4}X_{xyz} + \frac{1}{2}Y_{xyz} \right) = Ub(G),
\end{aligned}$$

where (14) and (15) hold according to Lemma 5 and Corollary 4, respectively.  $\square$

## A.5 Proofs of Lemmas 8–9 and Theorem 1

**Lemma 21** For any  $(x, y) \in V \times V$ , it holds that:

$$\sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6}X_{xyz} + \frac{1}{6}Y_{xyz}^{xz} + \frac{1}{6}Y_{xyz}^{yz} \right) \leq 1.$$

PROOF. The events in the collection  $\Omega_{xy}$  are disjoint (Lemma 4), hence it holds that:

$$\Pr \left[ \bigvee_{\omega \in \Omega_{xy}} \omega \right] = \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \Pr[T_z|x,y] + \sum_{\{x,y,z,w\} \in \mathcal{Q}_{xy}^d} \Pr[A_{xyzw}] \leq 1.$$

As  $\Pr[T_z|x,y] \geq \frac{1}{2} \Pr[A_{xyz}]$  (Lemma 1) and  $\Pr[A_{xyzw}] \geq \frac{1}{6} \Pr[A_{xyzw}]$  (Lemma 2), the latter leads to the following:

$$\sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6} \Pr[A_{xyzw}] \leq 1.$$

Recalling the reasoning exploited in Lemmas 19 and 20, it results that:

$$\sum_{\{x,y,z,w\} \in \mathcal{Q}_{xy}^d} \Pr[A_{xyzw}] = \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} (X_{xyz} + Y_{xyz}^{xz} + Y_{xyz}^{yz}).$$

Thus, we have that:

$$\begin{aligned}
&\sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6} \Pr[A_{xyzw}] \leq 1 \Leftrightarrow \\
&\Leftrightarrow \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6}X_{xyz} + \frac{1}{6}Y_{xyz}^{xz} + \frac{1}{6}Y_{xyz}^{yz} \right) \leq 1. \quad \square
\end{aligned}$$

**Lemma 22** For any B-triplet  $\{x, y, z\}$ , it holds that:

$$\frac{1}{6}Y_{xyz}^{xy} \leq 1, \quad \frac{1}{6}Y_{xyz}^{xz} \leq 1, \quad \frac{1}{6}Y_{xyz}^{yz} \leq 1.$$

PROOF. For a pair  $(x', y')$ , according to Lemma 21, it holds that:

$$\sum_{\{x',y',z'\} \in \mathcal{T}_{x'y'}} \left( \frac{1}{2} \Pr[A_{x'y'z'}] + \frac{1}{6}X_{x'y'z'} + \frac{1}{6}Y_{x'y'z'}^{x'z'} + \frac{1}{6}Y_{x'y'z'}^{y'z'} \right) \leq 1,$$

and, hence:

$$\frac{1}{6}Y_{x'y'z'}^{x'z'} \leq 1, \quad \frac{1}{6}Y_{x'y'z'}^{y'z'} \leq 1, \quad \forall \{x', y', z'\} \in \mathcal{T}_{x'y'}. \quad (16)$$

By definition, any B-triplet  $\{x, y, z\}$  hits pairs that can be defined over  $\{x, y, z\}$ ; then it follows that  $\{x, y, z\} \in \mathcal{T}_{xy}$ ,  $\{x, y, z\} \in \mathcal{T}_{xz}$ ,  $\{x, y, z\} \in \mathcal{T}_{yz}$ , and, therefore, according to Equation (16):

$$\{x, y, z\} \in \mathcal{T}_{xy} \Rightarrow \frac{1}{6}Y_{xyz}^{xz} \leq 1, \quad \frac{1}{6}Y_{xyz}^{yz} \leq 1,$$

$$\{x, y, z\} \in \mathcal{T}_{xz} \Rightarrow \frac{1}{6}Y_{xyz}^{xy} \leq 1, \quad \frac{1}{6}Y_{xyz}^{xz} \leq 1,$$

$$\{x, y, z\} \in \mathcal{T}_{yz} \Rightarrow \frac{1}{6}Y_{xyz}^{xy} \leq 1, \quad \frac{1}{6}Y_{xyz}^{xz} \leq 1. \quad \square$$

**Lemma 8** For any pair  $(x, y) \in V \times V$ , it holds that:

$$\sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \frac{1}{1+|\mathcal{T}_{xy}|} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6}X_{xyz} + \frac{1}{6}Y_{xyz} \right) \leq 1.$$

PROOF. As  $Y_{xyz} = Y_{xyz}^{xy} + Y_{xyz}^{xz} + Y_{xyz}^{yz}$  (Equations (12) and (13)), it holds that:

$$\begin{aligned}
&\sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6}X_{xyz} + \frac{1}{6}Y_{xyz} \right) = \\
&= \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6}X_{xyz} + \frac{1}{6}Y_{xyz}^{xz} + \frac{1}{6}Y_{xyz}^{yz} \right) + \\
&\quad \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \frac{1}{6}Y_{xyz}^{xy}.
\end{aligned}$$

The first sum in the latter expression is  $\leq 1$  according to Lemma 21. As far as the second sum, Lemma 22 states that  $\frac{1}{6}Y_{xyz}^{xy} \leq 1$ ; this implies that  $\sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \frac{1}{6}Y_{xyz}^{xy} \leq |\mathcal{T}_{xy}|$ . Hence:

$$\begin{aligned}
&\sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6}X_{xyz} + \frac{1}{6}Y_{xyz} \right) \leq 1 + |\mathcal{T}_{xy}| \Rightarrow \\
&\Rightarrow \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \frac{1}{1+|\mathcal{T}_{xy}|} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6}X_{xyz} + \frac{1}{6}Y_{xyz} \right) \leq 1. \quad \square
\end{aligned}$$

**Lemma 9** The cost  $c^*(G)$  of the optimum solution on any input instance  $G$  is lower bounded as follows:

$$\begin{aligned}
c^*(G) &\geq Lb(G) = \\
&= \sum_{\{x,y,z\} \in \mathcal{T}} \frac{1}{1+t_{max}} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6}X_{xyz} + \frac{1}{6}Y_{xyz} \right),
\end{aligned}$$

where  $t_{max} = \max_{(x,y) \in V \times V} |\mathcal{T}_{xy}|$  is the maximum number of B-triplets that hit a pair of objects.

PROOF. According to Lemma 8, it holds that (for all  $(x, y) \in V \times V$ ):

$$\begin{aligned} & \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \frac{1}{1+t_{max}} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6} X_{xyz} + \frac{1}{6} Y_{xyz} \right) \leq \\ & \leq \sum_{\{x,y,z\} \in \mathcal{T}_{xy}} \frac{1}{1+|\mathcal{T}_{xy}|} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6} X_{xyz} + \frac{1}{6} Y_{xyz} \right) \leq 1. \end{aligned}$$

Thus, we note that, setting

$$\alpha_{xyz} = \frac{1}{1+t_{max}} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6} X_{xyz} + \frac{1}{6} Y_{xyz} \right), \quad \forall \{x, y, z\} \in \mathcal{T}.$$

the condition about fractionally assigning each pair of objects within  $V \times V$  to the B-triplets in  $\mathcal{T}$  stated by Lemma 7 is satisfied. Thus, Lemma 7 can be applied here to derive the following:

$$\begin{aligned} c^*(G) & \geq \sum_{\{x,y,z\} \in \mathcal{T}} \alpha_{xyz} = \\ & = \sum_{\{x,y,z\} \in \mathcal{T}} \frac{1}{1+t_{max}} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6} X_{xyz} + \frac{1}{6} Y_{xyz} \right) = Lb(G). \end{aligned}$$

□

**Theorem 1** *The approximation ratio of the Chromatic Balls algorithm on any input instance  $G$  is*

$$r(G) = \frac{\mathbb{E}[c(G)]}{c^*(G)} \leq 6 (1 + t_{max}).$$

PROOF. Given the expressions for  $Ub(G)$  and  $Lb(G)$  derived in Lemmas 6 and 9, respectively, Equation (8) becomes:

$$\begin{aligned} \frac{\mathbb{E}[c(G)]}{c^*(G)} & \leq \frac{Ub(G)}{Lb(G)} = \\ & = \frac{\sum_{\{x,y,z\} \in \mathcal{T}} \left( 3 \Pr[A_{xyz}] + \frac{3}{4} X_{xyz} + \frac{1}{2} Y_{xyz} \right)}{\sum_{\{x,y,z\} \in \mathcal{T}} \frac{1}{1+t_{max}} \left( \frac{1}{2} \Pr[A_{xyz}] + \frac{1}{6} X_{xyz} + \frac{1}{6} Y_{xyz} \right)} \leq \\ & \leq 6 (1 + t_{max}). \end{aligned}$$

□

**Corollary 1** *The approximation ratio of the Chromatic Balls algorithm on any input instance  $G$  is*

$$r(G) \leq 6 (2 D_{max} - 1).$$

where  $D_{max} = \max_{x \in V} |\{y \mid y \in V \wedge \ell(x, y) \neq l_0\}|$ .

PROOF. By definition, at least two objects within any B-triplet must have a label other than  $l_0$ . Thus, the number of B-triplets hitting any pair  $(x, y)$  is upper bounded by the number of neighbors of  $x$  plus the neighbors of  $y$  minus 2, which is clearly  $\leq 2 \Delta - 2$ . This leads to  $t_{max} \leq 2 D_{max} - 2 \Rightarrow 6 (1 + t_{max}) \leq 6 (2 D_{max} - 1)$ , which proves the corollary. □