FPTAS for #BIS with Degree Bounds on One Side

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ABSTRACT

Counting the number of independent sets for a bipartite graph (#BIS) plays a crucial role in the study of approximate counting. It has been conjectured that there is no fully polynomial-time (randomized) approximation scheme (FP-TAS/FPRAS) for #BIS, and it was proved that the problem for instances with a maximum degree of 6 is already as hard as the general problem. In this paper, we obtain a surprising tractability result for a family of #BIS instances. We design a very simple deterministic fully polynomial-time approximation scheme (FPTAS) for #BIS when the maximum degree for one side is no larger than 5. There is no restriction for the degrees on the other side, which do not even have to be bounded by a constant. Previously, FPTAS was only known for instances with a maximum degree of 5 for both sides.

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General Terms

Algorithms, Theory

Keywords

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1. INTRODUCTION

Counting the number of independent sets in a bipartite graph (#BIS) is arguably the most important open question in the study of approximation algorithms for counting problems, which plays a similar role as the unique game for optimization problems, or the PPAD class for fixed points and Nash equilibria. We do not know if it admits a fully polynomial-time (randomized) approximation scheme (FP-TAS/FPRAS), and we do not know if it is as hard as counting the satisfying assignments for a satisfaction problem (#SAT) either. It is conjectured to be of intermediate complexity [6]. Similar to unique game, the approximability of #BIS is important not only because it is an interesting problem on its own, but mainly due to the fact that many other counting problems are proved to have the same complexity as #BIS. It is a complete problem for a family of logically defined problems called $\#RH\Pi_1$ as a subfamily of #P [6]. With the help of this intermediate class, a number of complete classifications for the approximability for various families of problems have been proved, such as the Boolean #CSP problems [7, 10, 3].

Without restricting input graphs to be bipartite, the approximability for counting the number of independent sets (#IS) is well understood. For general graphs, approximately counting the number of independent sets is as hard as finding the maximum independent set, which is NP-hard. This reduction was one of the very first proofs for inapproximability for counting problems. The hard instances used in the reduction have very large degrees, and as a result later research has been mainly focused on sparse graphs, such as graphs with a maximum degree bound. An FPRAS based on the Markov chain Monte Carlo (MCMC) method was obtained when the maximum degree is 3 in [8] and then 4 in [23]. Later, a deterministic FPTAS based on the correlation decay technique was obtained for graphs with a maximum degree of 5 by Weitz [30]. On the inapproximability side, it was proved that the problem is NP-hard as long as we allow the maximum degree to be 25 [9]. The hardness bound was eventually reduced to 6 and thus closed the gap

However, the approximability for #BIS is much more challenging. We do not know any NP-hardness result even if we do not have a degree bound. The previous proof for general graphs does not work because finding a maximum independent set for bipartite graph is equivalent to finding a maximum matching (Konig's theorem), which is polynomial time solvable rather than being NP-hard. The main reason to make #BIS extremely important in the study of approx-

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imate counting is that a large number of other problems are proved to have the same complexity as #BIS (#BISequivalent) or at least as hard as #BIS (#BIS-hard) under approximation-preserving reduction (AP-reduction) [6]. Examples include combinatorial counting problems such as #Downsets (counting the number of downsets of a partial order system), #Bipartite-q-COL, #Bipartite-MAX-IS (all in [6]) and #Stable-Matching [5], logical problems such as #1P1NSAT and #IM [6], problems from statistical physics such as computing the partition problems for ferromagnetic Ising model with mixed external fields [13] and Potts system [14], and many other counting problems. One recent interesting result on #BIS itself indicates that #BIS with maximum degree 6 is already as hard as general #BIS [4]. This restricted version of #BIS is more useful in some reductions and the new result has been used to prove #BIShardness for other problems such as ferromagnetic two-spin systems with a uniform external field [21]. Moreover, it was shown that if #BIS does not admit an FPRAS, then there is an infinite approximation hierarchy even within #BIS [1].

The main reason to make #BIS flexible in these reductions is indeed due to its bipartite structure, on which the vertices from two sides can encode (or be encoded by) two different objects for other problems. For example, a hypergraph can be represented as a bipartite graph (known as its incidence graph), with the left side being the vertex set and the right side being the edge set. In this new bipartite graph, the degrees on the left side are the same as the degrees in the hypergraph, while the degrees of the right side are sizes of hyperedges in the hypergraph. This nature makes it suitable to study #BIS with different degree constraints on two sides. For example, #Semi-regular-BIS studied in [14] has one side regularity requirement.

On the algorithmic side, it was shown in [9, 24] that any local MCMC algorithm that uses subsets of vertices as state space, mixes slowly even on a bipartite graph with a maximum degree of 6. More recently, an interesting attempted Markov chain by Ge and Stefankovic [11], which uses subsets of edges as state space and differs from previous MCMC methods, was also shown to mix slowly in [12]. Prior to our work, the best known FPRAS or FPTAS for #BIS was the same as that for #IS for graphs with a maximum degree of 5. There was no algorithmic evidence to distinguish #BIS from #IS.

Our Results

Our main result is an FPTAS for #BIS when the maximum degree for one side is no larger than 5. There is no restriction for the degrees on the other side, which do not even have to be bounded by a constant. Assuming that there is no FPTAS or FPRAS for general #BIS, our result is of the best possible in the sense that, if we allow degrees of 6 on both sides, the problem is already #BIS-hard. Our FPTAS can also be viewed as the first algorithmic evidence to distinguish #BIS from #IS.

Our algorithm is almost identical to Weitz's algorithm for general #IS with a maximum degree of 5, and the main technique is also correlation decay. We elaborate a bit on the ideas. Due to a standard argument, computing the number of independent sets is reduced to computing the marginal probability of a vertex to be chosen, if one samples an independent set uniformly at random from all possible independent sets of the input graph. Then, the main idea is

to estimate these marginal probabilities directly rather than through sampling, which is made possible by the remarkable self-avoiding walk (SAW) tree introduced by Weitz in [30]. For efficiency of computation, the marginal distribution of a vertex is estimated using only a local neighborhood around a vertex. To justify the precision of the estimation, we show that far-away vertices have little influence on the marginal distribution. This is done by analyzing the decay rate of correlation between two vertices in terms of their distance. In [30], it is proved that when the degree of each vertex is at most 5, this decay rate is exponentially small in the depth of the SAW tree. However, the same analysis does not apply to our case as the degrees of one side can be arbitrarily large. To overcome this, our main idea is to combine two recursion steps of the SAW for #BIS into one, and work with this two-layer recursion instead. As it turns out, it has the same effect as treating one side of vertices as variables, while the other side of vertices as constraints. Then we ensure that the degrees in the first layer, which are the variables' degree, are always no more than 5. The key is to formalize an observation that the larger the second layer degree (the constraint's degree), the faster the correlation decays. Such analysis is only possible for #BIS rather than general #IS.

Such a two-layer type recursion is similar to that for monotone CNF and hypergraph matching in [20]. As we have seen there, the analysis for these two-layer recursions is usually much more challenging and complicated. One additional complication here is due to the fact that the degrees for the other side are not even bounded by a constant. For these cases, we need to prove an even stronger notion of correlation decay called computationally efficient correlation decay as in [17, 19, 20], which says that the error decays by a super-constant factor if we go through a vertex with a super-constant degree. In order to prove the correlation decay property, we use a potential function to amortize the decay rate as in many previous works [25, 17, 18, 26, 22]. A good potential function is the key to these proofs. In this paper, the potential function is carefully constructed to not only make the decay rate less than one but also make the proof simpler. Effectively, the potential function we use in this paper makes the amortized decay rate of the two-layer recursion act as if it is a single layer. This dramatically simplifies the proof. We believe this simple idea can find applications in the analysis of other two-layer recursions.

Related work

The correlation decay based FPTAS for counting independent sets was extended to anti-ferromagnetic two-spin systems [17, 26, 18]. From a statistical physics point of view, the independent set problem is a special case of the hard-core model, where one introduces an activity parameter and counts weighted independent sets. To extend our result to weighted independent sets and anti-ferromagnetic two-spin systems in general is an interesting open question.

There are some other works that study counting problems for richer families of graphs other than a single maximum degree constraint. A beautiful direction is to replace the maximum degree constraint by the connective constant [28, 27], which can be viewed as a version of average degree. However, if one would like to apply this average degree type argument to the #BIS instances in our setting, the connective constant is unbounded since the degree of one side is unbounded. Our result also indicates that in the case of

bipartite graphs, the average degree may not be powerful enough to capture the complexity of the problem.

Such phenomena where larger degrees (the degrees of constraints) only make the problem easier, also come up in hypergraph independent sets. In particular, let d be the maximum degree, and m be the minimum edge size (which plays the role of constraint degree). As is shown in [2], if $m \geq d+2 \geq 5$, the problem of counting independent sets on such hypergraphs admits FPRAS. In contrast, if we only have maximum degree parameter d, then it only admits FPTAS when $d \leq 5$ [20].

Bipartiteness changing the complexity of a problem is also an interesting phenomenon in the study of approximate counting. Two other famous examples are graph colorings and perfect matchings. Counting the number of colorings for bipartite graphs is an important open question, which is known to be #BIS-hard, but not known to be #BIS-equivalent or not. There is an FPRAS for counting perfect matchings in a bipartite graph [15], while for general graphs it is a long-standing open question.

2. PRELIMINARY

For an undirected graph G=(V,E), a subset of vertices $I\subseteq V$ is an *independent set* of G if there is no edge between any two vertices within I. We denote I(G) as the set of independent sets of graph G, and $Z(G)\triangleq |I(G)|$. G is bipartite if there exists $U\subseteq V$ such that both $U\in I(G)$ and $U^C\in I(G)$. Hence it can be written as $G=(U\uplus U^C,E)$.

Given a graph G=(V,E), a vertex $u\in V$, a set of vertices $U\subseteq V$, we define the following:

- Removing a vertex u and its incident edges: $G-u=(V\setminus\{u\}\,,\{e\in E\mid e\text{ is not incident with }u\})\,.$
- \bullet Removing a set of vertices U and all incident edges:

$$G-U=(V\setminus U,$$

$$\{e\in E\mid e\text{ is not incident with any }u\in U\})\,.$$

We write $N_G(u)$ for an open neighborhood of a vertex u (which does not contain u), and $N_G[u]$ for a closed neighborhood of u (which includes u itself). Note that in the case of a bipartite graph $G = (U \uplus V, E)$, for every vertex $u \in U$, we have $N_G(u) \subseteq V$.

In general, we use u=0 to refer the vertex u is not chosen in an independent set, and u=1 for being chosen. With an independent set sampled uniformly at random, the probability that the vertex u is chosen is denoted by $\mathbb{P}_G(u=1)$. Similarly, $\mathbb{P}_G(u=0)$ is for the probability that the vertex u is not chosen.

As an easy observation, the number of independent sets without choosing u is Z(G-u), with those choosing u being $Z(G-N_G[u])$. Thus $\mathbb{P}_G(u=0)=\frac{Z(G-u)}{Z(G)}$, and similarly $\mathbb{P}_G(u=1)=\frac{Z(G-N_G[u])}{Z(G)}$.

3. THE ALGORITHM

The main result of this paper is the following algorithm.

Theorem 1. There is an FPTAS for counting the number of independent sets of a bipartite graph $G = (U \uplus V, E)$ with min $\{\Delta_U, \Delta_V\} \leq 5$, where Δ_U and Δ_V are the maximum degree over vertex set U and V respectively.

Without loss of generality, we assume $\Delta_U \leq \Delta_V$. Thus, we have $\Delta_U \leq 5$. We denote n = |U|, m = |V|.

3.1 Counting from Likelihood Ratios

We shall first reduce the problem of counting to computing likelihood ratios. This is a standard reduction, and was introduced as the self-reducibility structure in [16].

For vertices $u \in U$, let $R(G, u) \triangleq \frac{\mathbb{P}_G(u=1)}{\mathbb{P}_G(u=0)} = \frac{Z(G-N_G[u])}{Z(G-u)}$. Although we can similarly define R(G, v) for $v \in V$, our ultimate algorithm would only involve vertices $u \in U$ as variables.

Let u_1, u_2, \ldots, u_n be an arbitrary enumeration of vertices in U, and $G_i = G - \{u_1, \ldots, u_{i-1}\}$. In particular, $G_1 = G$, and $G_n - u_n = G - U$. Also recall that the vertex set of G - U is just V, which is an independent set of G by itself,

$$Z(G) = Z(G - u_1) + Z(G - N_G[u_1])$$

$$= Z(G_1 - u_1) \cdot (1 + R(G_1, u_1))$$

$$= (Z(G_2 - u_2) + Z(G_2 - N_{G_2}[u_2])) \cdot (1 + R(G_1, u_1))$$

$$= Z(G_2 - u_2) \cdot (1 + R(G_2, u_2)) \cdot (1 + R(G_1, u_1))$$

$$\vdots$$

$$= Z(G_n - u_n) \cdot \prod_{i=1}^{n} (1 + R(G_i, u_i))$$

$$= Z(G - U) \cdot \prod_{i=1}^{n} (1 + R(G_i, u_i))$$

$$= 2^m \prod_{i=1}^{n} (1 + R(G_i, u_i))$$

PROPOSITION 1. Provided an algorithm $R(G, u, \varepsilon)$ for estimating R(G, u) within an additive error ε , which runs in time $poly(n, 1/\varepsilon)$, and outputs \hat{R} such that $\left|\hat{R} - R(G, u)\right| \le \varepsilon$. There is an FPTAS for estimating Z(G) based on $R(G, u, \varepsilon)$.

PROOF. Let $G_i = G - \{u_1, \dots, u_{i-1}\}$. Given $0 < \varepsilon < 1$, let $\hat{R}_i \triangleq R\left(G_i, u_i, \frac{\varepsilon}{2n}\right)$ and $R_i \triangleq R\left(G_i, u_i\right)$. Consider the algorithm that returns $\hat{Z}(G) = 2^m \prod_{i=1}^n \left(1 + \hat{R}_i\right)$ as an approximation for $Z(G) = 2^m \prod_{i=1}^n (1 + R_i)$. We have

$$\frac{\left|\hat{R}_{i} - R_{i}\right|}{1 + R_{i}} \leq \left|\hat{R}_{i} - R_{i}\right| \leq \frac{\varepsilon}{2n}$$

$$\implies \left(1 - \frac{\varepsilon}{2n}\right) \leq \frac{1 + \hat{R}_{i}}{1 + R_{i}} \leq \left(1 + \frac{\varepsilon}{2n}\right).$$

Since $\frac{\hat{Z}(G)}{Z(G)} = \prod_{i=1}^{n} \frac{1+\hat{R_i}}{1+R_i}$, we have,

$$\left(1 - \frac{\varepsilon}{2n}\right)^n \le \prod_{i=1}^n \frac{1 + \hat{R}_i}{1 + R_i} = \frac{\hat{Z}(G)}{Z(G)} \le \left(1 + \frac{\varepsilon}{2n}\right)^n$$

$$\implies 1 - \varepsilon \le \frac{\hat{Z}(G)}{Z(G)} \le 1 + \varepsilon.$$

This concludes the proof. \Box

Therefore, the remaining task is to design an algorithm for $R(G, u, \varepsilon)$.

3.2 Tree Recursion from Self-Reducibility

Before implementing the algorithm required by Proposition 1, we will show a recursive relation for R(G, u) using the self-reducibility structure again, which gives an alternative derivation of Weitz's self-avoiding walk tree approach [30].

LEMMA 1. Let $d \triangleq \deg_{G}(u)$, and $N_{G}(u)$ be enumerated as $\{v_{i}\}_{i=1}^{d}$. Denote $G_{i} \triangleq (G-u) - \{v_{j}\}_{j=1}^{i-1}$, $w_{i} \triangleq \deg_{G_{i}}(v_{i})$. Let $N_{G_{i}}(v_{i})$ be enumerated as $\{u_{i,j}\}_{j=1}^{w_{i}}$, and $G_{i,j} \triangleq (G_{i} - v_{i}) - \{u_{i,k}\}_{k=1}^{j-1}$. Then

$$R(G, u) = \prod_{i=1}^{d} \left(1 + \prod_{i=1}^{w_i} \left(1 + R\left(G_{i,j}, u_{i,j}\right) \right)^{-1} \right)^{-1}.$$

We refer to d as the first-layer degree, and w_i as the second-layer degrees. If d = 0 or $w_i = 0$ for some i, we follow the convention that an empty product is 1.

The same recursion can be obtained by first constructing the self-avoiding walk tree for G from u and then combining two steps of the tree recursion at a time. Instead of explicitly constructing the whole SAW tree, we present an alternative derivation based only on a nontrivial partition scheme promised by the self-reducibility.

PROOF. Recall that $d \triangleq \deg_G(u)$, and $N_G(u)$ is enumerated as $\{v_i\}_{i=1}^d$, and $G_i - v_i = G_{i+1}$,

$$R(G, u) = \frac{\mathbb{P}_G(u = 1)}{\mathbb{P}_G(u = 0)} = \frac{\frac{Z(G - N_G[u])}{Z(G)}}{\frac{Z(G - u)}{Z(G)}} = \prod_{i=1}^d \frac{Z(G_i - v_i)}{Z(G_i)}$$

Next we use the self-reducibility structure of the problem, which gives the following partition scheme for free: $Z(G_i) = Z(G_i - v_i) + Z(G_i - N_{G_i}[v_i])$, thus

$$\begin{split} \frac{Z(G_i - v_i)}{Z(G_i)} &= \frac{Z(G_i - v_i)}{Z(G_i - v_i) + Z(G_i - N_{G_i}[v_i])} \\ &= \frac{1}{1 + \frac{Z(G_i - N_{G_i}[v_i])}{Z(G_i - v_i)}} \\ &= \frac{1}{1 + R(G_i, v_i)}. \end{split}$$

Similarly one could use self-reducibility again and show that $R(G_i, v_i) = \prod_{j=1}^{w_i} (1 + R(G_{i,j}, u_{i,j}))^{-1}$. Substituting these $R(G_i, v_i)$ into the above recursion, we conclude the proof. \square

It is worth noting that $\mathbb{P}_{G_i}\left(v_i=0\right)=\frac{Z(G_i-v_i)}{Z(G_i)}$. As an intuition, if $\mathbb{P}_{G_i}\left(v_i=0\right)\approx 1$, namely $Z(G_i-v_i)\approx Z(G_i)$, then one could safely ignore the vertex v_i and still get a good approximation. We will see a more quantitative version of this fact, and in particular how it relates to the one-sided maximum degree in Claim 4.

Since $0 < \frac{1}{\prod_{j=1}^{w_i} (1+R(G_{i,j},u_{i,j}))} \le 1$, we can get the following bound for R(G,u) from the recursion, which will be useful in the analysis:

Lemma 2.

$$2^{-\deg_G(u)} \le R(G, u) \le 1.$$

We can further expand $R(G_{i,j}, u_{i,j})$ s by the above recursion and get a tree recursion for R(G, u). Since $\Delta_U \leq 5$, except for the root of the recursion, we always have $d = \deg_{G_{i,j}}(u_{i,j}) \leq \Delta_U - 1 \leq 4$ for the first-layer degree. In these cases, the above bounds are $\frac{1}{16} \leq R(G, u) \leq 1$.

3.3 Computation Tree

Now we are ready to implement the algorithm as required by Proposition 1. We recursively define R(G,u,L) as follows. For base case $L=0,\ R(G,u,L)=2^{-\deg_G(u)}$. For $L>0,\$ let $L_i'=\max\big(0,L-\left\lceil\log_{45}(w_i+1)\right\rceil\big),$ then

$$R(G, u, L) = \prod_{i=1}^{d} \left(1 + \prod_{j=1}^{w_i} \left(1 + R\left(G_{i,j}, u_{i,j}, L'_i\right)\right)^{-1} \right)^{-1}.$$

The recursion depth L is used to control the accuracy of the estimation, and plays the same role as ε referred in Proposition 1. After one step of recursion, L is subtracted by $\lceil \log_{45}(w_i+1) \rceil$ rather than 1, which is known as M-based depth introduced in [17], with M=45 in our case.

As an intuition, the recursion depth L can be thought of as a computational budget, in which we replace every node with a branching degree greater than 45 with a 45-ary branching subtree. Then, it is clear that the size of this branching computation tree up to depth L is at most $O((45d)^L) = O(180^L)$, and for second-to-base-case nodes (that is, nodes with $0 < L \le \lceil \log_{45} (w+1) \rceil$) they involve at most O(n) extra base cases, so the running time for the algorithm to compute R(G, u, L) is $O(n180^L)$.

By definition, our estimation R(G, u, L) has the same bounds as R(G, u) in Lemma 2.

Lemma 3.

$$2^{-\deg_G(u)} < R(G, u, L) < 1.$$

Formally we have the following key lemma, for which the proof is laid out in Section 4.

LEMMA 4 (CORRELATION DECAY). Let $\alpha = 0.9616$. If $G = (U \uplus V, E)$ is a bipartite graph with $\Delta_U \leq 5$, then for any $u \in U$,

$$|R(G, u, L) - R(G, u)| \le O(\alpha^L). \tag{1}$$

With this lemma, it is easy to estimate R(G,u) by R(G,u,L) with an additive error of ϵ by choosing $L=O(\log \frac{1}{\epsilon})$. Then combined with Proposition 1, we get the proof for Theorem 1

4. ANALYSIS AND CORRELATION DECAY

In this section, we establish the key correlation decay Lemma 4. To do that, a natural approach is to use induction, and show that the error decreases by a constant factor along each recursion step. Unfortunately, this step-wise decay is not true in our case. Instead, we perform an amortized analysis on the decay rate by a potential function, and show that step-wise decay is recovered on the new domain under the potential function.

In Section 4.1, we outline the induction and give a derivation of the *amortized decay rate*. We show that it suffices to bound the amortized decay rates as in Claim 2 and 3, which are proved in Section 4.2. In particular, we show how our choice of the potential function simplifies the amortized decay rates and the proof.

4.1 Amortized Decay Rates

We use $\varphi(x) = \ln(\ln(1+x))$ to map the values R(G, u, L) and R(G, u) into a new domain, and prove the following:

$$|\varphi \circ R(G, u, L) - \varphi \circ R(G, u)| \le 12\alpha^{L}.$$
 (2)

The choice of this potential function will become clear in the next subsection.

Claim 1. The condition (2) implies (1).

PROOF. Note that $\varphi(x)$ is an increasing function. Let $R=R(G,u), \hat{R}=R(G,u,L)$, recall the bounds from Lemma 2 and 3, we have

$$\varphi(\frac{1}{2^5}) \le \varphi(R), \varphi(\hat{R}) \le \varphi(1).$$

As a result, by Mean Value Theorem, $\exists \bar{y}: \varphi(\frac{1}{32}) \leq \bar{y} \leq \varphi(1)$ such that

$$\begin{aligned} \left| \hat{R} - R \right| &= \frac{\mathrm{d}\varphi^{-1}(y)}{\mathrm{d}y} \Big|_{y = \bar{y}} \cdot \left| \varphi(\hat{R}) - \varphi(R) \right| \\ &\stackrel{(\clubsuit)}{\leq} 2\ln(2) \cdot 12\alpha^L = 24\ln(2) \cdot \alpha^L. \end{aligned}$$

where (\clubsuit) follows from the fact

$$\frac{\mathrm{d}\varphi^{-1}(y)}{\mathrm{d}y} = \left(1 + \varphi^{-1}(y)\right) \ln\left(1 + \varphi^{-1}(y)\right) \le 2 \ln 2,$$

together with condition (2). This completes the proof. \Box

Since the case d=5 is applied only once at the root, we first assume that $1\leq d\leq 4$ and show the following:

$$|\varphi \circ R(G, u, L) - \varphi \circ R(G, u)| \le 4\alpha^L.$$
 (3)

We prove it by induction on L. Let $R=R(G,u), \hat{R}=R(G,u,L),$ For the base case L=0, we have

$$\left| \varphi(\hat{R}) - \varphi(R) \right| \le \varphi(1) - \varphi(\frac{1}{32}) < 4.$$

Supposing the induction hypothesis holds for L < l, we prove that it also holds for L = l. If u is an isolated vertex, i.e. d = 0, $\hat{R} = R = 1$. Now we assume that d > 0. Let

$$h(\mathbf{x}) = \prod_{i=1}^{d} \left(1 + \prod_{i=1}^{w_i} (1 + x_{i,j})^{-1} \right)^{-1},$$

which is the analytic version of the recursion. We also write $h = h(\mathbf{x})$ for short.

Let \mathbf{v} be the accurate vector with

$$y_{i,j} = \varphi \circ R(G_{i,j}, u_{i,j}),$$

and $\hat{\mathbf{y}}$ be the estimated vector with

$$\hat{y}_{i,j} = \varphi \circ R\left(G_{i,j}, \ u_{i,j}, \ \max(0, L - \lceil \log_M(w_i + 1) \rceil)\right).$$

Define $\mathbf{x} \triangleq \varphi^{-1}(\mathbf{y})$ for $x_{i,j} = \varphi^{-1}(y_{i,j})$, which is applying φ^{-1} entry-wise to \mathbf{y} , similarly for $\hat{\mathbf{x}} \triangleq \varphi^{-1}(\hat{\mathbf{y}})$. Then

$$\varphi \circ R(G, u, L) = \varphi \circ h(\mathbf{x}),$$

and

$$\varphi \circ R(G, u) = \varphi \circ h(\mathbf{\hat{x}}).$$

Denote

$$\Phi(x) \triangleq \frac{\mathrm{d}\varphi(x)}{\mathrm{d}x} = \frac{1}{(1+x)\ln(1+x)}.$$

Now by Mean Value Theorem, $\exists \gamma : 0 \leq \gamma \leq 1, \tilde{\mathbf{y}} = \gamma \mathbf{y} + (1 - \gamma)\hat{\mathbf{y}}$ such that, let $\tilde{\mathbf{x}} \triangleq \varphi^{-1}(\tilde{\mathbf{y}})$,

$$\varphi \circ R(G, u, L) - \varphi \circ R(G, u)$$

$$= \sum_{i,j} \frac{\partial (\varphi \circ h \circ \varphi^{-1})}{\partial y_{i,j}} \Big|_{\mathbf{y} = \tilde{\mathbf{y}}} \cdot (\hat{y}_{i,j} - y_{i,j})$$

$$= \sum_{i,j} \left(\frac{\partial h}{\partial x_{i,j}} \Big|_{\mathbf{x} = \tilde{\mathbf{x}}} \right) \frac{\Phi(h(\tilde{\mathbf{x}}))}{\Phi(\tilde{x}_{i,j})} \cdot (\hat{y}_{i,j} - y_{i,j}).$$

By induction hypothesis, we have

$$\begin{aligned} |\hat{y}_{i,j} - y_{i,j}| \leq & 4\alpha^{\max(0, L - \lceil \log_M(w_i + 1) \rceil)} \\ \leq & 4\alpha^{L - \lceil \log_M(w_i + 1) \rceil}. \end{aligned}$$

Let $\alpha = 0.9616$, $\alpha_i = \alpha^{-\lceil \log_M(w_i+1) \rceil}$, by substitution we have

$$|\varphi \circ R(G, u, L) - \varphi \circ R(G, u)| \le 4\alpha^{L} \cdot \sum_{i,j} \left| \frac{\partial h}{\partial \tilde{x}_{i,j}} \right| \frac{\Phi(h(\tilde{\mathbf{x}}))}{\Phi(\tilde{x}_{i,j})} \alpha_{i}.$$

Therefore, the key is to bound the amortized decay rate defined as

$$\kappa_d(\mathbf{x}) \triangleq \sum_{i,j} \left| \frac{\partial h}{\partial x_{i,j}} \right| \frac{\Phi(h)}{\Phi(x_{i,j})} \alpha_i$$

In particular, the following Claim 2 completes the inductive proof for (3). Then condition (3) and Claim 3 implies condition (2), and concludes the proof of Lemma 4.

Claim 2. For
$$d \leq 4$$
, and $\frac{1}{16} \leq x_{i,j} \leq 1$, $\kappa_d(\mathbf{x}) \leq 1$.

Claim 3. For
$$\frac{1}{16} \le x_{i,j} \le 1$$
, $\kappa_5(\mathbf{x}) < 3$.

4.2 Choice of Potential Functions

In this section, we establish Claim 2 and 3 and thus conclude the key lemma and main theorem. The amortized decay rate $\kappa_d(\mathbf{x})$ is a double summation over variables $x_{i,j}$ of two layers. We first show that under our choice of the potential function, where $\varphi(x) = \ln (\ln(1+x))$ and thus $\Phi(x) = \frac{\mathrm{d}\varphi(x)}{\mathrm{d}x} = \frac{1}{(1+x)\ln(1+x)}$, the double summation can be simplified into a single summation after a suitable change of variables. Let

$$s_i = (1 + \prod_{j=1}^{w_i} (1 + x_{i,j})^{-1})^{-1}.$$

Now we have $h = \prod_{i=1}^d s_i$, and $\kappa_d(\mathbf{x})$ can be rewritten as

$$\begin{split} \kappa_{d} &= \sum_{i,j} \left| \frac{\partial h}{\partial x_{i,j}} \right| \frac{\Phi(h)}{\Phi(x_{i,j})} \alpha_{i} \\ &= \Phi(h) h \cdot \sum_{i=1}^{d} \frac{\alpha_{i} \prod_{j=1}^{w_{i}} \frac{1}{1+x_{i,j}}}{1+\prod_{j=1}^{w_{i}} \frac{1}{1+x_{i,j}}} \sum_{j=1}^{w_{i}} \left| \frac{\frac{1}{1+x_{i,j}}}{\Phi(x_{i,j})} \right| \\ &= \frac{h}{(1+h) \ln(1+h)} \cdot \sum_{i=1}^{d} \frac{\alpha_{i} \prod_{j=1}^{w_{i}} \frac{1}{1+x_{i,j}}}{1+\prod_{j=1}^{w_{i}} \frac{1}{1+x_{i,j}}} \sum_{j=1}^{w_{i}} \ln(1+x_{i,j}) \\ &= \frac{h}{(1+h) \ln(1+h)} \cdot \sum_{i=1}^{d} \frac{\alpha_{i} \prod_{j=1}^{w_{i}} \frac{1}{1+x_{i,j}}}{1+\prod_{j=1}^{w_{i}} \frac{1}{1+x_{i,j}}} \ln \left(\prod_{j=1}^{w_{i}} (1+x_{i,j}) \right) \\ &= \frac{h}{(1+h) \ln(1+h)} \cdot \sum_{i=1}^{d} \alpha_{i} (1-s_{i}) \ln \frac{s_{i}}{1-s_{i}}. \end{split}$$

Therefore, this specific potential function collapses the twolayer decay rate into a single layer one, which only depends on $\{s_i\}$. As a remark, s_i has a combinatorial meaning back in the original tree recursion, which corresponds exactly to an estimate of \mathbb{P}_{G_i} $(v_i = 0)$. In the following, we treat κ_d as a function of $\{s_i\}$ rather than $\{x_{i,j}\}$, which significantly simplifies the proof.

Next we convert the bounds for $x_{i,j}$ to bounds for s_i . Note that lowerbounding s_i here is essentially by giving a lowerbound of \mathbb{P}_{G_i} ($v_i = 0$).

Claim 4.

$$\frac{17^{w_i}}{16^{w_i} + 17^{w_i}} \le s_i < 1.$$

As an intuition, this claim says that larger w_i will only make \mathbb{P}_{G_i} ($v_i = 0$) closer to 1. In other words, if we treat v_i as a constraint, larger w_i will only weaken the overall "influence" of the constraint by making it almost always satisfied, which allows the correlation to decay faster.

PROOF. Since $s_i = (1 + \prod_{j=1}^{w_i} (1 + x_{i,j})^{-1})^{-1}$, it is clear that $s_i < 1$. As $x_{i,j} \ge \frac{1}{16}$, we have

$$s_i = (1 + \prod_{j=1}^{w_i} (1 + x_{i,j})^{-1})^{-1}$$

$$\geq (1 + \prod_{j=1}^{w_i} (1 + \frac{1}{16})^{-1})^{-1}$$

$$= \frac{17^{w_i}}{16^{w_i} + 17^{w_i}}.$$

The rate κ_d also involves parameters $\alpha_i = \alpha^{-\lceil \log_M(w_i+1) \rceil}$, which are discontinuous functions in degrees w_i . To handle this, we group the variables into the following two parts:

$$I_1 = \{i : w_i < M\}, \text{ and } I_2 = \{i : w_i \ge M\}.$$

From now on we set M=45. For α_i , we have the following:

- If $i \in I_1$, we have $\alpha_i = \frac{1}{\alpha}$ being a constant;
- If $i \in I_2$, we bound them as

$$\alpha_i = \alpha^{-\lceil \log_M(w_i+1) \rceil} < \alpha^{-\log_M(w_i+1)-1}$$
.

Let $d_1 = |I_1|$ and $d_2 = |I_2|$, then $d_1 + d_2 = d$. Next we divide the summation in κ_d into two parts, for $i \in I_1$ and $i \in I_2$ respectively.

• For $i \in I_2$, recall that s_i lies in a rather narrow range $\left[\frac{17^{w_i}}{16^{w_i}+17^{w_i}},1\right]$, and $w_i \geq M=45$, we have

$$\frac{17^{w_i}}{16^{w_i} + 17^{w_i}} \ge \frac{17^{45}}{16^{45} + 17^{45}} > \frac{9}{10}$$

As we will see, for $i \in I_2$ and s_i in these ranges, κ_d is a decreasing function in s_i . As a result, those terms in κ_d corresponding to I_2 can be upperbounded by $\frac{1}{5}$.

• For $i \in I_1$, we use Jensen's inequality to show that the maximum is achieved when s_i s are all equal to the same value \hat{s} .

Finally, we will bound the decay rate by a function in a single variable \hat{s} . Here is the formal definition and the proof. We consider the symmetrized version of κ_d as

$$\hat{\kappa}_d(\hat{s}) = \frac{\hat{s}^{d_1} \cdot d_1 \cdot (1 - \hat{s}) \ln\left(\frac{\hat{s}}{1 - \hat{s}}\right)}{\alpha \left(2^{d_2} + \hat{s}^{d_1}\right) \ln(1 + 2^{-d_2} \cdot \hat{s}^{d_1})} + \frac{d_2}{5},$$

Claim 5.

$$\max_{\mathbf{s}} \kappa_d \leq \max_{\hat{\mathbf{s}}} \hat{\kappa}_d.$$

PROOF. We begin with some elementary inequalities. Let $f(x) = (1 - e^x)(x - \ln(1 - e^x)),$

$$f''(x) = -\frac{e^x \left(1 + (1 - e^x) \ln\left(\frac{e^x}{1 - e^x}\right)\right)}{1 - e^x}.$$

Since $\frac{1}{2} \le e^x \le 1$ for $x \in [-\ln 2, 0]$, we have $f''(x) \le 0$ and f(x) is concave over $x \in [-\ln 2, 0]$. Let $\hat{s} = \left(\prod_{i \in I_1} s_i\right)^{1/d_1}$, by Jensen's inquality, we have

$$\sum_{i \in I_1} (1 - s_i) \ln \frac{s_i}{1 - s_i} = \sum_{i \in I_1} f(\ln s_i)$$

$$\leq \sum_{i \in I_1} f(\ln \hat{s})$$

$$= d_1 \cdot (1 - \hat{s}) \ln \frac{\hat{s}}{1 - \hat{s}}.$$
 (4)

Let $g(h) = \ln(1+h) - \frac{h}{1+h}$, since $g'(h) = \frac{h}{(1+h)^2} \ge 0$ for $0 \le h \le 1$, we have $g(h) \ge g(0) = 0$, namely

$$\frac{h}{(1+h)\ln(1+h)} \le 1$$
, for $0 \le h \le 1$ (5)

Also note that $\frac{h}{(1+h)\ln(1+h)}$ is decreasing in h, and $2^{-d_2} \cdot \hat{s}^{d_1} \leq h$ due to $s_i \geq \frac{1}{2}$, we have

$$\begin{split} \kappa_d \leq & \frac{h}{(1+h)\ln(1+h)} \left(\sum_{i \in I_1} \alpha_i (1-s_i) \ln \frac{s_i}{1-s_i} \right. \\ & + \sum_{i \in I_2} \alpha_i (1-s_i) \ln \frac{s_i}{1-s_i} \right) \\ \leq & \frac{2^{-d_2} \cdot \hat{s}^{d_1} \cdot \alpha^{-1}}{(1+2^{-d_2} \cdot \hat{s}^{d_1}) \ln(1+2^{-d_2} \cdot \hat{s}^{d_1})} \sum_{i \in I_1} (1-s_i) \ln \frac{s_i}{1-s_i} \\ & + \sum_{i \in I_2} \alpha^{-\lceil \log_M(w_i+1) \rceil} (1-s_i) \ln \frac{s_i}{1-s_i} \\ \leq & \frac{2^{-d_2} \cdot \hat{s}^{d_1} d_1}{\alpha \left(1+2^{-d_2} \cdot \hat{s}^{d_1}\right) \ln(1+2^{-d_2} \cdot \hat{s}^{d_1})} \cdot (1-\hat{s}) \ln \left(\frac{\hat{s}}{1-\hat{s}}\right) \\ & + \sum_{i \in I_2} \alpha^{-\lceil \log_M(w_i+1) \rceil} (1-s_i) \ln \frac{s_i}{1-s_i}. \end{split}$$

Finally, it remains to show that, for $i \in I_2$,

$$\alpha^{-\lceil \log_M(w_i+1) \rceil} (1-s_i) \ln \frac{s_i}{1-s_i} \le \frac{1}{5}. \tag{6}$$

Recall that for $i \in I_2$, $w_i \ge M = 45$, and by Claim 4, we have

$$\frac{1}{1 + \left(\frac{16}{17}\right)^{w_i}} \le s_i < 1.$$

Also note that $(1-s) \ln \frac{s}{1-s}$ is decreasing in s for $\frac{9}{10} \leq s < 1$. Let $\gamma(w) = w \left(\frac{16}{17}\right)^w \ln \left(\frac{17}{16}\right) \alpha^{-\log_M(w+1)-1}$, we have

$$\alpha^{-\lceil \log_M(w_i+1)\rceil} (1-s_i) \ln \frac{s_i}{1-s_i} \le \gamma(w_i).$$

It can be verified that $\gamma(w)$ is a decreasing function in w for $w \geq 45$, and as a result we have $\gamma(w) \leq \gamma(45) < \frac{1}{5}$. Hence the relation (6) follows and we conclude the proof. \square

Now it suffices to bound the decay rate with $\hat{\kappa}_d$. As it is a real function in a single variable \hat{s} with bounded domain $\left[\frac{1}{2},1\right]$, there is a standard calculus method to find their maximum values and we only need to verify that they satisfy Claim 2 and 3. We do that on a case-by-case basis.

Proof of Claim 2. Recall that $\hat{\kappa}_d$ is single-variate in \hat{s} with $\frac{1}{2} \leq \hat{s} \leq 1$, parameterized by d_1, d_2 .

Also note that $\hat{\kappa}_d$ is increasing in both d_1, d_2 , whereas $d_1 + d_2 = d \le 4$. So it suffices to check that $\hat{\kappa}_4 < 1$ for each case

Case $d_2 = 0$:

In this case, we have

$$\hat{\kappa}_4 = \frac{4(1-\hat{s})\hat{s}^4}{\alpha\left(\hat{s}^4+1\right)\ln\left(\hat{s}^4+1\right)}\ln\left(\frac{\hat{s}}{1-\hat{s}}\right).$$

It achieves a unique maximum at $\hat{s}^* \approx 0.758669$, and thus $\hat{\kappa}_4(\hat{s}^*) < 1$.

Case $d_2 = 1$:

In this case, we have

$$\hat{\kappa}_4 = \frac{3(1-\hat{s})\hat{s}^3}{\alpha \left(\hat{s}^3 + 2\right) \ln \left(\frac{1}{2} \left(\hat{s}^3 + 2\right)\right)} \ln \left(\frac{\hat{s}}{1-\hat{s}}\right) + \frac{1}{5}.$$

It achieves a unique maximum at $\hat{s}^*\approx 0.7691,$ and thus $\hat{\kappa}_4<\frac{4}{5}+\frac{1}{5}\leq 1.$

Case $d_2 = 2$:

In this case, we have

$$\hat{\kappa}_4 = \frac{2(1-\hat{s})\hat{s}^2}{\alpha(\hat{s}^2+4)\ln\left(\frac{1}{4}(\hat{s}^2+4)\right)}\ln\left(\frac{\hat{s}}{1-\hat{s}}\right) + \frac{2}{5}.$$

It achieves a unique maximum at $\hat{s}^* \approx 0.776043$, and thus $\hat{\kappa}_4 < 0.55 + \frac{2}{5} < 1$.

Case $d_2 = 3$:

In this case, we have

$$\hat{\kappa}_4 = \frac{(1-\hat{s})\hat{s}}{\alpha(\hat{s}+8)\ln\left(\frac{\hat{s}+8}{8}\right)}\ln\left(\frac{\hat{s}}{1-\hat{s}}\right) + \frac{3}{5}.$$

It achieves a unique maximum at $\hat{s}^* \approx 0.780104$, and thus $\hat{\kappa}_4 < 0.3 + \frac{3}{5} < 1$.

Case $d_2 = 4$:

In this case, we have $\hat{\kappa}_4 \leq \frac{4}{5} < 1$. \square

Proof of Claim 3. Let $f(\hat{s}) = (1 - \hat{s}) \ln \left(\frac{\hat{s}}{1 - \hat{s}}\right)$. For $\frac{1}{2} \leq \hat{s} \leq 1$, f achieves its unique maximum at $\hat{s}^* \approx 0.782188$, and $f(\hat{s}^*) < 0.3$. Hence $\hat{\kappa}_5 < \frac{5f(\hat{s}^*)}{\alpha} + 1 < 3$. \square

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