

A 0.51-Approximation of Maximum Matching in Sublinear $n^{1.5}$ Time

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Abstract

We study the problem of estimating the size of a maximum matching in sublinear time. The problem has been studied extensively in the literature and various algorithms and lower bounds are known for it. Our result is a 0.5109-approximation algorithm with a running time of $\tilde{O}(n\sqrt{n})$.

All previous algorithms either provide only a marginal improvement (e.g., 2^{-280}) over the 0.5-approximation that arises from estimating a *maximal* matching, or have a running time that is nearly n^2 . Our approach is also arguably much simpler than other algorithms beating 0.5-approximation.

1 Introduction

Given a graph $G = (V, E)$, a matching is a set of edges with no common endpoints, and the maximum matching problem asks for finding a largest such subset. Matching is a fundamental combinatorial optimization problem, and a benchmark for new algorithmic techniques in all major computational models. It also has a wide range of applications such as ad allocation, social network recommendations, and information retrieval, among others. Given that many of these applications need to handle large volumes of data, the study of sublinear time algorithms for *estimating* the maximum matching size has received considerable attention over the past two decades.¹ A sublinear time algorithm is not allowed to read the entire graph, which would take $\Omega(n^2)$ time where $n = |V|$; instead it is provided *oracle access* to the input graph and must run in $o(n^2)$ time. There are two main oracles for graph problems considered in the literature, which we also consider in this work.

- **Adjacency list oracle.** Here, the algorithm can query (v, i) , where $v \in V$ and $i \leq n$, and the oracle reports the i -th neighbor of the vertex v in its adjacency list, or NULL if i is larger than the number of v 's neighbors.
- **Adjacency matrix oracle.** Here, the algorithm can query (u, v) , where $u, v \in V$, and the oracle reports whether there exists an edge between u and v .

Earlier results on estimating the size of the maximum matching in sublinear time mostly focused on graphs with bounded degree Δ , starting with the pioneering work of Parnas and Ron [PR07],

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¹It is impossible to *find the edges* of any constant-approximate matching in time sublinear with respect to the size of the input [PR07].

and later works of [NO08, YYI09, RTVX11, ARVX12, LRY17, Gha22]. However, for $\Delta = \Omega(n)$ these do not lead to sublinear time algorithms. Thus, later works focused on general graphs with arbitrary maximum degree and managed to obtain sublinear time algorithms for them [CKK20, KMNFT20, Beh21, BRRS23, BKS23b, BKS23a]. In particular, the state of the art results can be categorized into two regimes:

- Algorithms that run in slightly sublinear time, i.e., $n^{2-\Omega_\varepsilon(1)}$. For example, the works of [BRR23b, BKS23b] gave a $(2/3 - \varepsilon)$ -approximation algorithm that runs in time $n^{2-\Omega_\varepsilon(1)}$. Later, [BKS23a] improved the approximation factor to $1 - \varepsilon$.
- Algorithms whose approximation factor is $0.5 + \varepsilon$. The state of the art result in this category is the work of [BRRS23], whose running time is $n^{1+\varepsilon}$ for an approximation factor of $0.5 + \Omega_\varepsilon(1)$. However, the best approximation factor one can get using their trade-off is only $0.5 + 2^{-280}$.² Indeed, they mention: *We do not expect our techniques to lead to a better than, say, 0.51-approximation in $n^{2-\Omega(1)}$ time.*

Thus, all known algorithms either give only a marginal improvement over the 0.5-approximation arising from estimating the size of a *maximal* matching (which can be done even in $\tilde{O}(n)$ time [Beh21]), or have a running time that is nearly n^2 . Making a significant improvement on both fronts simultaneously has remained an elusive open question.

Our results. In this work, we show how to beat the factor 0.5 in time that is strongly sublinear. Specifically, we present an algorithm that runs in time $O(n\sqrt{n})$ and achieves an approximation factor of 0.5109 for estimating the size of the maximum matching.

It is worth noting that our algorithm is much simpler, both in terms of implementation and analysis, compared to the prior works that beat 0.5-approximation [BRRS23].

Theorem 1.1. *There exists an algorithm that, given access to the adjacency list of a graph, estimates the size of the maximum matching with a multiplicative approximation factor of 0.5109 and runs in $\tilde{O}(n\sqrt{n})$ time with high probability.*

Theorem 1.2. *There exists an algorithm that, given access to the adjacency matrix of a graph, estimates the size of the maximum matching with a multiplicative-additive approximation factor of $(0.5109, o(n))$ and runs in $\tilde{O}(n\sqrt{n})$ time with high probability.*

Moreover, our algorithm can be employed as a subroutine for Theorem 2 of [Beh23] to obtain an improved approximation ratio in the dynamic setting. More precisely, that result requires a subroutine for estimating the maximum matching size in $\tilde{O}(n\sqrt{n})$ time, for which it uses the 0.5-approximation of [Beh21]. Our algorithm can be used instead, resulting in a very slight improvement to the overall approximation guarantee for [Beh23].

We note that the framework of [BKS23a] can also be used to obtain a similar result. Their algorithm, which performs a single iteration to find a constant fraction of augmenting paths of length three on top of a maximal matching, likewise yields a better-than-2 approximate matching with $n^{2-\Omega(1)}$ running time. However, the trade-off in this approach is worse in terms of both the approximation ratio and the running time.

²More specifically, for $\varepsilon \in (0, 1/4)$, they get an algorithm with approximation factor of $0.5 + 2^{-70/\varepsilon}$ that runs in time $O(n^{1+\varepsilon})$.

Related work. On the lower bound front, Parnas and Ron [PR07] demonstrated that any algorithm getting a constant approximation of the maximum matching size needs at least $\Omega(n)$ time. More recently, the work of [BRR23b] established that any algorithm providing a $(2/3 + \Omega(1), \varepsilon n)$ -multiplicative-additive approximation requires at least $n^{6/5-o(1)}$ time. For sparse graphs, a lower bound of $\Delta^{\Omega(1/\varepsilon)}$ was shown for any εn additive approximation [BRR23a]. For dense graphs, [BRR24] showed a lower bound of $n^{2-O_\varepsilon(1)}$ for the runtime of algorithms achieving such additive approximations.

Paper organization. In Section 2, we provide an overview of the challenges encountered while designing our algorithm and the techniques used to address them. We first develop an algorithm for bipartite graphs with a multiplicative-additive error in Section 4, avoiding additional challenges that arise from general graphs, trying to obtain multiplicative error (in the adjacency list model), or working with the adjacency matrix. In Section 5, we extend our algorithm to handle general graphs. In Section 6, we demonstrate how to achieve a multiplicative approximation guarantee. Finally, in Section 7, we present a simple reduction showing that our algorithm also works in the adjacency matrix model with a multiplicative-additive error.

2 Technical Overview

In this section, we provide an overview of the techniques used in this paper to design our algorithm. We begin with the two-pass semi-streaming algorithm of Konrad and Naidu [KN21] for bipartite graphs. In the first pass, the algorithm constructs a maximal matching M . In the second pass, it constructs a maximal b -matching between vertices matched in M and those unmatched in M . More specifically, each vertex in $V(M)$ has a capacity of k , while each vertex in $V \setminus V(M)$ has a capacity of kb , where $b = 1 + \sqrt{2}$ and k is a large constant. The idea is that if M is far from maximum, the b -matching will contain many length-3 augmenting paths that can be used to augment M . This algorithm obtains a $(2 - \sqrt{2}) \approx 0.585$ -approximation.

Our goal is to develop a sublinear-time algorithm by translating this semi-streaming two-pass algorithm to the sublinear time model. When trying to do so, several challenges arise. In this section we describe them step by step, and show how to overcome them.

Challenge (1): constructing a maximal matching in sublinear time is not possible. In fact, finding all edges of any constant-factor approximation of the maximum matching is impossible in sublinear time due to [PR07]. Dynamic algorithms for maximum matching [Beh23, BKS23] use the following approach: they maintain a maximal matching M and then apply the sublinear-time **random greedy maximal matching (RGMM)** algorithm of Behnezhad [Beh21] to estimate the size of the maximal b -matching. In our setting, we cannot afford to explicitly construct M . However, we can obtain oracle access to M using the sublinear-time RGMM algorithm of [Beh21]. More specifically, we can query whether a vertex v is matched in M or not in $\tilde{O}(n)$ time. Therefore, a possible solution to address the first challenge is to design two nested oracles: the outer oracle attempts to build a maximal b -matching, whereas the inner oracle checks the status of vertices (matched or not in M) to correctly filter edges and assign capacities to each vertex.

Challenge (2): two nested oracles require $\Omega(n^2)$ time. The algorithm of [Beh21] runs in $\tilde{O}(\bar{d}(G))$ time, where $\bar{d}(G)$ denotes the average degree of the graph G . Additionally, for the outer oracle, it requires $\tilde{O}(\bar{d}(G[V(M), V \setminus V(M)]))$ time (i.e., queries to the inner oracle). Unfortunately, it is possible for both $\bar{d}(G)$ and $\bar{d}(G[V(M), V \setminus V(M)])$ to be as large as $\Omega(n)$. For example,

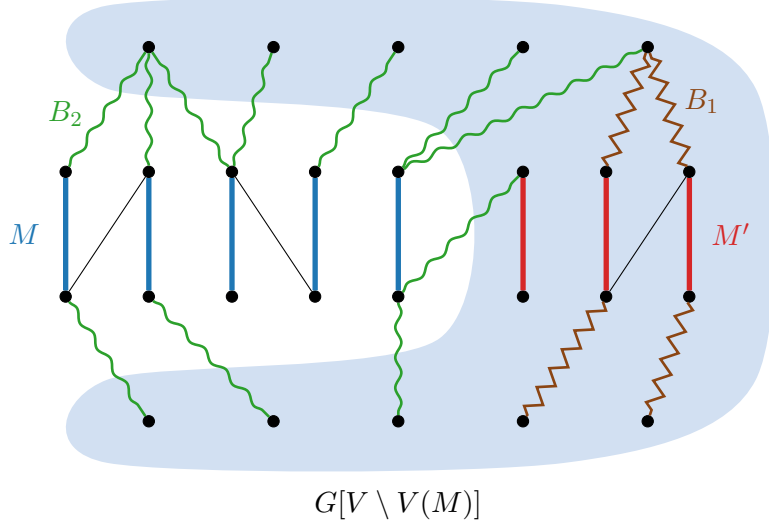


Figure 1: Our algorithm explicitly constructs a matching M (blue), which need not be maximal in G . We extend it with another matching M' (red), such that $M \cup M'$ is maximal. The highlighted (light blue) subgraph $G[V \setminus V(M)]$ has degree at most \sqrt{n} with high probability. In case 1, our algorithm augments M' using a b -matching B_1 (zigzag edges, brown). In case 2, our algorithm augments M using a b -matching B_2 (swirly edges, green).

consider a graph with a vertex set $A \cup B$, where $|A| = |B| = n/2$. The edges within A form a complete bipartite graph, while there is an $\varepsilon n/2$ -regular graph between A and B . After running the RGMM algorithm, most edges in the maximal matching belong to $G[A]$, and most vertices in B are unmatched. Consequently, we have $\bar{d}(G[V(M), V \setminus V(M)]) = \Omega(n)$.

To address this issue, we sparsify the original graph while manually constructing a matching M . In a preprocessing step, starting from an empty M , for each unmatched vertex in the graph, we sample $\Theta(\sqrt{n})$ neighbors uniformly at random. If an unmatched neighbor exists, we match the two vertices and add this edge to M . Using this preprocessing step, we show that after spending $\tilde{O}(n\sqrt{n})$ time, the induced subgraph of vertices that remain unmatched in M has a maximum degree of \sqrt{n} with high probability. Moreover, since we explicitly materialize M , we are able to check if any vertex is matched in M in $O(1)$ time, eliminating the need for costly oracle calls. Note that M need not be maximal in G ; therefore we next extend it to a maximal matching.

Let M' be a maximal matching in $G[V \setminus V(M)]$ obtained by running the sublinear time RGMM algorithm of [Beh21]. Now $M \cup M'$ is a maximal matching for G . Inspired by the two-pass semi-streaming algorithm of [KN21], we attempt to augment the maximal matching $M \cup M'$ in two possible ways (see also Figure 1):

1. Augment M' using a b -matching between $V(M')$ and $V \setminus V(M) \setminus V(M')$. The algorithm then outputs the size of the augmented matching, plus the size of the previously constructed matching M .
2. Augment M using a b -matching between $V(M)$ and $V \setminus V(M)$. The algorithm then outputs the size of the augmented matching.

The key intuition here (think of the case when $M \cup M'$ yields only a 0.5-approximation) is that either M' is sufficiently large, making $|M| + (2 - \sqrt{2}) \cdot 2|M'|$ larger than the approximation

guarantee, or M itself is large enough so that $(2 - \sqrt{2}) \cdot 2|M|$ meets our approximation guarantee (note that $(2 - \sqrt{2})$ is the approximation guarantee of [KN21]). Augmenting M using a b -matching is easier since we have explicit access to M and only need to run a single RGMM oracle to estimate the size of the b -matching. Our first estimate, which requires finding a b -matching between $V(M')$ and $V \setminus V(M) \setminus V(M')$, is more challenging since we do not have explicit access to M' . To avoid the $\Omega(n^2)$ running time of the two nested oracles, we make crucial use of the aforementioned property that the subgraph of vertices unmatched in M has low induced degree (at most \sqrt{n}); this is the reason why we only try to augment M' rather than $M \cup M'$. We will discuss this in the next paragraphs.

Challenge (3): the algorithm does not have access to the adjacency list of $G[V \setminus V(M)]$.

After the sparsification step, the average degree d of $G[V \setminus V(M)]$ is at most \sqrt{n} . Hence, if the algorithm had access to the adjacency list of $G[V \setminus V(M)]$, it could run the nested oracles in $\tilde{O}(d^2) = \tilde{O}(n)$ time by executing two RGMM algorithms: inner oracle for computing M' and outer oracle for the b -matching to augment M' . But, since the nested oracles may visit up to n vertices, and retrieving the full adjacency list of a vertex in $G[V \setminus V(M)]$ requires $\Omega(n)$ time, it seems that the overall running time of the algorithm could still be as high as $\Omega(n^2)$.

Here, we leverage two key properties of the RGMM algorithm to refine the runtime analysis. The first property is that at each step, the algorithm requires only a random neighbor of the currently visited vertex. Intuitively, if a vertex has degree $\Theta(n)$ in G , in expectation it takes $O(n/d)$ samples from the adjacency list of the original graph to encounter a vertex from $G[V \setminus V(M)]$. Thus, if all vertices in $G[V \setminus V(M)]$ had degree d , one could easily argue that the running time of the algorithm is $\tilde{O}(d^2 \cdot n/d) = \tilde{O}(n\sqrt{n})$. However, vertex degrees can vary, and for a vertex with a constant degree, we would need $\Omega(n)$ samples from the adjacency list of G to find a single neighbor in $G[V \setminus V(M)]$. To address this challenge, we utilize another property of the RGMM algorithm, recently proven by [MRTV25]. Informally, this result shows that during oracle calls for RGMM, a vertex is visited proportionally to its degree, implying that low-degree vertices are visited only a small number of times.

Challenge (4): outer oracle creates biased inner oracle queries. The final main challenge we discuss here is that the simple $\tilde{O}(n\sqrt{n})$ bound, which we informally proved in the previous paragraph, relies on the tacit assumption that the inner oracle queries generated by the outer oracle correspond to $\tilde{O}(\sqrt{n})$ uniformly random calls to the inner oracle. Indeed, the running time of the algorithm of [Beh21] is analyzed for a uniformly random query vertex; however, there may exist a vertex v in the graph for which calling the inner oracle takes significantly more than $\tilde{O}(d)$ time. Consequently, if all outer oracle calls end up querying v , the running time could be significantly worse than $\tilde{O}(n\sqrt{n})$. To overcome this issue, we use the result of [MRTV25] along with the fact that the maximum degree of $G[V \setminus V(M)]$ is $\tilde{O}(\sqrt{n})$. We show that for any vertex v , the outer oracle queries the inner oracle for v at most $\tilde{O}(\deg_{G[V \setminus V(M)]}(v)/\sqrt{n})$ times in expectation. This enables us to formally prove that the total running time of the algorithm remains at most $\tilde{O}(n\sqrt{n})$.

General graphs and the adjacency matrix model. There are additional challenges when dealing with general graphs as opposed to bipartite graphs, such as the fact that the sizes of the maximal matching and the b -matching alone are insufficient to achieve a good approximation ratio. For general graphs, our algorithm estimates the maximum matching in the union of the maximal matching and the b -matching, which requires using the $(1 - \varepsilon)$ -approximate local computation

algorithm (LCA) by [LRY17] on the subgraph formed by this union, to which we only have oracle access. We encourage readers to refer to [Section 5](#) for more details about the techniques used there.

Additionally, for more information on the extension of the algorithm that operates in the adjacency matrix model, we recommend readers to check [Section 7](#).

3 Preliminaries

Given a graph G , we use $V(G)$ to denote its set of vertices and use $E(G)$ to refer to its set of edges. For a vertex $v \in V(G)$, we use $\deg_G(v)$ to denote the degree of the vertex, i.e., the number of edges with one endpoint equal to v . We use $\Delta(G)$ to denote the maximum degree over all vertices in the graph, and $\bar{d}(G)$ to denote the average degree of the graph. Further, we use $\mu(G)$ to denote the size of the maximum matching in G .

Given a graph $G = (V, E)$, and a subset of vertices $A \subseteq V$, $G[A]$ is defined to be the induced subgraph consisting of all edges with both endpoints in A . Further, given disjoint subsets $A, B \subset V$ of vertices, $G[A, B]$ is defined to be the bipartite subgraph of G consisting of all edges between A and B .

Given a matching M in G , an *augmenting path* is a simple path starting and ending at different vertices, such that the first and the last vertices are unmatched in M , and the edges of the path alternate between not belonging to M and belonging to M .

Given a vector b of integer capacities of dimension $|V(G)|$, a *b-matching* in G is a *multi-set* F of edges in G such that each vertex $v \in V$ appears no more than b_v times as an endpoint of an edge in F .

Given a graph G and a permutation π over its edges, $\text{GMM}(G, \pi)$ is used to refer to the unique matching M obtained by the following process. We let M be initialized as empty, and consider the edges of G one by one according to the permutation π . We add an edge to the matching M if none of its endpoints are already matched in M . A random greedy maximal matching, i.e., $\text{RGMM}(G)$ is the matching obtained by picking a permutation π uniformly at random and outputting $\text{GMM}(G, \pi)$.

Proposition 3.1 ([Beh21]). *There exists an algorithm that, given adjacency list access to a graph G of average degree d , for a random vertex v and a random permutation π , determines if v is matched in $\text{GMM}(G, \pi)$ in $\tilde{O}(d)$ expected time.*

Given the problem of maximizing a function $f : D \rightarrow \mathbb{R}$ defined over a domain D , with optimal value f^* , an (α, β) -multiplicative-additive approximation of f^* is a solution $s \in D$ such that $(f^*/\alpha) - \beta \leq f(s) \leq f^*$.

4 Algorithm for Bipartite Graphs

We begin by describing our algorithm for bipartite graphs. We focus on implementing an algorithm with a multiplicative-additive approximation guarantee. Also, we assume that we have access to the adjacency list of the graph. These assumptions will help us avoid certain complications and challenges that arise when working with general graphs, the adjacency matrix model, or when trying to obtain a multiplicative approximation guarantee. To lift these assumptions, we can leverage strong tools and methods from the literature, which, with slight modifications, can be applied here. This section contains the main novelties of our approach and proofs. Our algorithm for bipartite graphs can be seen as a translation and implementation of a two-pass streaming algorithm, which we discuss in the next subsection.

4.1 Two-Pass Streaming Algorithm for Bipartite Graphs

Our starting point is the two-pass streaming algorithm which is described in [Algorithm 1](#). This algorithm, or its variations, has appeared in previous works on designing streaming or dynamic algorithms for maximum matching [[KN21](#), [BKS23](#), [Beh23](#)]. In words, the first pass of the algorithm only finds a maximal matching M . In the second pass, the algorithm finds a maximal b -matching B in $G[V(M), V \setminus V(M)]$, where $V(M)$ is the set of vertices matched by M . The capacities of vertices in $V(M)$ and in $V \setminus V(M)$ for the b -matching are k and kb , respectively. Moreover, in the second pass of the algorithm, when an edge (u, v) arrives in the stream, we add multiple copies of the edge to the subgraph B , as long as doing so does not violate the capacity constraints.

Algorithm 1: Two-pass Streaming Algorithm for Bipartite Graphs

```

1 Parameter: let  $b = 1 + \sqrt{2}$  and  $k$  be an integer larger than  $\frac{1}{b\varepsilon^3}$ .
2 First Pass:  $M \leftarrow$  maximal matching of  $G$ . ▷ Finding maximal matching
3 Second Pass: ▷ Finding  $b$ -matching
4 Let  $B = \emptyset$ .
5 for  $(u, v) \in G[V(M), \overline{V(M)}]$  where  $u \in V(M)$  do
6   while  $\deg_B(u) < k$  and  $\deg_B(v) < \lceil kb \rceil$  do
7      $B \leftarrow B \cup (u, v)$ . ▷ We allow multi edges
8 return  $(1 - 1/b) \cdot |M| + 1/(kb) \cdot |B|$ .
```

Intuitively, the algorithm tries to find length-3 augmenting paths using the b -matching that it finds in the second pass. The following lemma shows the approximation guarantee of [Algorithm 1](#).

Lemma 4.1 (Lemma 3.3 in [[BKS23](#)]). *For any $\varepsilon \in (0, 1)$, the output of [Algorithm 1](#) is a $(2 - \sqrt{2} - \varepsilon)$ -approximation for maximum matching of G .*

4.2 Sublinear Time Implementation of [Algorithm 1](#)

In this section, we demonstrate how to implement a modification of [Algorithm 1](#) in the sublinear time model, as outlined in [Section 2](#).

Sparsification: In order to be able to use two levels of recursive oracle calls, we need to sparsify the graph. We first sample $\tilde{O}(n\sqrt{n})$ edges and construct a maximal matching on the sampled edges to sparsify the induced subgraph of unmatched vertices. This sparsification step is formalized in [Algorithm 2](#). In [Lemma 4.3](#), we show that if we sample enough edges, then vertices that remain unmatched after this phase have an induced degree of \sqrt{n} . This step is very similar to the algorithm of [[CKK20](#), Appendix A] for approximating a maximal matching.

Claim 4.2. *[Algorithm 2](#) runs in $\tilde{O}(n\sqrt{n})$ time.*

Proof. [TOPROVE 1](#) □

Lemma 4.3. *With high probability, we have $\Delta(G[V \setminus V(M)]) \leq \sqrt{n}$.*

Proof. [TOPROVE 2](#) □

Algorithm 2: Sparsification of the Induced Subgraph of Unmatched Vertices

```
1 Parameter: let  $c = 2\sqrt{n} \cdot \log n$  be the sparsification parameter that controls the number
   of edges that the algorithm samples.
2  $M \leftarrow \emptyset$ 
3 for  $v \in V$  do
4   if  $v \notin V(M)$  then
5     Sample  $c$  vertices  $u_1, \dots, u_c$  from  $N(v)$ .
6     for  $i \leftarrow 1 \dots c$  do
7       if  $u_i \notin V(M)$  then
8          $M \leftarrow M \cup \{(v, u_i)\}$ 
9         break;
10 return  $M$ 
```

Augmenting M using nested oracles: Now we are ready to present our sublinear algorithm. After sparsifying the graph by finding a partially maximal matching M , we try to augment M in two different ways that we have outlined in [Section 2](#) and which are formalized in [Algorithm 3](#). See also [Figure 1](#).

For simplicity, we pretend that $kb \in \mathbb{Z}$ throughout the paper. Since k is an arbitrarily large constant, using kb instead of $\lceil kb \rceil$ leads to an arbitrarily small error in the calculations. We also note that a maximal b -matching can be viewed as maximal matching if we duplicate vertices multiple times.

First, we try to augment the matching by designing a maximal matching oracle for $G[V \setminus V(M)]$ vertices and then another oracle for finding a b -matching between the vertices newly matched using the new oracle and unmatched vertices. Let M' be the maximal matching of $G[V \setminus V(M)]$ that can be obtained by the oracle. We try to augment it with a b -matching B_1 .

However, it is also possible that $|M'|$ is small compared to $|M|$, which implies that in the previous case, the b -matching does not help to find many augmenting paths, as the size of the maximal matching that we try to augment is too small. To account for this case, the algorithm also finds a b -matching B_2 between $V(M)$ and $V \setminus V(M)$.

Note that because the algorithm finds the initial matching M explicitly, checking whether a vertex belongs to $V(M)$ or not can be done in $O(1)$ time.

Algorithm 3: Sublinear Time Algorithm for Bipartite Graphs with Access to the Adjacency List (see [Figure 1](#))

```
1 Run Algorithm 2 with  $c = 2\sqrt{n} \log n$  and let  $M$  be its output.
2 Let  $\mu_{M'}$  and  $\mu_{B_1}$  be the estimate of the size of a random greedy maximal matching  $M'$  in
    $G[V \setminus V(M)]$  and the estimate of the size of a random greedy maximal  $b$ -matching  $B_1$  in
    $G[V(M'), V \setminus V(M) \setminus V(M')]$  by running Algorithm 4. ▷ Case 1
3 Let  $\mu_1 := |M| + (1 - \frac{1}{b})\mu_{M'} + \frac{1}{kb}\mu_{B_1}$ . ▷ Case 1
4 Let  $\mu_{B_2}$  be the estimate of the size of a random greedy maximal  $b$ -matching  $B_2$  in
    $G[V(M), V \setminus V(M)]$  by running Algorithm 5. ▷ Case 2
5 Let  $\mu_2 := (1 - \frac{1}{b})|M| + \frac{1}{kb}\mu_{B_2}$ . ▷ Case 2
6 return  $\max(\mu_1, \mu_2)$ .
```

Algorithm 4: Algorithm for the First Case

- 1 Let $b = 1 + \sqrt{2}$ and k be an integer larger than $\frac{1}{b\epsilon^3}$.
 - 2 Let π be a random permutation over edges of $G[V \setminus V(M)]$ and let M' be its corresponding random greedy maximal matching.
 - 3 Let $G_1 := G[A, B]$ where $A = V(M')$ and $B = V \setminus V(M) \setminus V(M')$.
 - 4 Let G'_1 be a bipartite graph obtained from G_1 by adding k copies of vertices in A and kb copies of vertices in B . Further, if there exists an edge between $u \in A$ and $v \in B$ in G_1 , we add edges between all copies of u and v in G'_1 .
 - 5 $r \leftarrow 6 \log^3 n$.
 - 6 Run the algorithm of [Proposition 3.1](#) for r random vertices and fixed permutation π in $G[V \setminus V(M)]$ and let X_i be the indicator if the i -th vertex is matched.
 - 7 Let $X \leftarrow \sum_{i=1}^r X_i$ and $\mu_{M'} \leftarrow \frac{nX}{2r} - \frac{n}{2 \log n}$.
 - 8 Run nested oracles of [Proposition 3.1](#) for r random vertices and fixed permutation π in G'_1 and let Y_i be the indicator if the i -th vertex is matched.
 - 9 Let $Y \leftarrow \sum_{i=1}^r Y_i$ and $\mu_{B_1} \leftarrow \frac{nY}{2r} - \frac{n}{2 \log n}$.
 - 10 **return** $\mu_{M'}$ and μ_{B_1} .
-

Algorithm 5: Algorithm for the Second Case

- 1 Let $b = 1 + \sqrt{2}$ and k be an integer larger than $\frac{1}{b\epsilon^3}$.
 - 2 Let $G_2 := G[A, B]$ where $A = V(M)$ and $B = V \setminus V(M)$.
 - 3 Let G'_2 be a bipartite graph obtained from G_2 by adding k copies of vertices in A and kb copies of vertices in B . Further, if there exists an edge between $u \in A$ and $v \in B$ in G_2 , we add edges between all copies of u and v in G'_2 .
 - 4 $r \leftarrow 6 \log^3 n$.
 - 5 Run the algorithm of [Proposition 3.1](#) for r random vertices and permutations in G'_2 and let Z_i be the indicator that shows if the i -th vertex is matched.
 - 6 Let $Z \leftarrow \sum_{i=1}^r Z_i$ and $\mu_{B_2} \leftarrow \frac{nZ}{2r} - \frac{n}{2 \log n}$.
 - 7 **return** μ_{B_2} .
-

Implementation details of the algorithm: There are some technical details in the implementation of the algorithm that are not included in the pseudocode:

- **Access to the adjacency list of an induced subgraph:** Both in [Algorithm 4](#) and [Algorithm 5](#), we run the algorithm of [Proposition 3.1](#) for some induced subgraph of G (for example, line 5 of [Algorithm 5](#)). However, [Proposition 3.1](#) works with access to the adjacency list of the input graph. To address this issue, we leverage an important property of the algorithm in [Proposition 3.1](#), namely that it only needs to find a random neighbor of a given vertex at each step of its execution. Now, whenever the algorithm requires a random neighbor of vertex v in a subgraph H , it queries random neighbors in the original graph G until it finds one that belongs to H . This increases the running time of the algorithm, as it may take $\omega(1)$ time to locate a valid neighbor in H , which we will formally bound in our runtime analysis.
- **Nested oracles in line 8 of [Algorithm 4](#):** Unlike M , we do not explicitly construct the maximal matching M' in [Algorithm 4](#). Moreover, the edges of the subgraph G'_1 connect vertices matched by M' with those that remain unmatched in either M or M' . Hence, to verify whether an edge belongs to G'_1 , we need to determine whether its endpoints are matched

or unmatched in M' by accessing the algorithm of [Proposition 3.1](#). This again increases the algorithm's runtime, which we will also formally bound in our runtime analysis.

4.3 Analysis of the Approximation Ratio

The following lemma, an analogue of Observation 3.1 in [\[BKSW23\]](#), substantiates the soundness of the estimates μ_1 and μ_2 produced in [Algorithm 3](#).

Lemma 4.4. *Let M , M' , B_1 and B_2 be as in the description of [Algorithm 3](#). Then*

- $\mu(G) \geq \mu(M \cup M' \cup B_1) \geq |M| + (1 - \frac{1}{b})|M'| + \frac{1}{kb}|B_1|$,
- $\mu(G) \geq \mu(M \cup B_2) \geq (1 - \frac{1}{b})|M| + \frac{1}{kb}|B_2|$.

Proof. [TOPROVE 3](#) □

The following lemma states the $(2 - \sqrt{2})$ -approximation guarantee of the "maximal matching plus b -matching" approach obtained in prior work, for both bipartite and general graphs. We will invoke it for appropriate subgraphs of G to obtain our guarantee.

Lemma 4.5. *Let G' be a graph, M' be any maximal matching in G' , and B be a maximal b -matching in $G'[V(M'), V(G') \setminus V(M')]$ for vertex capacities k for vertices in $V(M')$ and kb for vertices in $V(G') \setminus V(M')$, where $k > \frac{1}{b\varepsilon^3}$ and $b = 1 + \sqrt{2}$. Then:*

- for bipartite G' , we have $\mu(M' \cup B) \geq (1 - \frac{1}{b})|M'| + \frac{1}{kb}|B| \geq (2 - \sqrt{2} - \varepsilon)\mu(G')$,
- for general G' , if B is a random greedy maximal b -matching, we still have $\mathbf{E}[\mu(M' \cup B)] \geq (2 - \sqrt{2} - \varepsilon)\mu(G')$.

Proof. [TOPROVE 4](#) □

The following lemma is the crux of our approximation ratio analysis.

Lemma 4.6. *In a bipartite graph G , let M , M' , B_1 and B_2 be as in the description of [Algorithm 3](#). Then*

$$\max \left[|M| + (1 - \frac{1}{b})|M'| + \frac{1}{kb}|B_1|, (1 - \frac{1}{b})|M| + \frac{1}{kb}|B_2| \right] \geq 0.5109 \cdot \mu(G).$$

Proof. [TOPROVE 5](#) □

Lemma 4.7. *Let $\max(\mu_1, \mu_2)$ be the output of [Algorithm 3](#). With high probability, it holds that*

$$0.5109 \cdot \mu(G) - o(n) \leq \max(\mu_1, \mu_2) \leq \mu(G).$$

The proof of [Lemma 4.7](#) is routine.

Proof. [TOPROVE 6](#) □

4.4 Running Time Analysis

For the analysis of the running time, we use a crucial property of random greedy maximal matching algorithm that was proved recently in [MRTV25].

Proposition 4.8 (Lemma 5.14 of [MRTV25]). *Let $Q(v)$ be the expected number of times that the oracle queries an adjacent edge of v if we start the oracle calls from a random vertex, for a random permutation over the edges of the graph G when running random greedy maximal matching. It holds that $Q(v) = \tilde{O}(\deg_G(v)/|V(G)|)$.*

Proposition 4.9 (Corollary 5.15 of [MRTV25]). *Let $T(v)$ be the expected time needed to return a random neighbor of vertex v . Then, the expected time to run the random greedy maximal matching oracle for a random vertex and a random permutation in graph G is $\sum_{v \in V(G)} \tilde{O}(T(v) \cdot \deg_G(v)/|V(G)|)$.*

Lemma 4.10. *Algorithm 4 runs in $\tilde{O}(\bar{d}(G) \cdot \sqrt{n})$ time in expectation.*

Proof. TOPROVE 7 □

Lemma 4.11. *Algorithm 5 runs in $\tilde{O}(\bar{d}(G))$ time in expectation.*

Proof. TOPROVE 8 □

Lemma 4.12. *Algorithm 3 runs in $\tilde{O}(n\sqrt{n})$ time with high probability.*

Proof. TOPROVE 9 □

Now we are ready to prove the final theorem of this section.

Theorem 4.13. *There exists an algorithm that, given access to the adjacency list of a bipartite graph, estimates the size of the maximum matching with a multiplicative-additive approximation factor of $(0.5109, o(n))$ and runs in $\tilde{O}(n\sqrt{n})$ time with high probability.*

Proof. TOPROVE 10 □

5 Algorithm for General Graphs

The following is an analogue of Lemma 4.6 for general graphs.

Lemma 5.1. *In a general graph G , let M , M' , B_1 and B_2 be as in the description of Algorithm 3. Then*

$$(1 - \varepsilon) \max [|M| + \mathbf{E}[\mu(M' \cup B_1)], \mathbf{E}[\mu(M \cup B_2)]] \geq 0.5109 \cdot \mu(G).$$

Proof. TOPROVE 11 □

In this section, we show how to extend our algorithm to work for general graphs. The main difference between bipartite and general graphs is that the estimate based on the sizes of $|M|$, $|M'|$, $|B_1|$, and $|B_2|$ is insufficient to achieve a 0.5109 approximation guarantee. In Lemma 5.1, we show that we can achieve this approximation ratio by estimating $\mu(M \cup B_2)$ and $\mu(M' \cup B_1)$. More formally, to produce μ_1 and μ_2 in Algorithm 3, we use $|M| + \mu(M' \cup B_1)$ and $\mu(M \cup B_2)$, respectively. Also, Algorithm 4 and Algorithm 5 output $\mu(M' \cup B_1)$ and $\mu(M \cup B_2)$, respectively.

In both Algorithm 4 and Algorithm 5 we have access to oracles that can return whether a vertex is matched in M' , B_1 , or B_2 for a fixed permutation π . These oracles can also be used to return the edge of the matching if the vertex is matched, which is a corollary of Proposition 3.1 since the algorithm of Proposition 3.1 can be also used to return the edge of the matching.

Lemma 5.2. *For a vertex v , there exists an algorithm that returns the edges of v in M' and B_1 in $\tilde{O}(\bar{d}(G) \cdot \sqrt{n})$ expected time.*

Proof. **TOPROVE 12** □

Lemma 5.3. *For a vertex v , there exists an algorithm that returns the edges of v in M and B_2 in $\tilde{O}(\bar{d}(G))$ expected time.*

Proof. **TOPROVE 13** □

Local computation algorithms (LCA) is a model of computation, also motivated by large data sets, in which the algorithm is not expected to produce the entire output at once. Instead, the algorithm is queried for parts of the output, and must produce a consistent and approximately optimal output. We use the following local computation algorithm (LCA) by [LRY17] to design our algorithm.

Proposition 5.4 ([LRY17]). *There exists a $(1 - \varepsilon)$ -approximate local computation algorithm for maximum matching of graph G in $\tilde{O}(\Delta(G)^{1/\varepsilon^2})$ time with access to the adjacency list of G .*

Now we prove the main technical part of this section that can be used to estimate $\mu(M' \cup B_1)$ and $\mu(M \cup B_2)$.

Lemma 5.5. *Let H be a subgraph of graph G . Suppose that H is the union of a constant number of random greedy maximal matching on different subsets of vertices. Also, we have oracle access to edges of random greedy maximal matching. We can query a vertex to obtain the matching edge of vertex v in $\tilde{O}(T)$ expected time. Moreover, the maximum degree of H is constant. Then, there exists a $(1 - \varepsilon)$ -approximate algorithm that estimates the size of the maximum matching of H in $\tilde{O}(T)$ expected time.*

Proof. **TOPROVE 14** □

Lemma 5.6. *There exists an algorithm that outputs a $(1 - \varepsilon)$ -approximate estimation of the value of $\mu(M \cup B_2)$ in $\tilde{O}(\bar{d}(G))$ expected time.*

Proof. **TOPROVE 15** □

Lemma 5.7. *There exists an algorithm that outputs a $(1 - \varepsilon)$ -approximate estimation of the value of $\mu(M' \cup B_1)$ in $\tilde{O}(\bar{d}(G) \cdot \sqrt{n})$ expected time.*

Proof. **TOPROVE 16** □

Theorem 5.8. *There exists an algorithm that, given access to the adjacency list of a graph, estimates the size of the maximum matching with a multiplicative-additive approximation factor of $(0.5109, o(n))$ and runs in $\tilde{O}(n\sqrt{n})$ time with high probability.*

Proof. **TOPROVE 17** □

6 Multiplicative Approximation

In this section, we show that we can achieve a multiplicative approximation guarantee by slightly increasing the number of samples in [Algorithm 4](#) and [Algorithm 5](#). First, we prove a simple lower bound for the size of the maximum matching of a graph based on its maximum and average degree.

Claim 6.1. *For any graph G , it holds that $\mu(G) \geq n\bar{d}(G)/(4\Delta(G))$.*

Proof. [TOPROVE 18](#) □

The goal is to obtain a multiplicative approximation guarantee of $(0.5109 - \varepsilon)\mu(G)$. It is important to note that if any of $|M|$, $|M'|$, $|B_1|$, or $|B_2|$ is not a constant fraction of the others, it can be omitted from the equation in the statement of [Lemma 4.6](#) without affecting the approximation by more than a function of ε . Thus, without loss of generality, we can assume that $|M| = \Omega(\mu(G))$, $|M'| = \Omega(\mu(G))$, $|B_1| = \Omega(\mu(G))$, and $|B_2| = \Omega(\mu(G))$. Consequently, using [Claim 6.1](#) and as an application of Chernoff bound, we can use $\tilde{O}_\varepsilon(\Delta(G)/\bar{d}(G))$ samples in [Algorithm 4](#) and [Algorithm 5](#) to obtain multiplicative estimation of $\mu_{M'}$, μ_{B_1} , and μ_{B_2} .

By [Lemma 4.10](#), [Algorithm 4](#) runs in $\tilde{O}(\bar{d}(G) \cdot \sqrt{n})$ time when we have $r = \tilde{O}(1)$ samples. By increasing the number of samples to $\tilde{O}_\varepsilon(\Delta(G)/\bar{d}(G))$, the running time of [Algorithm 4](#) increases to $\tilde{O}(\Delta(G) \cdot \sqrt{n})$. Moreover, by [Lemma 4.11](#), [Algorithm 5](#) runs in $\tilde{O}(\bar{d}(G))$ time when we have $r = \tilde{O}(1)$ samples. By increasing the number of samples to $\tilde{O}_\varepsilon(\Delta(G)/\bar{d}(G))$, the running time of [Algorithm 5](#) increases to $\tilde{O}(\Delta(G))$. Therefore, the total running time of the algorithm is within $\tilde{O}(n\sqrt{n})$.

Finally, we can obtain the degree of each vertex in the graph using binary search. Therefore, we can assume that we have access to $\Delta(G)$ and $\bar{d}(G)$ by spending $\tilde{O}(n)$ time. Thus we get:

Theorem 1.1. *There exists an algorithm that, given access to the adjacency list of a graph, estimates the size of the maximum matching with a multiplicative approximation factor of 0.5109 and runs in $\tilde{O}(n\sqrt{n})$ time with high probability.*

7 Algorithm with Access to the Adjacency Matrix

In this section, we use a simple reduction to show that with a small modification, our algorithm can be adapted to the setting where we have access to the graph's adjacency matrix. A slightly different version of this kind of reduction appeared in previous works on sublinear time algorithms for maximum matching [[Beh21](#), [BRRS23](#)].

It is important to note that obtaining a constant-factor multiplicative approximation is impossible when the algorithm only has access to the adjacency matrix of the graph. This is because if the graph is guaranteed to contain either a single edge or be completely empty, any algorithm would require $\Omega(n^2)$ adjacency matrix queries to distinguish between these two cases. Consequently, we allow the algorithm to have an additive error of $o(n)$ in addition to the multiplicative approximation ratio.

We build an auxiliary graph H with the following vertex set and edge set:

- **Vertex set:** $V(H)$ contains $n + 2$ disjoint sets of n vertices V_1, V_2 , and U_1, \dots, U_n . Each V_i is a copy of the vertices of the original graph. Each U_i contains $n \log^2 n$ vertices.
- **Edge set:** For each vertex $v \in V_1$, the i -th neighbor of v is the i -th vertex in V_1 if $(v, i) \in E(G)$, and otherwise it is the i -th vertex in V_2 . Similarly, for each vertex $v \in V_2$, the i -th neighbor of v is i -th vertex in V_2 if $(v, i) \in E(G)$, and otherwise it is the i -th vertex in V_1 .

Also, each $v \in V_2$ is connected to all vertices of U_v . As a result, the degree of each vertex in $U_1 \cup U_2 \cup \dots \cup U_n$ is 1, the degree of each vertex in V_2 is n , and the degree of each vertex in V_2 is $n + n \log^2 n$. Therefore, we have $\Delta(H) = \tilde{O}(n)$.

Because of the way we constructed the graph, it is not hard to see that we can find the i -th neighbor of the adjacency list of each vertex in H using only a single query to the adjacency matrix of G .

Observation 7.1. *For each vertex v in graph H , the i -th neighbor of H can be found using at most a single adjacency matrix query in G .*

Proof. **TOPROVE 19** □

Modification to the algorithm: Since the graph contains $\tilde{\Theta}(n^2)$ vertices, we cannot afford to apply the sparsification step to all vertices. However, vertices in $U_1 \cup \dots \cup U_n$ have degree 1. Therefore, we apply the sparsification step only to vertices in V_1 and V_2 . Since we have $|V_1| + |V_2| = 2n$, we can apply the sparsification for these sets in $\tilde{O}(n\sqrt{n})$ time. We first iterate over the vertices in V_2 and apply the sparsification step, and then we apply it to the vertices in V_1 . This ordering ensures that most vertices in V_2 get matched to vertices in $U_1 \cup \dots \cup U_n$ in this step, which is desirable for our application.

Claim 7.2. *After the sparsification step, each vertex in V_2 is matched by M with high probability. Moreover, at most $n/\log n$ vertices in V_2 are matched to vertices in $V_1 \cup V_2$ with high probability.*

Proof. **TOPROVE 20** □

Equipped with this reduction, we can now simply run the rest of the algorithm for vertices in V_1 . The only difference is that we exclude the edges of M that lie between V_2 and $U_1 \cup \dots \cup U_n$ in the estimation. Additionally, in the final estimation, the algorithm returns the previous estimate minus $n/\log n$, accounting for the vertices in V_2 that are not matched within $V_1 \cup V_2$, which introduces an additional $o(n)$ additive error. Thus we obtain:

Theorem 1.2. *There exists an algorithm that, given access to the adjacency matrix of a graph, estimates the size of the maximum matching with a multiplicative-additive approximation factor of $(0.5109, o(n))$ and runs in $\tilde{O}(n\sqrt{n})$ time with high probability.*

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