



# Exploring the LANDSCAPE of Distributed Graph Sketching

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## Abstract

Recent work has initiated the study of dense graph processing using graph sketching methods, which drastically reduce space costs by lossily compressing information about the input graph. In this paper, we explore the strange and surprising performance landscape of sketching algorithms. We highlight both their surprising advantages for processing dense graphs that were previously prohibitively expensive to study, as well as the current limitations of the technique. Most notably, we show how sketching can avoid bottlenecks that limit conventional graph processing methods.

Single-machine streaming graph processing systems are typically bottlenecked by CPU performance, and distributed graph processing systems are typically bottlenecked by network latency. We present LANDSCAPE, a distributed graph-stream processing system that uses linear sketching to distribute the CPU work of computing graph properties to distributed workers with no need for worker-to-worker communication. As a result, it overcomes the CPU and network bottlenecks that limit other systems. In fact, for the connected components problem, LANDSCAPE achieves a stream ingestion rate one-fourth that of maximum sustained RAM bandwidth, and is four times faster than random access RAM bandwidth. Additionally, we prove that for any sequence of graph updates and queries LANDSCAPE consumes at most a constant factor more network bandwidth than is required to receive the input stream. We show that this system can ingest up to 332 million stream updates per second on a graph with  $2^{17}$  vertices. We show that it scales well with more distributed compute power: given a cluster of 40 distributed worker machines, it can ingest updates 35 times as fast as with 1 distributed worker machine. Graph sketching algorithms tend to incur high computational costs when answering queries; to address this LANDSCAPE uses heuristics to reduce its query latency by up to four orders of magnitude over the prior state of the art.

The full version of the paper can be accessed at  
<https://arxiv.org/abs/2410.07518>

Our code and experiments can be found at <https://github.com/GraphStreamingProject/Landscape> and  
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## 1 Introduction

Computing connected components is a fundamental graph-processing task with uses throughout computer science and engineering. It has applications in relational databases [80], scientific computing [62, 72], pattern recognition [31, 40], graph partitioning [49, 50], random walks [38], social network community detection [46], graph compression [39, 48], medical imaging [33], flow simulation [73], genomics [27, 57], identifying protein families [54, 78], microbiology [3], and object recognition [32]. Strictly harder problems such as edge/vertex connectivity, shortest paths, and  $k$ -cores often use it as a subroutine. Connected Components is also used as a heuristic for clustering problems [22, 23, 24, 61, 76, 77], pathfinding algorithms (such as Dijkstra and  $A^*$ ), and some minimum spanning tree algorithms. A survey by Sahu et al. [66] of database applications of graph algorithms reports that, for both practitioners and academic researchers, connected components was the most frequently performed computation from a list of 13 fundamental graph problems that includes shortest paths, triangle counting, and minimum spanning trees.

Computing the minimum cut of a graph (or equivalently its edge connectivity) is a closely related problem to connected components. It has applications in clusterings on similarity graphs [79, 30], community detection [12], graph drawing [41], network reliability [64, 43], and VLSI design [45].

The task of computing connected components or minimum cut becomes more difficult when graphs are *dynamic*, meaning the edge set changes over time subject to a stream of edge insertions and deletions, and this task becomes harder still when the graphs are very large. Applications using dynamic graphs include identifying objects from a video feed rather than a static image [44] or tracking communities in social networks that change as users add or delete friends [9, 10]. Applications on large graphs include metagenome assembly tasks that may include gene

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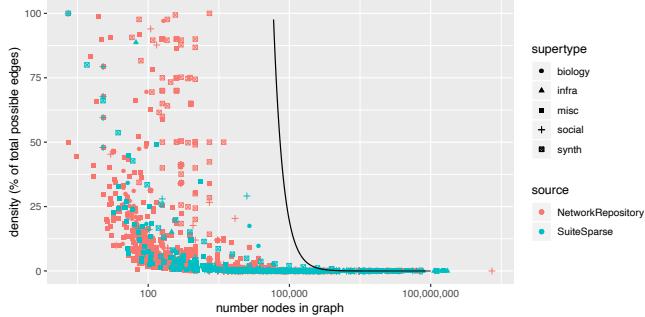


Figure 1: Graphs studied in academic works exhibit a selection effect. Any point to the left of the dark line indicates a dataset which can be represented as an adjacency list in 16GB of RAM.

databases with hundreds of millions of entries with complex relations [27], and large-scale clustering (a common machine learning challenge [24]). And of course graphs can be both large and dynamic. Indeed, Sahu et al.’s [66] database applications survey reports that a majority of industry respondents work with large graphs ( $> 1$  million vertices or  $> 1$  billion edges) and a majority work with dynamic graphs.

**Dense-graph processing.** The task of computing connected components is especially difficult for dynamic *dense graphs*. The conventional wisdom is that massive graphs are always sparse, meaning that they have few edges per vertex. Tench et al. [71] contend that instead large, dense graphs do not appear in academic publications due to a selection effect: since we lack the tools to process these graphs, they are not studied. We expand on their survey of graph datasets to further support this claim. Figure 1 plots all graph datasets from the NetworkRepository [65] and SuiteSparse [17] collections. Note that nearly all the graphs can be stored as an adjacency lists using less than 16GB, and this pattern holds across repositories and across different types of graph dataset - including a variety of biological data, social networks, and infrastructure(e.g., road and computer networks). See Appendix F.1 for an expanded analysis which shows this selection effect even more strongly.

Despite this effect, there is evidence of dense graphs emerging in practical applications. Tench et al. note that “Facebook works with graphs with 40 million nodes and 360 billion edges. These graphs are processed at great cost on large high-performance clusters, and are consequently not released for general study” [71]. As another example, bipartite projection methods [56], commonly used in social sciences and bioinformatics, naturally generate large, dense graphs. Current techniques require storing these graphs in RAM, limiting the size of datasets analyzed this way — an explicit example of the selection effect [58]. The only way to conclusively determine whether there exist more applications

that would benefit from dense-graph processing systems is to build such systems and see what applications emerge.

Tench et al. [71] demonstrate that *linear sketching* techniques [53, 1, 2, 29] can be used to process large, dense, and dynamic graphs. Linear sketching saves the most space when graphs are dense—this is because the size of a connectivity sketch of a  $V$ -vertex graph is  $O(V \text{ polylog } V)$  and therefore is independent of the number of edges. This algorithm is representative of a large number of graph-sketching algorithms [53, 1, 2, 42, 29, 6] in the database theory literature that all share the same general structure.

However, the price for the small space of the sketches is high CPU cost: processing each update requires  $O(\log^2 V)$  work, and the constants hidden by the asymptotic notation are large. Concretely, for a million-vertex graph, processing each edge *update* (an insertion or deletion) requires evaluating roughly 500 hash functions (in addition to other costs). Tench et al.’s implementation, GRAPHZEPPELIN, introduces some techniques for mitigating this high computational cost but ultimately their implementation is bottlenecked by CPU.

**This paper.** We design and implement distributed sketching algorithms for connected components and  $k$ -connectivity (or bounded  $k$  mincut) on dynamic graph streams. Using our implementation, we explore the surprising performance landscape of graph-sketching algorithms. Thus, we call this graph streaming system LANDSCAPE.

The peaks and valleys of this performance landscape will seem unfamiliar to graph-processing practitioners. The highest-order observation is that these techniques work well when graphs are dense and work poorly when they’re sparse. This is in contrast to most techniques which work better when graphs are sparse. We also show these sketching algorithms can avoid computational bottlenecks that are unavoidable using conventional graph-processing techniques, but these algorithms also struggle in some cases where traditional algorithms excel.

We summarize these findings in Claim 1 in Section 2 but for context we first briefly survey the bottlenecks in graph-stream processing and how they affect existing graph-processing systems (which are designed for sparse graphs).

### 1.1 Three Bottlenecks in Graph-Stream Processing

Graph-stream processing systems can be bottlenecked on any of three resources: *space*, *CPU*, and *communication*. Even a single-worker system requires some network communication; at minimum it must receive the input stream from a network link.

Some systems, such as Aspen [18] and Terrace [60], are bottlenecked on *space*. These systems maintain lossless representations of the input graph in RAM on a single machine, and such lossless representations can be large. These systems are optimized for processing sparse graphs

Table 1: Summary of ingestion bottlenecks. Existing graph systems are bottlenecked by space, CPU, or network communication costs when processing dense graph streams. In contrast, in this paper we present a system that overcomes these three bottlenecks.

	space	CPU	Network
single-machine lossless [18, 60]	★ $\Theta(V^2)$	★★	N/A
single-machine sketching [71]	★★★ $\Theta(V \log^3(V))$	★	N/A
distributed lossless [13, 34, 74, 35, 36, 26]	★★ $\Theta(V^2)$	★★	★
distributed sketching (this paper)	★★★ $\Theta(V \log^3(V))$	★★	★★★

\* = bottleneck; \*\* = good; \*\*\* = optimal

where this memory burden is less onerous but struggle when processing large, dense graphs.

Other systems, such as GRAPHZEPPELIN [71], are bottlenecked on **computation** (as we explain earlier). GRAPHZEPPELIN maintains a lossily-compressed graph representation, reducing storage requirements (and improving performance on dense graphs) at the cost of higher CPU load for processing updates and answering queries.

Distributed graph-stream processing systems are bottlenecked by network **communication**. For dynamic systems [74], the input stream is received at a central node called the **main node** and information about the graph is distributed across the cluster to **worker nodes**. For non-dynamic systems the data is distributed among worker nodes before computation begins. Since the graph data is spread across many worker nodes, and graphs often have poor data locality, most computation on graphs require worker nodes to send lots of information to each other [13, 34, 74, 35, 36, 26]. This is an example of a general challenge in distributed database systems. These systems scale by spreading data among the aggregated memory of nodes in a cluster at the cost of high inter-node communication, which can increase query and transaction latencies [81, 19, 20].

Table 1 summarizes the bottlenecks for each of these approaches.

**Objective standard for update performance: RAM bandwidth.** One issue with evaluating the performance of a distributed system to compute fully-dynamic (i.e., edges can be inserted or deleted) graph connectivity is that there are no other open-source distributed systems that solve this precise problem (see Appendix C). However, for stream ingestion rate we can compare against objective upper bound.

The update performance of any graph-streaming system is limited by the **data acquisition cost**, that

is, the cost for the main node to simply read the entire input stream. Even if we ignore the cost to update the connectivity information in the graph data structures, we still need to read the input. Thus, an objective standard describing the ideal performance is simply the RAM bandwidth. In fact, there are two notions of RAM bandwidth: **random-access RAM bandwidth** (the speed at which we can write words to random locations) and **sequential-access RAM bandwidth** (the speed at which we can write words to sequential locations).

Since graphs have notoriously poor data locality, an update rate close to random-access RAM bandwidth is a natural goal for a graph stream-processing system. Sequential-access RAM bandwidth is truly a bound on the best possible update rate because any stream-processing system must write the input data into memory, that is, it must pay the data acquisition cost.

## 2 Results

In this paper, we build the LANDSCAPE graph-processing system, which is optimized for dense, dynamic graphs. We use this system to establish how graph sketching can overcome classical graph processing bottlenecks.

### 2.1 The Landscape Graph-Processing System

We build LANDSCAPE, a linear-sketch-based distributed graph-stream processing system that computes connected components and k-connectivity of dynamic graphs.

LANDSCAPE keeps its sketch (a lossily compressed graph representation) on the main node and distributes the CPU work of processing updates.

Because these linear sketches are small (size  $\Theta(V \log^3 V)$ ), they fit on a single node, even when the graph is dense. The work to maintain these sketches can be chunked off into large batches that can be computed independently by worker nodes. As a result, LANDSCAPE avoids the CPU bottleneck because it has lots of worker nodes to help, and LANDSCAPE avoids the communication bottleneck because the communication cost is amortized away by the CPU cost to process a batch.

In fact, we prove theoretically that for any sequence of graph updates and queries LANDSCAPE’s total communication cost is only a small, user-configurable constant (by default, four) factor larger than the cost for the main node to receive the input stream. Additionally, we introduce a new sketching algorithm, CAMEOSKETCH, which requires only  $O(\log V)$  distributed work per update (compared to  $O(\log^2 V)$  for the prior state of the art), which allows the algorithm to scale more rapidly with limited cluster resources.

**Performance.** LANDSCAPE achieves the following:

- LANDSCAPE is able to process graph streams only 4.5× slower than the multi-threaded RAM sequential-write bandwidth, the objective upper bound on insertion

performance for any system that receives the input stream at the main node. This is more than four times faster than random access RAM bandwidth.

- We show that LANDSCAPE can ingest up to 332 million stream updates per second on a graph with  $2^{17}$  vertices.
- We show that it scales well with more distributed compute power: given a cluster of 40 distributed worker machines, it can ingest updates 35 times as fast as with 1 distributed worker machine.
- We experimentally verify that LANDSCAPE uses at most  $4\times$  the network bandwidth required to read the input stream.
- LANDSCAPE’s GREEDYCC query heuristic reuses partial information from prior query results, achieving up to a four orders-of-magnitude reduction in query latency.

**Outperforming lossless representations on dense graphs.** To put this performance in context, consider a simpler task: maintaining an adjacency matrix of the graph defined by the input stream<sup>1</sup>. If the graph is dense and edges are random, an adjacency matrix is essentially the space-optimal lossless graph representation. We ignore the cost of answering queries, which an adjacency matrix does not efficiently support.

LANDSCAPE’s graph-sketch representation is smaller than this adjacency matrix even when the input graph has only 310,000 vertices. Even more interestingly, LANDSCAPE’s update throughput is also faster than the update throughput of the adjacency-matrix representation—which is just a single bit flip per edge. We emphasize: one of the most dramatic advantages of distributed graph sketching is that updates are faster than adjacency-matrix updates even when the entire adjacency matrix fits in RAM.

LANDSCAPE’s updates are fast not because they are small (you cannot beat a single bit flip), but because the CAMEOSKETCHES have good data locality—and the edge updates result in primarily sequential accesses to RAM on the main node, rather than random access. Said differently, LANDSCAPE processes more edge updates per second than it is possible to flip bits in random locations in RAM.

**2.2 Circumventing the Classical Bottlenecks for Graph-Stream Processing** The performance implications of sketching for dense graph processing are encapsulated in the following claim. Throughout this paper, as we present theoretical and experimental results, we will refer to the element of the claim they support.

### CLAIM 1. (Dense graph processing)

1. **Space consumption.** Graph sketches for dynamic, massive, dense graphs can be maintained so that they use less space than traditional graph-storage methods. E.g., LANDSCAPE is asymptotically space-optimal and its new CAMEOSKETCH algorithm uses 29% of the space of the prior state-of-the-art [71].
2. **CPU cost.** Graph sketches have high CPU cost to update, but this cost can be distributed away. E.g., LANDSCAPE’s CAMEOSKETCH algorithm reduces the asymptotic work per update from  $O(\log^2 V)$  (the prior SOTA) to  $O(\log V)$ . We show that this yields a  $7\times$  increase to update throughput in experiments.
3. **Communication costs.** The above distribution of CPU work can be done with nearly optimal communication: specifically, the total communication cost is only a small constant number of times larger than the data acquisition cost. We prove this theoretically and validate it experimentally.
4. **Stream-ingestion can be blindingly fast—nearly the universal speed limit.** There is a universal speed limit for stream ingestion, which is simply the cost to write the stream sequentially into RAM. A graph-sketch based system for connectivity and  $k$ -connectivity can match this bound within a remarkably small constant factor. LANDSCAPE ingests graph data at a rate that is within a factor 4 of sequential RAM bandwidth.

Table I summarizes the strengths of the sketching approach: LANDSCAPE ingestion is not bottlenecked on space, CPU, or communication.

The following observations are nearly immediate and unsurprising, and we include them for completeness.

**Queries.** Similar to the trade-off between reading and writing in (nongraph) databases [50, 8, 67, 7], ingesting a graph stream and processing it into a sketch is faster than querying the sketch. LANDSCAPE answers each query in single digit seconds, even when there are 37 billion edges and  $2^{19}$  vertices, using a combination of provable worst-case query algorithms, accelerated with powerful heuristics.

**Sparse graph-processing via sketching.** Graph sketching is not the best solution for graphs that are relatively sparse. On sparse graphs, the LANDSCAPE approach retains its asymptotic guarantee of low communication but is not space-efficient and may not be able to distribute away its CPU costs. For completeness we evaluate LANDSCAPE’s performance on sparse graphs and validate these theoretical predictions.

<sup>1</sup>Note that while adjacency matrices may be compact, they have high query latency. Even disregarding this limitation, we see that adjacency matrices are outperformed by sketching on dense graphs.

### 3 Preliminaries & Definitions

**The Graph Streaming Model.** In the *graph semi-streaming* model [25] (sometimes just called the *graph streaming* model), an algorithm is presented with a *stream*  $S$  of updates (each an edge insertion or deletion) where the length of the stream is  $N$ . Stream  $S$  defines an input graph  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  with  $V = |\mathcal{V}|$  and  $E = |\mathcal{E}|$ . The challenge in this model is to compute (perhaps approximately) some property of  $\mathcal{G}$  given a single pass over  $S$  and at most  $O(V \text{polylog } V)$  words of memory. Each update has the form  $((u, v), \Delta)$  where  $u, v \in \mathcal{E}$ ,  $u \neq v$  and  $\Delta \in \{-1, 1\}$  where 1 indicates an edge insertion and  $-1$  indicates an edge deletion. Let  $s_i$  denote the  $i$ th element of  $S$ , and let  $S_i$  denote the first  $i$  elements of  $S$ . Let  $\mathcal{E}_i$  be the edge set defined by  $S_i$ , i.e., those edges which have been inserted and not subsequently deleted by step  $i$ . The stream may only insert edge  $e$  at time  $i$  if  $e \notin \mathcal{E}_{i-1}$ , and may only delete edge  $e$  at time  $i$  if  $e \in \mathcal{E}_{i-1}$ .

In this paper, we consider the following two problems in this model:

**PROBLEM 1. (Streaming Connected Components.)**

*Given an insert/delete edge stream of length  $N$  that defines a graph  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$ , return a spanning forest of  $\mathcal{G}$ .*

**PROBLEM 2. (Streaming  $k$ -Edge Connectivity.)**

*Given an insert/delete edge stream of length  $N$  that defines a graph  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$ , for any cut  $C \subset \mathcal{V}$ , return the cardinality of cut  $C$  (denoted by  $w(C)$ ) if  $w(C) < k$ , and return  $\infty$  otherwise.*

Prior work [1, 7] has also considered these problems in the graph streaming model.

**A note on the query model for Landscape.** The graph streaming model above requires computing an answer at the end of the stream. In contrast, LANDSCAPE can answer connectivity queries interspersed arbitrarily among insertions and deletions during the stream. It answers both *global connectivity* queries, where the task is to return the connected components or  $k$ -edge connectivity of the graph, and *batched reachability* queries, where the query consists of a set of vertex pairs  $(u_1, v_1), (u_2, v_2) \dots (u_k, v_k)$  and the task is to determine whether  $u_i$  is in the same connected component as  $v_i$  for each  $i \in [k]$ . The goal is to minimize query latency in addition to the goals of minimizing space use and maximizing stream ingestion considered by prior work. For more information see Section 5.3. We assume that the stream is not adaptive, meaning that edge updates and connectivity queries do not depend on the results of prior queries.

**Other models in streaming connected components systems.** Aspen [18] and Terrace [60] are graph-stream processing systems which support connectivity queries. Unlike the semi-streaming model proposed by [1] above, these systems work in the *batch-dynamic model*,

where updates are applied to a non-empty graph in batches exclusively containing insertions or deletions. The minimum size of batches is a parameter which affects system performance. Connectivity queries may be issued in between batches, but not during a batch. Aspen is capable of computing queries concurrently with processing updates while Terrace is not. In contrast, LANDSCAPE is designed to handle arbitrarily interspersed updates and queries with no notion of batched input. Like Terrace, it does not compute queries concurrently with processing updates.

### 4 Sketching Graphs

In this section we briefly review the prior work on connectivity sketching, and then present CAMEOSKETCH, an improved sketching subroutine that reduces the update cost from  $O(\log^2 V)$  to  $O(\log V)$ , and reduces the sketch size by a significant constant factor (73%).

**4.1 Prior Work** Ahn et al. [1] initiate the field of graph sketching with their connected components sketch, which solves the streaming connected components problem in  $O(V \log^3 V)$  space. A key subproblem in their algorithm is  *$\ell_0$ -sampling*: a vector  $x$  of length  $n$  is defined by an input stream of updates of the form  $(i, \Delta)$  where value  $\Delta$  is added to  $x_i$ , and the task is to sample a nonzero element of  $x$  using  $o(n)$  space. They use an  $\ell_0$ -sampler (also called an  $\ell_0$ -sketch) due to Cormode et al.:

**THEOREM 4.1.** (Adapted from [16], Theorem 1): *Given a 2-wise independent hash family  $\mathcal{F}$  and an input vector  $x \in \mathbb{Z}^n$ , there is an  $\ell_0$ -sampler using  $O(\log^2(n) \log(1/\delta))$  space that succeeds with probability at least  $1 - \delta$ .*

We denote the  $\ell_0$  sketch of a vector  $x$  as  $\mathcal{S}(x)$ . The sketch is a linear function, i.e.,  $\mathcal{S}(x) + \mathcal{S}(y) = \mathcal{S}(x + y)$  for any vectors  $x$  and  $y$ . Ahn et al. define a *characteristic vector*  $f_u \in \mathbb{Z}_2^{\binom{V}{2}}$  of each vertex  $u \in \mathcal{V}$  such that each nonzero element of  $f_u$  denotes an edge incident to  $u$ . That is,  $f_u \in \mathbb{Z}_2^{\binom{V}{2}}$  s.t. for all vertices  $0 \leq i < j < V$ :

$$f_u[(i, j)] = \begin{cases} 1 & u \in \{i, j\} \text{ and } (i, j) \in \mathcal{E} \\ 0 & \text{otherwise} \end{cases}$$

Further, for any  $S \subset \mathcal{V}$ , the nonzero elements of  $f_S = \sum_{u \in S} f_u$  are precisely the set of edges crossing the cut  $S, \mathcal{V} \setminus S$ . For each  $u \in \mathcal{V}$ , the algorithm computes  $\mathcal{S}(f_u)$ : for each edge update  $(u, v, \Delta)$  it computes  $\mathcal{S}(e_{(u, v)})$  where  $e_{(u, v)}$  denotes the vector with 1 in position  $\text{idx} = (u, v)$  and 0 in all other positions. It maintains  $\mathcal{S}(f_u) = \sum_i \mathcal{S}(x_i)$  for each  $u$ . Then for arbitrary  $X \subset \mathcal{V}$  they can sample an edge across the cut  $X, \mathcal{V} \setminus X$  by computing  $\mathcal{S}(f_X) = \sum_{u \in X} \mathcal{S}(f_u)$ .

This allows them to perform Boruvka's algorithm using the sketches: they form  $O(\log V)$   $\ell_0$ -sketches

$\mathcal{S}_0(f_u), \mathcal{S}_1(f_u), \dots, \mathcal{S}_{O(\log V)}(f_u)$  for each  $u$ . We call  $\mathcal{S}(f_u) = \bigcup_{i \in [O(\log V)]} \mathcal{S}_i(f_u)$  the **vertex sketch** of  $u$ , and it has size  $O(\log^3 V)$ . Then  $\forall u \in \mathcal{V}$  they query  $\mathcal{S}_1(f_u)$  to sample an edge incident to  $u$ . For each resulting component  $X$  they compute  $\mathcal{S}_2(f_X)$  and repeat until all connected components are found. Since each vertex sketch  $\mathcal{S}(f_u)$  has size  $O(\log^3 V)$  bits, the entire data structure has size  $O(V \log^3 V)$ . See Appendix A for a more complete description of this algorithm, along with an illustrative example.

**Testing k-connectivity.** Ahn et al. [1] also show how to test k-connectivity of a graph  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  (that is, to exactly compute the minimum cut of  $\mathcal{G}$  provided this value is  $\leq k$ ) by constructing a k-connectivity certificate  $H = \bigcup_{i \in [k]} F_i$  where  $F_0, F_1, \dots, F_{k-1}$  are edge-disjoint spanning forests of  $\mathcal{G}$ .  $H$  has the property that it is  $k'$ -edge connected iff  $G$  is  $k'$ -edge connected for all  $k' \leq k$ . They find each  $F_i$  by computing  $k$  connectivity sketches of  $\mathcal{G}$  in a single pass over the stream. After the stream, the first connectivity sketch is used to find  $F_0$  and the edges of  $F_0$  are deleted from the remaining  $k - 1$  connectivity sketches. The second connectivity sketch can now be used to get  $F_1$ , whose edges are subsequently deleted from the remaining  $k - 2$  sketches and so on. The size of the sketches is  $O(kV \log^3 V)$  bits.

**GraphZeppelin.** Tench et al. [71] present GRAPHZEPPELIN, the first implementation of Ahn et al.’s connected components algorithm which uses a faster  $\ell_0$ -sketch algorithm which they call CUBESKETCH. GRAPHZEPPELIN also uses an external-memory-optimized data structure called a *gutter tree* to I/O-efficiently collect updates to be processed, allowing the algorithm to run quickly even when the sketches are stored on disk.

To achieve a failure probability of  $\delta$ , GRAPHZEPPELIN's CUBESKETCH uses  $O(\log^2(n) \log(1/\delta))$  bits of space and has worst-case update time  $O(\log(n) \log(1/\delta))$ . As in Ahn et al., they use  $O(\log V)$  of these sketches for each vertex and set  $\delta$  to be a small constant. So the space cost per vertex is  $O(\log^3 V)$  bits and the worst-case update time is  $O(\log^2 V)$ . See Appendix B for a detailed description of CUBESKETCH.

**4.2 Landscape’s new sketch: CameoSketch** We develop a new  $\ell_0$  sampler called CAMEOSKETCH for use in LANDSCAPE. CAMEOSKETCH improves upon CUBESKETCH with a new update procedure that is a  $O(\log V)$  factor faster to update and reduces space usage by a constant factor via a refined analysis. All other details, including the query procedure, remain unchanged. We present the full details of CAMEOSKETCH in Appendix B and report the asymptotic performance here.

**Update procedure.** CAMEOSKETCH uses a simpler and faster update procedure than CUBESKETCH.

THEOREM 4.2. CAMEOSKETCH is an  $\ell_0$ -sampler that, for vector  $x \in \mathbb{Z}_2^n$ , uses  $O(\log^2(n) \log(1/\delta))$  space, has

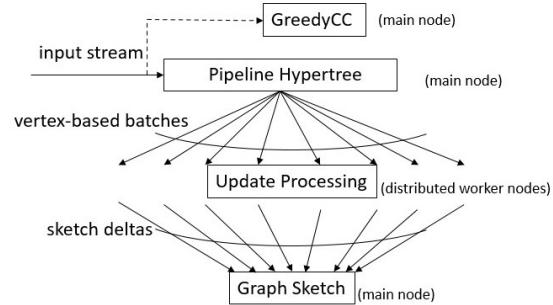


Figure 2: Data flow diagram for LANDSCAPE’s ingestion algorithm. Connectivity information from the input stream is compressed into the graph sketch and added to GREEDYCC.

worst-case update time  $O(\log(1/\delta))$ , and succeeds with probability  $1 - \delta$ .

Theorem 4.2 demonstrates that CAMEOSKETCH reduces the CPU burden of performing updates and supports Claim 1.2. The proof of this theorem can be found in Appendix B.

**Reduced constant factors.** In GRAPHZEPPELIN, Tench et al. use  $56 \log(1/\delta) \log n$  bytes of space to guarantee a failure probability of at most  $\delta$  when sketching a vector of length  $n < 2^{64}$  using CUBESKETCH. Via a careful constant-factor analysis, we can show that CAMEOSKETCH can match this failure probability with significantly less space:

**THEOREM 4.3.** *Using 3-wise independent hash functions, CAMEOSKETCH requires  $8 \log_3(1/\delta)(\log n + 5)$  bytes of space to return a nonzero element of a length  $n < 2^{64}$  input vector  $w/p$  at least  $1 - \delta$ .*

Theorem 4.3 immediately implies a space savings of up to 90% compared to CUBESKETCH and thus supports Claim 11. In our implementation, we conservatively choose to use slightly more space than this theorem requires to reduce the failure probability further. Still, our implementation requires only 2/7ths of the space used in GRAPHZEPPELIN [71] (see Section 6 for details).

See Appendix H for the proof of this theorem.

## 5 Landscape Design

LANDSCAPE uses CAMEOSKETCHES to compute connectivity. The CPU work of computing updates to  $\mathcal{S}(\mathcal{G})$  is done by distributed workers, while  $\mathcal{S}(\mathcal{G})$  itself is stored on the main node. To answer queries, the main node computes a spanning forest via Borůvka's algorithm using  $\mathcal{S}(\mathcal{G})$  as described in Section 4 or via a heuristic algorithm which we call GREEDYCC and describe in Appendix E.4.

In this section we describe how LANDSCAPE’s main node efficiently collects updates into vertex-based batches to minimize communication, how the distributed workers process these batches of updates into sketch form, and how LANDSCAPE answers connectivity queries. We also prove the asymptotic upper bounds on the CPU and communication costs incurred by these operations. This analysis explains LANDSCAPE’s surprising performance profile.

Figure 2 summarizes the LANDSCAPE data flow. The input stream, consisting of edge insertions, edge deletions, and queries arrives at the main node. Updates (insertions and deletions) are inserted into the *pipeline hypertree* (Section 5.1.2) and also into GREEDYCC (Section E.4). The pipeline hypertree collects updates into *vertex-based batches* (Section 5.1.1) which are sent to distributed worker nodes. These worker nodes process batches by running the CAMEOSKETCH algorithm, producing *sketch deltas* (Section 5.2) which are applied to the CAMEOSKETCHES on the main node.

## 5.1 Ingesting Stream Updates on the Main Node

**5.1.1 Vertex-Based Batching** The core technique that makes distributed sketch processing communication-efficient (and therefore feasible) is *vertex-based batching*, where many updates with a common endpoint are collected into a batch. Intuitively, because one or many updates to the same endpoint can be represented as a single sketch delta of fixed size, batching updates by endpoint drastically reduces the communication cost of sending sketch deltas from worker nodes to the main node.

Specifically, any updates for  $u \in \mathcal{V}$  are collected into a batch  $B_u \subseteq \{(x, y) \in \mathcal{E} \mid x = u \vee y = u\}$ .  $B_u$  is sent to a single distributed worker, which returns a sketch of the updates. As we will see in Section 5.2 this sketch has size  $\phi = O(\log^3 V)$  bits. During update processing, LANDSCAPE only sends  $B_u$  when  $|B_u| \geq \alpha\phi/\log V$  for some constant  $\alpha \geq 1$ . Each update requires  $\log V$  bits to represent, so a buffer that contains  $\alpha\phi/\log V$  updates has size  $\alpha\phi$  bits.

As a result of this policy, the amortized communication cost per update is small. Say the stream contains  $N$  updates. The bandwidth cost to receive the input stream is  $N$ . Since each update is included in two vertex-based batches (one per endpoint) and each batch is sent to a distributed worker once, the total bandwidth cost of sending vertex-based batches is  $2N$ . Finally, each vertex-based batch induces the distributed worker that receives it to respond with a sketch delta (which is  $1/\alpha$  of the batch’s size). This means that as long as LANDSCAPE is processing full vertex-based batches, the network bandwidth cost of processing  $N$  updates is at most  $(3 + 1/\alpha)N$ . This technique is simple, but crucial for good performance.

**5.1.2 Pipeline Hypertree** To make vertex-based batching fast, we design the *pipeline hypertree*, which is a simplified and parallel variant of the buffer tree [5] designed to minimize cache line misses and thread contention. The pipeline hypertree receives arbitrarily ordered stream updates and consolidates them into vertex-based batches. Each update inserted into the pipeline hypertree is moved  $O(\log_{C/\mathcal{L}} V)$  times before being returned in a vertex-based batch, where  $C$  denotes the size of L3 cache and  $\mathcal{L}$  denotes the size of an L3 cache line. The total size of the data structure is  $O(V \log^3 V)$  bits. We defer description of the design and implementation of the pipeline hypertree to Appendix C.

## 5.2 Distributed Sketch Processing

CAMEOSKETCHES have strong data locality, which we exploit for parallelism: processing a graph edge update  $(u, v, \Delta)$  requires updating only  $\mathcal{S}(f_u)$  and  $\mathcal{S}(f_v)$  and these updates can be performed independently of each other. Because the sketches are linear, the sketch update for  $(u, v)$  can be computed on its own and later summed to the sketch of  $u$ . This means that the vast majority of the computation required to process  $(u, v)$  can be performed before accessing the sketches for  $u$  and  $v$ . LANDSCAPE exploits this independence to distribute the computational cost of these updates while storing the sketches on a single worker.

When a batch of updates  $(e_1, e_2, \dots)$  for vertex  $u$  is sent to a worker node, the worker node computes  $\sum_j \mathcal{S}(e_j)$ , which we call a *sketch delta* (since it is a sketch encoding the change in the neighborhood of vertex  $u$ ). Note that it has size  $O(\log^3 V)$ —equal to the size of a vertex sketch. This immediately gives the following result:

**THEOREM 5.1. (DISTRIBUTED COST.)** *The distributed CPU cost of processing a batch of  $x$  updates for vertex  $u \in \mathcal{V}$  into a sketch delta  $\mathcal{S}_u$  is  $O(x \log(V))$ .*

**Sketch merging.** LANDSCAPE’s main node maintains the *graph sketch*:  $\mathcal{S}(\mathcal{G}) = \bigcup_{u \in \mathcal{V}} \mathcal{S}(f_u)$ . After a sketch delta for vertex  $u$  is created by a distributed worker is then sent to the main node where it is added to  $\mathcal{S}(f_u)$ . The graph sketch is stored in RAM which is feasible since it has total size  $O(V \log^3(V))$  bits.

**5.3 Finding Connected Components** LANDSCAPE is designed to efficiently answer two types of connected component queries: *global connectivity queries* where the task is to map each vertex to its connected component, and *batched reachability queries* where the query consists of a set of vertex pairs  $(u_1, v_1), (u_2, v_2), \dots, (u_k, v_k)$  and the task is to determine whether  $u_i$  is in the same connected component as  $v_i$  for each  $i \in [k]$ . At a high level, LANDSCAPE answers these queries by producing a spanning forest of the

graph defined by the input stream. This is done as described in Section 4 via Boruvka's algorithm on the vertex sketches.

Processing queries can increase network bandwidth costs if not handled carefully. Computing the spanning forest from the sketches can only be done after all pending stream updates have been processed. We say the graph sketch stored on the main node is *current* with respect to query  $q$  at time  $t$  if there are no pending updates; i.e., all updates that arrived prior to time  $t$  have been processed and merged into the graph sketch. If for some vertex  $i$  there are a small number of updates for  $i$  in the pipeline hypertree, then by the reasoning in Section 5.1.1 this could incur an average communication cost per update of  $O(\log^2 V)$ . In the worst case, the input stream could insert a perfect matching into an empty graph and then issue a query; the average communication cost per update would be  $O(\log^2 V)$ .

LANDSCAPE avoids this problem by adopting a hybrid distribution policy for pending updates when a query is issued. When a query is issued, LANDSCAPE first flushes the pipeline hypertree so all pending updates are stored in the leaves. For each leaf, if the leaf is at least a  $\gamma$ -fraction full, it is sent as a vertex-based batch to a distributed worker, where  $\gamma \in (0, \frac{1}{2}]$  is a parameter chosen by the user. All leaves that are less than a  $\gamma$ -fraction full are processed locally on the main node, costing no additional network bandwidth.

As a consequence of this policy, we have the following theorems.

**THEOREM 5.2. (COMMUNICATION COST.)** *The communication cost of ingesting  $N$  updates and answering  $Q$  queries is at most  $(3 + 1/(\gamma\alpha))N$ , where  $\gamma$  and  $\alpha$  are constants.*

*Proof.* See Appendix D.  $\square$

Importantly, this means that LANDSCAPE never uses more than a constant multiple of the network bandwidth required to receive the input stream, *regardless of the number or distribution of queries*.

**THEOREM 5.3. (COMPUTATIONAL COST ON MAIN NODE.)** *Given a input stream of length  $N$  defining  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  and a series of  $Q$  connectivity queries issued throughout the stream, let  $N_i$  denote the number of edge updates that arrive after query  $Q_i$  but before query  $Q_{i+1}$ . LANDSCAPE never uses more than  $O(V \log^3 V)$  bits of space on the main node while processing the stream, and the amortized cost per update to process  $N_i$  is  $O(\log_{C/\mathcal{L}}(V))$  if  $N_i = \Omega(V \log^2(V))$  and  $O(\log(V))$  otherwise. Computing each query  $Q_i$  takes  $O(V \log^2(V))$  time.*

*Proof.* See Appendix D.  $\square$

As a result, when  $N_i = \Omega(V \log^2 V)$  the amortized CPU cost for all computation on the main node is  $O(\log_{C/\mathcal{L}}(V))$  (the amortized cost to process updates with

the pipeline hypertree). For reasonable values of  $C$  and  $\mathcal{L}$  this logarithm evaluates to a small constant, typically 3. Moreover, these operations are just data movement operations, so they can be done at near RAM bandwidth.

**5.4 Computing k-connectivity** LANDSCAPE's architecture can in principle accomodate many other graph sketch algorithms that use connectivity as a subroutine. For instance, since the  $k$ -connectivity sketch requires maintaining  $k$  independent copies of the connectivity sketch, we can achieve comparable computation and communication upper bounds by slightly modifying the procedure used to distribute connectivity sketching. Set the size of a vertex-based batch and the size of leaf node buffers in the pipeline hypertree to  $\alpha \cdot k \log^3 V$  (the per-vertex sketch size for  $k$ -connectivity). When a distributed worker receives a vertex-based batch, it computes the sketch delta of the batch for all  $k$  copies of the connectivity sketch and sends this back to the main node. To answer  $k$ -connectivity queries, LANDSCAPE produces a  $k$ -connectivity certificate using the query algorithm summarized in Section 4. This immediately gives the following result:

**THEOREM 5.4. (K-CONNECTIVITY.)** *Given a input stream of length  $N$  defining  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  and a series of  $Q$  connectivity queries issued throughout the stream, let  $N_i$  denote the number of edge updates that arrive after query  $Q_i$  but before query  $Q_{i+1}$ . LANDSCAPE never uses more than  $O(kV \log^3 V)$  bits of space on the main node while processing the stream. The main node amortized cost per update to process  $N_i$  is  $O(\log_{C/\mathcal{L}}(V))$  if  $N_i = \Omega(kV \log^2 V)$  and  $O(k \log V)$  otherwise. Computing each query  $Q_i$  takes  $O(k^2 V \log^2 V)$  time. The distributed CPU cost of processing a batch of  $x$  updates for vertex  $u \in \mathcal{V}$  into a sketch delta  $\mathcal{S}_u$  is  $O(xk \log V)$ . The communication cost of ingesting  $N$  updates and answering  $Q$  queries is at most  $(3 + 1/(\gamma\alpha))N$ , where  $\gamma$  and  $\alpha$  are constants.*

Note that the network communication cost does not increase above that of connectivity, and for sufficiently infrequent queries the cost to the main node is also independent of  $k$ . See Section 7.4 for experimental confirmation of these results.

## 6 Landscape Implementation

Processing stream updates into the graph sketch is a computationally intensive process: for a moderately sized data-set with  $2^{18}$  vertices, applying a single edge update requires evaluating 184 hash functions. LANDSCAPE farms out this computationally intensive portion of the workload to worker nodes while the other portions of stream ingestion, including update buffering and sketch storage, remain the responsibility of the main node. LANDSCAPE uses  $164V * (\log^2 V - \log V)B$  of space on the main node

to store the sketches and the pipeline hypertree. On the worker nodes, LANDSCAPE requires storage for a single sketch and batch per CPU. Thus, a worker node with  $t$  threads requires  $t \cdot 164(\log^2 V - \log V)$  bytes. On a billion-vertex graph, each worker thread requires only 64 KiB of RAM. When computing  $k$ -connectivity, all of the above costs are multiplied by a factor  $k$ .

LANDSCAPE must handle two tasks: stream ingestion, where edge updates from the input stream are compressed into the graph sketch; and query processing, where connectivity queries are computed from the graph sketch (or sometimes from auxiliary query-accelerating data structures, described below). We defer a full description of the implementation, including auxiliary data structures, parameter choices, and software tools, to Appendix E.

## 7 Experiments

**Experiment setup.** We implemented LANDSCAPE in C++14 and compiled using g++ version 9.3 with openmpi 4.1.3 for Linux. We ran our experiments on an AWS cluster composed of an c5n.18xlarge instance for a main node and 40 c5.4xlarge instances as worker nodes. These instances have respectively 36 and 8 2-way hyperthreaded Intel(R) Xeon(R) Platinum 8124M CPU @ 3.00GHz cores with respectively 196 GB and 32 GB of RAM.

Each of our worker nodes only requires 2 GB of RAM because sketch deltas are small and workers are stateless. However, AWS workers with sufficient CPU power come with more RAM than we need.

**A note about experimental comparisons.** Ideally, we would include experimental comparisons against existing distributed systems that solve connectivity or  $k$ -connectivity on graph streams with edge insertions and deletions. However, only one such system (KickStarter [74]) exists in the literature, and its source code is not available (see Appendix G). Instead, we compare against a theoretical upper bound for stream ingestion: the data acquisition cost.

We define the **data acquisition cost** to be the cost of receiving the input stream at the main node over a network link and logging it in RAM. Sequential RAM bandwidth is a trivial bound on this cost.

- **RAM sequential bandwidth** is the maximum rate at which the CPU can write consecutive words in RAM on any number of threads.
- **RAM random access bandwidth** is the maximum worst-case rate at which the CPU can write any sequence of words in RAM on any number of threads.

**Experiment Metrics.** Our experiments process the entire stream of updates and then perform a single query at the end of the stream. Where specified experiments also

Table 2: Datasets used in our experiments.

Name	Vertices	Edges	Stream Updates
<b>kron13</b>	$2^{13}$	$1.7 \times 10^7$	$1.2 \times 10^8$
<b>kron15</b>	$2^{15}$	$2.7 \times 10^8$	$1.9 \times 10^9$
<b>kron16</b>	$2^{16}$	$1.1 \times 10^9$	$7.7 \times 10^9$
<b>kron17</b>	$2^{17}$	$4.3 \times 10^9$	$3.1 \times 10^{10}$
<b>ca-citeseer</b>	$2.3 \times 10^6$	$8.1 \times 10^5$	$1.1 \times 10^8$
<b>p2p-gnutella</b>	$6.3 \times 10^4$	$1.5 \times 10^5$	$1.9 \times 10^6$
<b>rec-amazon</b>	$9.2 \times 10^4$	$1.3 \times 10^5$	$1.7 \times 10^6$
<b>google-plus</b>	$1.1 \times 10^5$	$1.4 \times 10^7$	$1.9 \times 10^8$
<b>web-uk-2005</b>	$1.3 \times 10^6$	$1.2 \times 10^8$	$1.6 \times 10^9$
<b>erdos18</b>	$2^{18}$	$1.7 \times 10^{10}$	$4 \times 10^{10}$
<b>erdos19</b>	$2^{19}$	$3.4 \times 10^{10}$	$4 \times 10^{10}$
<b>erdos20</b>	$2^{20}$	$8 \times 10^{10}$	$1 \times 10^{11}$

perform additional queries throughout the processing of the stream.

We report the update throughput by measuring wall-clock time from the beginning of the stream until all updates have been applied to the sketches. We also measure the total amount of network communication to/from the main node and the amount of RAM usage on the main node.

When performing a query we measure the wall-clock latency from the moment the query is issued to the time the answer is returned to the user. This time includes the latency of flushing all pending updates from the pipeline hypertree and then the query computation itself.

**7.1 Datasets** In many of the experiments below, we use the synthetically generated graph streams used in the evaluation of GRAPHZEPPELIN [71]. These graphs were generated using the Graph500 Kronecker graph generator specification [4], and are very dense: each graph contains approximately 1/4 of all possible edges.

For larger-scale experiments, we evaluate LANDSCAPE on randomly generated Erdos-Renyi graphs (with  $2^{18}$ ,  $2^{19}$ , and  $2^{20}$  vertices) with edge probability set to 1/4.

Finally, we evaluate LANDSCAPE on real-world graph datasets from the SNAP graph repository [47] and NetworkRepository [65].

All of the above graphs were transformed into a random streams of edge insertions and deletions using the method described in [71]. We additionally inserted and removed all edges seven times to increase stream length. Each random stream, once insert/delete pairs for the same edge are removed, is exactly the edge list of the graph used to generate it. These datasets are summarized in Table 2.

**7.2 Landscape is Highly Scalable.** We measured LANDSCAPE’s stream ingestion rate and network bandwidth usage on kron17 given varying numbers of distributed

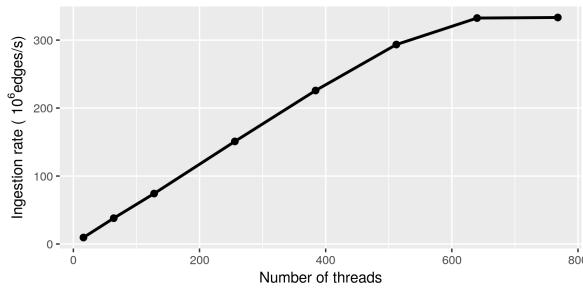


Figure 3: LANDSCAPE ingestion rate scales to one-fourth of sequential RAM bandwidth.

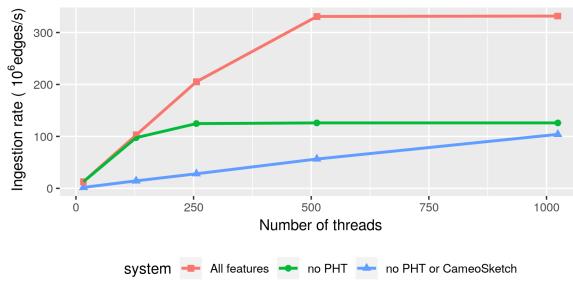


Figure 4: CAMEOSKETCH and pipeline hypertree are vital for good ingestion performance. Without CAMEOSKETCH, LANDSCAPE’s ingestion rate scales slowly as the number of threads increases. Without pipeline hypertree, the system bottlenecks at slightly over 100 million updates/sec.

workers. Figure 3 demonstrates a near-linear increase in ingestion rate as more distributed threads are added, until throughput levels off as it approaches 340 million updates/sec. The ingestion rate on one worker node (with 16 threads) is about 9.6 million updates/sec, and the ingestion rate on 40 worker nodes (with a total of  $40 \cdot 16 = 640$  threads) is 332 million updates/sec, a  $35\times$  speedup. We note that at this point LANDSCAPE is observably **not** CPU-bound: adding more worker nodes no longer increases throughput, and we measure instructions per cycle on the main node CPU to be 0.8, indicating that the CPU is not instruction bound, but RAM bandwidth bound [28].

Since graph stream updates are 9 bytes, LANDSCAPE ingestion bandwidth (2.8 GiB/sec) is four times that of random-access RAM bandwidth (759 MiB/sec), and roughly one-fourth of sequential RAM access bandwidth (12.4 GiB/sec) on the (c5n.18xlarge) main node.

Throughout these tests we observe a constant amount of network communication used; approximately 1.6 times the size of the input stream for dense graphs. See Table 5 for a more complete evaluation.

**Performance impact of CameoSketch and pipeline hypertree.** Figure 4 illustrates the performance im-

Table 3: LANDSCAPE has a high ingestion rate on sufficiently dense graphs and has low communication overhead.

Dataset	Ingestion Rate ( $\times 10^6$ updates/sec)	Communication (as a factor of stream size)
<b>kron13</b>	231	1.6
<b>kron15</b>	336	1.6
<b>kron16</b>	334	1.6
<b>kron17</b>	335	1.6
<b>ca-citeseer</b>	17.3	1.7
<b>p2p-gnutella</b>	13.5	0
<b>rec-amazon</b>	12.5	0
<b>google-plus</b>	134	2.6
<b>web-uk-2005</b>	91.5	3.4
<b>erdos18</b>	291	1.6
<b>erdos19</b>	226	1.6
<b>erdos20</b>	236	1.6

pact of CAMEOSKETCH and the pipeline hypertree on LANDSCAPE’s ingestion rate with the **kron17** dataset. When LANDSCAPE uses GRAPHZEPPELIN’s buffering data structure and its CUBESKETCH sketch algorithm, its ingestion rate increases as more distributed workers are added, but at a slow rate. Using CAMEOSKETCH alongside GRAPHZEPPELIN’s buffering system results in a much faster increase in ingestion rate, but the system bottlenecks at roughly 120 million updates/sec. In contrast, the full LANDSCAPE system continues to increase its ingestion rate dramatically as more workers are added, and bottlenecks at over 300 million updates/sec. We conclude that the  $O(\log V)$  decrease in ingestion cost for CAMEOSKETCH and the improved design of the pipeline hypertree are vital for LANDSCAPE’s performance. See Appendix F.4 for a direct experimental comparison of LANDSCAPE to GRAPHZEPPELIN on a single machine.

**More datasets.** Table 3 summarizes LANDSCAPE’s stream ingestion rate and network costs using 640 worker threads on a variety of synthetic and real-world datasets. Its ingestion rate is very high on dense graph streams and on real-world streams **google-plus** and **web-uk-2005**. It is lower on **ca-citeseer**, **p2p-gnutella** and **rec-amazon** because these datasets do not contain enough stream updates to surpass LANDSCAPE’s 4% leaf fullness threshold. So, for these streams, a large portion (or all) of update processing occurs on the main node.

**Circumventing bottlenecks.** These experiments support our claims that sketching can avoid the traditional bottlenecks in distributed graph stream processing. They demonstrate that CPU cost can be distributed away, supporting Claim 12 that network communication can be only a constant factor of the data acquisition cost, supporting Claim 13, and that stream processing

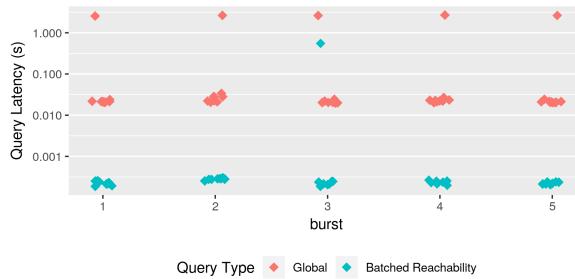


Figure 5: GREEDYCC dramatically decreases query latency.

bandwidth can reach to within a factor 4 of sequential RAM bandwidth, supporting Claim 14.

**7.3 Landscape Answers Queries Quickly.** We measured LANDSCAPE’s query latency for both global connectivity queries and batched reachability queries on an extended kron17 stream. For batched reachability queries, we uniformly sample the pairs to query  $(v_1, v_2), (v_3, v_4), \dots$  from the set of all vertices. We issue queries periodically and record (1) the time to flush the pipeline hypertree and update the graph sketch and (2) to perform Boruvka’s algorithm using the graph sketch. The sum of these times is the total query latency, though only the Boruvka computation is additional work induced by the query; the flushing work would have happened anyway as part of stream ingestion even if the query was never issued. We found that flushing takes roughly 2.3 seconds, while Boruvka’s algorithm takes 0.3 seconds.

We also evaluate how query latency and ingestion rate are affected by the use of GREEDYCC. We measure the performance impact of this optimization by running LANDSCAPE on extended kron17 and issuing *query bursts*: multiple connectivity queries issued close together (within 50 thousand stream updates). We track the latency of each query in the burst. The results are summarized in Figure 5. The first query in each burst, which requires flushing and running Boruvka’s algorithm to answer, has high (multi-second) latency. However, we see that for the remaining queries in the burst the latency is much lower: two orders of magnitude lower for global connectivity queries, and up to four orders of magnitude lower for batched reachability queries.

**7.4  $k$ -connectivity Performance** We repeat our scaling, network communication, and query latency experiments on the  $k$ -connectivity problem for a variety of datasets and values of  $k$ . Table 4 highlights LANDSCAPE’s performance for computing  $k$ -connectivity on the kron17 dataset. Note that increasing  $k$  incurs a linear decrease in ingestion rate, a linear increase in sketch size, a quadratic increase in query

Table 4: When computing  $k$ -connectivity, increasing  $k$  incurs a linear decrease in ingestion rate, a linear increase in sketch size, a quadratic increase in query latency, and does not affect total network communication. Reported values obtained by running LANDSCAPE on the kron17 dataset.

	Ingestion Rate ( $10^6$ u/s)	Sketch Size (GiB)	Query (seconds)	Network (GiB)
$k = 1$	338.5	15.25	1.27	425.2
$k = 2$	200	24.40	5.02412	464.981
$k = 4$	101.5	46.49	16.1608	468.886
$k = 8$	50.76	83.39	65.5716	476.598

latency, and has no significant effect on network communication. These experimental results support the asymptotic conclusions in Theorem 5.4. See Appendix F.3 for a summary of LANDSCAPE’s performance on more datasets.

## 8 Conclusion

This paper demonstrates how to use linear sketching to avoid the space, CPU & network bottlenecks faced by existing graph processing systems. We make our argument in the context of the LANDSCAPE system for finding connected components of dense, dynamic graphs.

By avoiding these bottlenecks, LANDSCAPE achieves remarkable performance. Specifically, it supports a stream ingestion rate one-fourth of sequential RAM bandwidth.

## 9 Appendix

The text of the appendix can be found in the full version of this paper<sup>2</sup>.

## Acknowledgments

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