

On the Depth of Oblivious Parallel RAM

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Abstract

Oblivious Parallel RAM (OPRAM), first proposed by Boyle, Chung, and Pass, is the natural parallel extension of Oblivious RAM (ORAM). OPRAM provides a powerful cryptographic building block for hiding the access patterns of programs to sensitive data, while preserving the parallelism inherent in the original program. All prior OPRAM schemes adopt a single metric of “simulation overhead” that characterizes the blowup in parallel runtime, assuming that oblivious simulation is constrained to using the *same* number of CPUs as the original PRAM.

In this paper, we ask whether oblivious simulation of PRAM programs can be further sped up if the OPRAM is allowed to have *more* CPUs than the original PRAM. We thus initiate a study to understand the true depth of OPRAM schemes (i.e., when the OPRAM may have access to unbounded number of CPUs). On the upper bound front, we construct a new OPRAM scheme that gains a logarithmic factor in depth and without incurring extra blowup in total work in comparison with the state-of-the-art OPRAM scheme. On the lower bound side, we demonstrate fundamental limits on the depth any OPRAM scheme — even when the OPRAM is allowed to have an unbounded number of CPUs and blow up total work arbitrarily. We further show that our upper bound result is optimal in depth for a reasonably large parameter regime that is of particular interest in practice.

Keywords: Oblivious Parallel RAM, depth complexity, parallel algorithms

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1 Introduction

Oblivious RAM (ORAM), originally proposed in the seminal works of Goldreich and Ostrovsky [13, 14], is a powerful cryptographic building block that allows a program to hide access patterns to sensitive data. Since Goldreich and Ostrovsky’s ground-breaking results, numerous subsequent works showed improved ORAM constructions [16, 19, 26, 28, 29] with better asymptotics and/or practical performance. ORAM has also been used in various practical and theoretical applications such as multi-party computation [17, 30], secure processor design [11, 12, 20, 21, 24], and secure storage outsourcing [27, 31].

Since most modern computing architectures inherently support parallelism (e.g., cloud compute clusters and modern CPU designs), a natural problem is how to hide sensitive access patterns in such a parallel computing environment. In a recent seminal work, Boyle, Chung, and Pass [2] first propose the notion of Oblivious Parallel RAM (OPRAM), which is a natural extension of ORAM to the parallel setting. Since then, several subsequent works have constructed efficient OPRAM schemes [5, 6, 22]. One central question in this line of research is whether there is an OPRAM scheme whose *simulation overhead* matches that of the best known ORAM scheme. Specifically, an OPRAM scheme with simulation overhead X means that if the original PRAM consumes m CPUs and runs in parallel time T , then we can obliviously simulate this PRAM also with m CPUs, and in parallel runtime $X \cdot T$. In a recent companion paper called Circuit OPRAM [5], we answered this question in the affirmative. In particular, if N is the number of distinct blocks that the CPUs can request, then Circuit OPRAM proposed a unifying framework where we can obtain statistically secure OPRAMs with $O(\log^2 N)$ simulation overhead, and computationally secure OPRAMs with $(\log^2 N / \log \log N)$ simulation overhead — thus matching the best known ORAM schemes in both settings [19, 29].

All previous OPRAM schemes consider a single performance metric referred to as simulation overhead as mentioned above. It is immediate that an OPRAM scheme with X simulation overhead also immediately implies an ORAM construction with X simulation overhead. Thus, the recent Circuit OPRAM [5] also suggests that we have hit some road-block for constructing more efficient OPRAM schemes — unless we knew how to asymptotically improve the efficiency of sequential ORAM. Note also that in the regime of sufficiently large block sizes, Circuit OPRAM achieves $O(\alpha \log N)$ simulation overhead for any super-constant function α , and this is (almost) tight in light of Goldreich and Ostrovsky’s logarithmic ORAM lower bound [13, 14].

1.1 Our Results and Contributions

In this paper, we rethink the performance metrics for an OPRAM scheme. We argue that while adopting a single simulation overhead metric is intuitive, this single metric fails to capture the true “work-depth” of the oblivious simulation. In particular, we ask the questions:

1. *If the OPRAM is allowed to access more CPUs than the original PRAM, can we have oblivious simulations with smaller parallel runtime blowup than existing OPRAM schemes?*
2. *Are there any fundamental limits to an OPRAM’s work-depth, assuming that the OPRAM can have access to an unbounded number of CPUs?*

To answer the above questions, we turn to the parallel algorithms literature, and adopt two classical metrics, that is, *total work* and *parallel runtime* in the study of OPRAMs. Like the parallel algorithms literature, we also refer to a(n) PRAM/OPRAM’s parallel runtime as its *work-depth* (or *depth*). The depth metric represents the runtime of a PRAM when given ample CPUs — thus the depth is the inherently sequential part of a PRAM that cannot be further parallelized even

with an arbitrarily large number of CPUs. The depth metric is commonly used in conjunction with total work — since we would like to design low-depth parallel algorithms that do not blow up total work by too much in comparison with the sequential setting (e.g., by repeating computations too many times). Using these classical metrics from the parallel algorithms literature, we can re-interpret the single “simulation overhead” metric adopted by previous OPRAM works as follows: an OPRAM with simulation overhead X has both total work blowup and parallel runtime blowup X in comparison with the original PRAM.

Note that when the OPRAM is constrained to using the same number of CPUs as the original PRAM, its parallel runtime blowup must be at least as large as the total work blowup. In this paper, however, we show that this need not be the case when the OPRAM can access more CPUs than the original PRAM. We design a new OPRAM scheme that gains a logarithmic factor in speed (i.e., depth) in comparison with the state-of-the-art [5] when given logarithmically many more CPUs than the original PRAM. In some sense, our new OPRAM scheme shows that the blowup in total work incurred due to obliviousness can be parallelized further (albeit through non-trivial techniques). Additionally, we prove new lower bounds that shed light on the inherent limits on any OPRAM scheme’s depth. In light of our lower bounds, our new OPRAM scheme is optimal in depth for a wide range of parameters. We now present an informal overview of our results and contributions.

1.1.1 Upper Bounds

First, we show that for any PRAM running in time T and consuming W amount of total work, there exists a statistically secure oblivious simulation that consumes logarithmically many more CPUs than the original PRAM, and runs in parallel runtime $O(T \log N \log \log N)$ and total work $O(W \log^2 N)$.

In comparison, the best known (statistically secure) OPRAM scheme incurs both $O(\log^2 N)$ blowup in both total work and parallel runtime (i.e., $O(\log^2 N)$ simulation overhead). In this sense, while preserving the total work blowup, we improve existing OPRAMs’ depth by a logarithmic factor.

We then extend our construction to the computationally secure setting by adapting an elegant trick originally proposed by Fletcher et al. [10], and show how to shave another $\log \log N$ factor off both the total work and parallel runtime, assuming that one-way functions exist. Our results are summarized in the following informal theorem.

Theorem 1 (Small-depth OPRAMs: Informal). *The following results are possible for small-depth OPRAMs where N denotes the original PRAM’s total memory size, m denotes the original PRAM’s number of CPUs, and the security failure must be negligible in N .*

- **Statistically secure, general block size.** *There exists a statistically secure OPRAM that achieves $O(\log^2 N)$ blowup in total work and $O(\log N \log \log N)$ blowup in parallel runtime for general block sizes of $\Omega(\log N)$ bits.*
- **Computationally secure, general block size.** *Assume the existence of one-way functions, then there exists a computationally secure OPRAM that achieves $O(\frac{\log^2 N}{\log \log N})$ total work blowup and $O(\log N)$ parallel runtime blowup for general block sizes of $\Omega(\log N)$ bits.*
- **Statistically secure, large block size.** *For any super-constant function $\alpha(N) = \omega(1)$, for any constant $\epsilon > 0$, there exists a statistically secure OPRAM that achieves $O(\alpha \log N \log \log N)$ total work blowup and $O(\log m + \log \log N)$ parallel runtime blowup for blocks of N^ϵ bits or larger.*

1.1.2 Lower Bounds

Next, we consider if there are any fundamental limits to an OPRAM scheme’s work-depth. We prove a non-trivial lower bound showing that any *online* OPRAM scheme (i.e., with no a-priori knowledge of future requests) that does not perform encoding of data blocks and does not duplicate data blocks too extensively must suffer from at least $\Omega(\log m)$ depth blowup where m is the number of CPUs — and this lower bound holds even when the OPRAM scheme may access arbitrarily many CPUs and have arbitrarily large total work blowup. We stress that our lower bound employs techniques that are different in nature from those of Goldreich and Ostrovsky’s classical ORAM lower bound [13, 14] — in particular, theirs bounds total work rather than depth. Furthermore, our lower bound holds even for computational security.

Theorem 2 (Lower bound for an OPRAM’s depth). *Any computationally or statistically secure online OPRAM scheme must incur at least $\Omega(\log m)$ blowup in parallel runtime, as long as the OPRAM 1) does not perform encoding of data blocks (i.e., in the “balls-and-bins” model); and 2) does not make more than $m^{0.1}$ copies of each data block.*

We note that the conditions our lower bound assumes (online, balls-and-bins, and bounded duplication) hold for all ORAM and OPRAM constructions.

1.1.3 On the Tightness of Our Upper and Lower Bounds

In light of our lower bound, our OPRAM constructions are optimal in depth in a reasonably large parameter regime. Specifically, our (computationally secure) OPRAM scheme is depth-optimal when $m = N^\epsilon$ for any constant $\epsilon > 0$ for general block sizes. For larger block sizes, our OPRAM scheme is depth-optimal for a larger range of m — in particular, when the block size is sufficiently large, our (statistically secure) OPRAM scheme is tight for m as small as $m = \text{poly log } N$.

1.1.4 Technical Highlights

Both our lower bounds and upper bounds introduce non-trivial new techniques. Since our lower bound studies the depth of parallel algorithms, it is of a very different nature than Goldreich and Ostrovsky’s ORAM lower bounds for total work [13, 14]. To prove the depth lower bound, we also depart significantly in technique from Goldreich and Ostrovsky [13, 14]. In particular, our lower bound is of an *online* nature and considers the possible batches of requests that a low-depth access pattern can support in a single PRAM step; whereas in comparison, Goldreich and Ostrovsky’s lower bound applies even to offline ORAM/OPRAM algorithms, and they perform a counting argument over many steps of the ORAM/OPRAM. The most difficult challenge in proving our lower bound is how to offset the large number of possibilities introduced by “preprocessing”, i.e., the number of possible memory configurations before the PRAM step of concern starts. To deal with this challenge, our core idea is to devise a new method of counting that is *agnostic to preprocessing*.

For our new small-depth OPRAM, the main challenge we cope with is of a very different nature from known ORAM and OPRAM works. In particular, all previous ORAMs and OPRAMs that follow the tree-based paradigm [26] adopt a standard recursion technique such that the CPU need not store a large amount of metadata (referred to as the position map). Known schemes treat this recursion as a blackbox technique. Unfortunately, in our work, it turns out that this recursion becomes the main limiting factor to an OPRAM’s depth. Thus, we open up the recursion, and our core technique for achieving small-depth OPRAM is to devise a novel offline/online paradigm, such that the online phase that is inherently sequential across recursion levels has small (i.e., $O(\log \log N)$) depth per recursion level; whereas all work that incurs logarithmic depth is performed

in an offline phase in parallel across all recursion levels. Designing such an offline/online algorithm incurs several challenges which we explain in Section 5.2. We hope that these new techniques can also lend to the design of oblivious parallel algorithms in general.

Another way to view our small-depth OPRAM’s contributions is the following. In our setting, we must address two challenges: 1) concurrency, i.e., how to coordinate a batch of m requests such that they can be served simultaneously without causing write conflicts; and 2) parallelism, i.e., how to make each request parallel by using more CPUs. Note that the concurrency aspect is applicable only to OPRAMs where multiple concurrent requests are involved, whereas the parallelism aspect is applicable even for parallelizing the operations of a sequential ORAM. Previous OPRAM constructions [2, 6] are concerned only about the former concurrency aspect, but we need to take both into account — in this sense, we are in fact the *first* to investigate the “parallelism” aspect of ORAMs/OPRAMs.¹ In particular, in our fetch phase algorithm, the two aspects are intertwined for the case of general m , in the sense that we cannot separate our techniques into two phases involving one “concurrent compilation” and one “parallel compilation” — such intertwining allows us to construct more efficient algorithms. In the maintain phase, our divide-and-conquer strategy for eviction indeed can be used to parallelize a sequential ORAM.

1.2 Related Work

Oblivious RAM (ORAM) was initially proposed in a ground-breaking work by Goldreich and Ostrovsky [13, 14], who showed that assuming the existence of one-way functions, any RAM-program can be obliviously simulated with $O(\alpha \log^3 N)$ overhead where α is any super constant function. Subsequently, numerous works constructed asymptotically better ORAM schemes [16, 26, 28, 29] and as well as ORAM schemes with statistical or perfect security [7, 9, 26, 28, 29]. As far as (sequential) ORAM is concerned, state-of-the-art results show how to construct 1) computationally secure ORAMs with $O(\log^2 N / \log \log N)$ simulation overhead [16, 19]; 2) statistically secure ORAM schemes with $O(\log^2 N)$ simulation overhead [29]; and 3) perfectly secure ORAM schemes with $O(\log^3 N)$ simulation overhead. All of the aforementioned results work for any block size (as long as the block is large enough to store its own memory address), as well as $O(1)$ blocks of CPU private cache. For sufficiently large block sizes, we know that ORAM schemes with $O(\alpha \log N)$ overhead can be constructed for an arbitrarily small super-constant function α — this is almost optimal in light of Goldreich and Ostrovsky’s logarithmic ORAM lower bound [13, 14].

Boyle, Chung, and Pass recently initiated the study of Oblivious Parallel RAM (OPRAM) [2]. They were also the first to phrase the simulation overhead metric for OPRAMs, i.e., the parallel runtime blowup of the OPRAM in comparison with the original PRAM, assuming that the OPRAM consumes the same number of CPUs as the original PRAM. Several subsequent works [2, 5, 6, 22] have improved Boyle et al. [2]’s OPRAM construction. Most recently, Chan and Shi [5] show that we can construct statistically secure and computationally secure OPRAMs whose asymptotical performance match the best known sequential ORAM; and their approach is based on the tree-based paradigm [26]. A similar asymptotical result (but for the computationally secure setting only) was also shown by Chan et al. [4] using the hierarchical framework originally proposed by Goldreich and Ostrovsky [13, 14]. In the OPRAM context, Goldreich and Ostrovsky’s logarithmic lower bound [13, 14] immediately implies that any OPRAM with constant blocks of CPU cache must suffer from at least logarithmic *total work* blowup. Thus far there is no other known OPRAM lower bound (and our depth lower bound departs significantly in techniques from Goldreich and Ostrovsky’s lower bound).

Besides the more standard notion of OPRAM, several other works have considered the parallel oblivious simulation of restricted forms of parallel programs [23]. Dachman-Soled [8] also consider

¹We gratefully acknowledge the Asiacrypt reviewers for pointing out this aspect of our contribution.

the parallel oblivious simulation in a special RAM model where addresses within a memory bank are not observable by the adversary, but the adversary can observe which CPUs access which memory banks.

While our upper bound results build on top of the OPRAM frameworks proposed in prior work [2, 5, 6], we show that several fundamentally new techniques are necessary to asymptotically improve the depth of OPRAMs.

2 Definitions

2.1 Parallel Random-Access Machines

A *parallel random-access machine* (PRAM) consists of a set of CPUs and a shared memory denoted mem indexed by the address space $[N] := \{1, 2, \dots, N\}$. In this paper, we refer to each memory word also as a *block*, and we use B to denote the bit-length of each block.

We use m to denote the number of CPUs. In each step t , each CPU executes a next instruction circuit denoted Π , updates its CPU state; and further, CPUs interact with memory through request instructions $\vec{I}^{(t)} := (I_i^{(t)} : i \in [m])$. Specifically, at time step t , CPU i 's instruction is of the form $I_i^{(t)} := (\text{op}, \text{addr}, \text{data})$, where the operation is $\text{op} \in \{\text{read}, \text{write}\}$ performed on the virtual memory block with address addr and block value $\text{data} \in \{0, 1\}^B \cup \{\perp\}$. If $\text{op} = \text{read}$, then we have $\text{data} = \perp$ and the CPU issuing the instruction should receive the content of block $\text{mem}[\text{addr}]$ at the initial state of step t . If $\text{op} = \text{write}$, then we have $\text{data} \neq \perp$; in this case, the CPU still receives the initial state of $\text{mem}[\text{addr}]$ in this step, and at the end of step t , the content of virtual memory $\text{mem}[\text{addr}]$ should be updated to data .

Write conflict resolution. By definition, multiple **read** operations can be executed concurrently with other operations even if they visit the same address. However, if multiple concurrent **write** operations visit the same address, a conflict resolution rule will be necessary for our PRAM be well-defined. In this paper, we assume the following:

- The original PRAM supports concurrent reads and concurrent writes (CRCW) with an arbitrary, parametrizable rule for write conflict resolution. In other words, there exists some priority rule to determine which **write** operation takes effect if there are multiple concurrent writes in some time step t .
- The compiled, oblivious PRAM (defined below) is a “concurrent read, exclusive write” PRAM (CREW). In other words, the design of our OPRAM construction must ensure that there are no concurrent writes at any time.

We note that a CRCW-PRAM with a parametrizable conflict resolution rule is among the most powerful CRCW-PRAM model, whereas CREW is a much weaker model. Our results are stronger if we allow the underlying PRAM to be more powerful but the compiled OPRAM uses a weaker PRAM model. For a detailed explanation on how stronger PRAM models can emulate weaker ones, we refer the reader to the work by Hagerup [18].

CPU-to-CPU communication. In the remainder of the paper, we sometimes describe our algorithms using CPU-to-CPU communication. For our OPRAM algorithm to be oblivious, the inter-CPU communication pattern must be oblivious too. We stress that such inter-CPU communication can be emulated using shared memory reads and writes. Therefore, when we express our performance metrics, we assume that all inter-CPU communication is implemented with shared

memory reads and writes. In this sense, our performance metrics already account for any inter-CPU communication, and there is no need to have separate metrics that characterize inter-CPU communication. In contrast, Chen et al. [6] defines separate metrics for inter-CPU communication.

Additional assumptions and notations. Henceforth, we assume that each CPU can only store $O(1)$ memory blocks. Further, we assume for simplicity that the runtime of the PRAM is *fixed* a priori and *publicly known*. Therefore, we can consider a PRAM to be a tuple

$$\text{PRAM} := (\Pi, N, m, T),$$

where Π denotes the next instruction circuit, N denotes the total memory size (in terms of number of blocks), m denotes the number of CPUs, and T denotes the PRAM's parallel time steps. *Without loss of generality, we assume that $N \geq m$.* We stress that henceforth in the paper, the notations N and m denote the number of memory blocks and the number of CPUs for the original PRAM — our OPRAM construction will consume $O(1)$ factor more memory and possibly more than m CPUs.

2.2 Oblivious Parallel Random-Access Machines

Randomized PRAM. A *randomized PRAM* is a special PRAM where the CPUs are allowed to generate private, random numbers. For simplicity, we assume that a randomized PRAM has a priori known, deterministic runtime.

Oblivious PRAM (OPRAM). A randomized PRAM parametrized with total memory size N is said to be *statistically oblivious*, iff there exists a negligible function $\epsilon(\cdot)$ such that for any inputs $x_0, x_1 \in \{0, 1\}^*$,

$$\text{Addresses}(\text{PRAM}, x_0) \stackrel{\epsilon(N)}{\equiv} \text{Addresses}(\text{PRAM}, x_1),$$

where $\text{Addresses}(\text{PRAM}, x)$ denotes the joint distribution of memory accesses made by PRAM upon input x and the notation $\stackrel{\epsilon(N)}{\equiv}$ means the statistical distance is bounded by $\epsilon(N)$. More specifically, for each time step $t \in [T]$, $\text{Addresses}(\text{PRAM}, x)$ includes the memory addresses requested by the CPUs in time step t , as well as whether each memory request is a read or write operation. Henceforth we often use the notation OPRAM to denote a PRAM that satisfies statistical obliviousness.

Similarly, a randomized PRAM parametrized with total memory size N is said to be *computationally oblivious*, iff there exists a negligible function $\epsilon(\cdot)$ such that for any inputs $x_0, x_1 \in \{0, 1\}^*$,

$$\text{Addresses}(\text{PRAM}, x_0) \stackrel{\epsilon(N)}{\equiv_c} \text{Addresses}(\text{PRAM}, x_1)$$

Note the only difference from statistical security is that here the access patterns only need to be indistinguishable to computationally bounded adversaries, denoted by the notation $\stackrel{\epsilon(N)}{\equiv_c}$.

Following the convention of most existing ORAM and OPRAM works [13, 14, 19, 28, 29], we will require that the security failure probability to be negligible in the N , i.e., the PRAM's total memory size.

Oblivious simulation. We say that a given OPRAM *simulates* a PRAM if for every input $x \in \{0, 1\}^*$, $\Pr[\text{OPRAM}(x) = \text{PRAM}(x)] = 1 - \mu(N)$ where the completeness error μ is a negligible function and the probability is taken over the randomness consumed by the OPRAM — in other words, we require that the OPRAM and PRAM output the same outcome on any input x .

Online OPRAM. In this paper we focus on *online* OPRAM that simulates a PRAM by processing memory request of each PRAM step in an online fashion. Namely, each PRAM memory request is processed by the OPRAM without knowing the future requests. Note that all known ORAM and OPRAM constructions satisfy the online property.

Performance measures. For an online OPRAM simulates a certain PRAM, we measure its performance by its *work-depth* and *total work* overhead. The work-depth overhead is defined to be the number of time steps d for OPRAM to simulate each PRAM step. Let W denote the total number of blocks accessed by OPRAM to simulate a PRAM step. The total work overhead is defined to be W/m , which captures the overhead to simulate a batch of memory request in a PRAM step. Note that both d and W are random variables.

Remark 1 (Static vs. adaptive security). Our above OPRAM definition works in the static security setting where the input is chosen upfront without having observed the earlier accesses. Later we shall prove our lower bound for static security since this makes our lower bound stronger. However, we point out that our upper bound in fact satisfies a stronger, adaptive and composable notion of security which we provide in the appendices.

3 Lower Bound on Work-Depth

We show a lower bound on the work-depth in terms of the number m of CPUs. We establish a $\Omega(\log m)$ depth lower bound for OPRAMs satisfying the following properties. We remark that our construction in Section 5 as well as all existing ORAM and OPRAM constructions satisfy these properties.

1. **Balls-and-bins storage.** As coined in the ORAM lower bound of Goldreich and Ostrovsky [3], data blocks are modeled as “balls,” while shared memory locations and CPU registers are modeled as “bins”. In particular, this means that every memory location stores at most one data block and the content of the data block can be retrieved from that location independent of other storage.
2. **Online OPRAM.** As defined in Section 2.2, we consider online OPRAM that only learns the logic memory request at the beginning of a PRAM step.
3. **s -bounded duplication.** We also need a technical condition on the bound of data duplication. Namely, there is a bound s such that every data block has at most s copies stored on the memory. All known ORAM and OPRAM constructions do not store duplications on the memory², i.e., $s = 1$.

It is worth comparing our depth lower bound for OPRAM with the ORAM lower bound of [3]. Both lower bounds assume the balls-and-bins model, but establish lower bound for different metrics and rely on very different arguments (in particular, as we discussed below, counting arguments do not work in our setting). We additionally require online and bounded duplication properties, which are not needed in [3]. On the other hand, our lower bound holds even for OPRAM with *computational security*. In contrast, the lower bound of [3] only holds for statistical security.

The setting for the lower bound. For simplicity, we consider the following setting for establishing the lower bound. First, we consider OPRAM with initialization, where n logical data blocks of the original PRAM are initialized with certain distinct content. This is not essential as we can

²In some hierarchical ORAMs [16, 19], there might be several copies of the same block on the server, but only one copy is regarded as *fresh*, while other copies are *stale* and may contain old contents.

view the initialization as the first n steps of the PRAM program. We also assume that the logical data size n is sufficiently larger than the total CPUs register size. Specifically, let α be a constant in $(0, 1/3)$ and r be the register size of a CPU. We assume $n \geq \Omega(r \cdot m^{1+(\alpha/4)})$. For any OPRAM satisfying the above three properties with $s \leq m^{(1/3)-\alpha}$, we show that the work-depth is at least $(\alpha/3) \cdot \log m$ with probability at least $1 - m^{-\alpha/4}$ for every PRAM step. In particular, the expected work-depth per step is at least $\Omega(\log m)$ as long as $s \leq m^{1/3-\Omega(1)}$.

Theorem 3 (Lower Bound on Work-Depth). *Let Π be a computationally-secure online OPRAM that satisfies the balls-and-bins model with s -bounded duplication for $s < m^{(1/3)-\alpha}$ for constant $\alpha \in (0, 1/3)$, where the number N of blocks is at least m . Let r be the register size of each CPU. Assume that $n \geq 4r \cdot m^{1+(\alpha/4)}$ and Π has correctness error $\mu \leq m^{-\alpha/4}/4$. Then for each PRAM step t , let $\text{depth}(\Pi, t)$ denote the work-depth of Π for PRAM step t ,*

$$\Pr[\text{depth}(\Pi, t) \leq (\alpha/3) \cdot \log m] \leq m^{-\alpha/4},$$

where the probability is over the randomness of the OPRAM compiler Π .

Before proving Theorem 3, we first discuss the intuition behind the lower bound proof in Section 3.1, where under simplifying assumptions, we reduce the OPRAM lower bound to solving a “user-movie problem” that captures the main argument of our lower bound proof. We discuss how to remove the simplifying assumptions in the end of the section. We then present the formal proof of Theorem 3 in Section 3.2

3.1 Intuition: A User-Movie Problem

As a warmup, we first present an intuitive proof making a few simplifying assumptions: 1) the OPRAM compiler must be perfectly correct and perfectly secure; and 2) there is no data block duplication in memory. Later in our formal proofs in Section 3.2, these assumptions will be relaxed.

Let us consider how to prove the depth lower bound for a PRAM step t for an OPRAM. Recall that we consider online OPRAM that learn the logical memory requests at the beginning of the step. We can view what happened before the step t as a preprocessing phase that stores the logical memory blocks in different memory locations, and the step t corresponds to an online phase where the CPUs fetch the requested memory blocks with certain observed access pattern. Since the access pattern should hide the logical memory request, any fixed access pattern should allow the CPUs to complete any possible batch of m requests (assuming perfect correctness and perfect security). We say that an access pattern can support a batch of m requests, if there *exists a pre-processing* (i.e., packing of data blocks into memory), such that each CPU can “reach” its desired data block through this access pattern. Our goal is to show that if the access pattern is low depth, then it is impossible to satisfy every batch of m requests — even when one is allowed to enumerate all possible pre-processings to identify one that best matches the requests (given the fixed access pattern). To show this, our argument involves two main steps.

1. First, we show that for any access pattern of low depth, say, d , each CPU can reach at most 2^d memory locations.
2. Second, we show that if an access pattern can satisfy all possible batches of m requests (with possibly different pre-processing), then it must be that some CPU can reach many physical locations in memory.

The former is relatively easy to show. Informally speaking, consider the balls-and-bins model as mentioned earlier: in every PRAM step, each CPU can only access a single memory location (although each memory location can be accessed by many CPUs). This means that at the end of the PRAM step, the block held by each CPU can only be one of two choices: 1) the block previously

held by the CPU; or 2) the block in the memory location the CPU just accessed. This means that the access pattern graph must have a small fan-in of 2 (although the fan-out may be unbounded). It is not difficult to formalize this intuition, and show that given any depth- d access pattern, only 2^d memory locations can “flow into” any given CPU. Henceforth, we focus on arguing why the latter is also true — and this requires a much more involved argument.

For ease of understanding, henceforth we shall refer to CPUs as *users*, and refer to data blocks in physical memory as *movies*. There are n distinct movies stored in a database of size N (without duplications) and m users. Each user wants to watch a movie and can access to certain 2^d locations in the database, but the locations the users access to cannot depend on the movies they want to watch. On the other hand, we can decide which location to store each movie to help the users to fetch their movies from the locations they access to. In other words, we first decide which 2^d locations each user access to, then learn which movie each user wants to watch. Then we decide the location to store each movie to help the users to fetch their movies. Is it possible to find a strategy to satisfy all possible movie requests?

We now discuss how to prove the impossibility for the user-movie problem. We first note that a simple counting argument does not work, since there are n^m possible movie requests but roughly $N^n \gg n^m$ possible ways to store the movies in physical memory. To prove the impossibility, we first observe that since we do not allow duplications, when two users request the same movie, they must have access to the same location that stores the movie. Thus, any pair of users must be able to reach a common movie location — henceforth we say that the two users “share” a movie location. This observation alone is not enough, since the users may all share some (dummy) location. If, however, two sets of users request two different movies, then not only must each set share a movie location, the two sets must share two *distinct* locations. More generally, the m users’ movie requests induce a *partition* among users where all users requesting the same movie are in the same *part* (i.e., equivalence class), and users in two different parts request different movies. This observation together with carefully chosen partitions allow us to show the existence of a user that needs to access to a large number of locations, which implies an impossibility for the user-movie problem for sufficiently small depth d . We stress that this idea of “partitioning” captures the essence of what pre-processing *cannot* help with, and this explains why our proof works even when there are a large number of possible pre-processings.

Specifically, let $k = m/2$ and label the m users with the set $M := [2] \times [k]$. We consider the following k partitions that partition the users into k pairs. For each $i \in [k]$, we define partition $P_i = \{(1, a), (2, a + i)\} : a \in [k]\}$, where the addition is performed modulo k . Note that all k^2 pairs in the k partitions are distinct. By the above observation, for each partition P_i , there are k *distinct* locations $\ell_{i,1}, \dots, \ell_{i,k} \in [N]$ such that for each pair $\{(1, a), (2, a + i)\}$ for $a \in [k]$, both users $(1, a), (2, a + i)$ access to the location $\ell_{i,a}$. Now, for each location $\ell \in [N]$, let w_ℓ denote the number of $\ell_{i,a} = \ell$ and d_ℓ denote the number of users access to the location ℓ . Note that $w_\ell \leq k$ since user pairs in a partition access to distinct locations (i.e., $\ell_{i,a} \neq \ell_{i,a'}$ for every $i \in [k]$ and $a \neq a' \in [k]$). Also note that $d_\ell \geq \sqrt{2w_\ell}$ since there are only $\binom{d_\ell}{2}$ distinct pairs of users access to the location ℓ .

To summarize, we have (i) $\sum_\ell w_\ell = k^2$, (ii) $w_\ell \leq k$ for all $\ell \in [N]$, and (iii) $d_\ell \geq \sqrt{2w_\ell}$ for all $\ell \in [N]$, which implies $\sum_\ell d_\ell \geq k \cdot \sqrt{2k} = \sqrt{k/2} \cdot m$. Recall that d_ℓ denote the number of users access to the location ℓ and there are m users. By averaging, there must exist a user who needs to access to at least $\sqrt{k/2}$ locations. Therefore, the user-movie problem is impossible for $d \leq 0.5 \cdot \log m - 2$. Note that the distinctness of the $\ell_{i,a}$ ’s induced by the partitions plays a crucial role to drive a non-trivial lower bound on the summation $\sum_\ell d_\ell$.

Removing the simplifying assumptions. In above intuitive proof we make several simplifying assumptions such as perfect security and perfect correctness. We now briefly discuss how to remove these assumptions. The main non-trivial step is to handle computational security, which requires

two additional observations. Following the above argument, let us say that an access pattern is compatible with a CPU/user partition if it can support a logic memory request with corresponding induces CPU/user partition.

- First, the above impossibility argument for the user-movie problem can be refined to show that if an access pattern has depth d , then it can be compatible with at most $2^{2(d+1)}$ partitions in P_1, \dots, P_k defined above.
- Second, whether an access pattern is compatible with a partition can be verified in polynomial time.

Based on these two observations, we show that if $d \leq 0.5 \cdot \log m - 4$ (with noticeable probability), then we can identify two efficiently distinguishable CPU partitions, which implies a depth lower bound for computationally-secure OPRAM. First, we consider the access pattern of partition P_1 . Since $d \leq 0.5 \cdot \log m - 4$, it can only be compatible with at most $k/2$ partitions. By an averaging argument, there exists some partition P_i such that P_i is not compatible with the access pattern of P_1 with probability at least $1/2$. On the other hand, by perfect correctness, the access pattern of P_i is always compatible with P_i . Therefore, the access patterns of P_1 and P_i are efficiently distinguishable by an efficient distinguisher D that simply verifies if the access pattern is compatible with P_i .

We now briefly discuss how to remove the remaining assumptions. First, it is not hard to see that the above argument does not require perfect correctness and can tolerate a small correctness error. Second, we make an implicit assumption that the requested data blocks are not stored in the CPU registers so that the CPUs must fetch the requested data blocks from physical locations on the server. This can be handled by considering logic access requests with random logical address and assuming that the logic memory size n is sufficiently larger than the total CPU register size (as in the theorem statement).

We also implicitly assume that we can observe the beginning and end of the access pattern of a PRAM step t . For this, we note that by the online property, we can without loss of generality consider t as the last step so that we know the end of the access pattern for free. Furthermore, we observe that we do not need to know the beginning of the access pattern since the compatibility property is monotone in the following sense. If a partition P_i is compatible with the access pattern of the last d accesses, it is also compatible with the access pattern of the last $d + 1$ accesses. Thus, we can consider the the access pattern of the last d accesses for certain appropriately chosen d .

Finally, to handle s -bounded duplication with $s > 1$, we consider CPU partitions where each part is a set of size $s + 1$, instead of a pair. By the pigeonhole principle, each part can still certify a pair of CPUs with a shared memory location. However, some extra care is needed for defining the partitions to make sure that different partitions do not certify the same pair of CPUs, and the depth lower bound degrades when s increases. Nevertheless, the lower bound remains $\Omega(\log m)$ for $s \leq m^{1/3 - \Omega(1)}$.

3.2 Proof of Theorem 3

We now proceed with a formal proof. We first note that for proving lower bound of the PRAM step t , we can consider PRAM programs where t is the last step, since the behavior of an online OPRAM does not depend on the future PRAM steps. Thus, we can focus on proving lower bound of the last PRAM step. We prove the theorem by contradiction. Suppose that

$$\Pr[\text{depth}(\Pi, t) \leq (\alpha/3) \cdot \log m] > m^{-\alpha/4}, \quad (1)$$

we show two PRAM programs P_1, P_2 with identical first $t - 1$ steps and different logic access request at step t such that the access pattern of $\Pi(P_1)$ and $\Pi(P_2)$, which denote the OPRAM simulation

of P_1, P_2 respectively, are efficiently distinguishable. Towards this, we define the CPU partition of a memory request.

Definition 1 (CPU Partition). Let $\text{addr} = (\text{addr}_1, \dots, \text{addr}_m) \in [n]^m$ be a memory request. addr induces a partition P on the CPUs, where two CPUs c_1, c_2 are in the same part *iff* they request for the same logical address $\text{addr}_{c_1} = \text{addr}_{c_2}$. In other words, P partitions the CPUs according to the requested logical addresses.

Recall that s is the bound on the number of duplication. We assume $m = (s + 1) \cdot k$ for some prime k . This is without loss of generality, because any integer has a prime number that is within a multiplicative factor of 2. We label the m CPUs with the set $M := [s + 1] \times [k]$. We consider the following set of partitions P_1, \dots, P_k : For $i \in [k]$, the partition $P_i := \{S_i(a) : a \in [k]\}$ is defined such that each part has the form $S_i(a) := \{(b, a + bi) : b \in [s + 1]\}$, where addition is performed modulo k . In other words, the parts in the partitions can be viewed as all possible distinct line segments in the \mathbb{Z}_k^2 plane.

We will show two programs where their last memory requests have induced partitions P_1 and P_i for some $i \in [k]$ such that their compiled access patterns are efficiently distinguishable. To show this, we model the view of the adversary with an *access pattern graph* and consider a *compatibility* property between an access pattern graph and a CPU partition, defined as follows.

Access pattern graphs and compatibility. Given the access pattern of $\Pi(P)$ for a PRAM program P and a depth parameter $d \in \mathbb{N}$, we define an access pattern graph G as follows.

- (a) **Nodes.** The nodes are partitioned into $d + 1$ *layers*. In layer 0, each node represents a physical location in the memory at the beginning of the last d -th time step of $\Pi(P)$.

For $1 \leq i \leq d$, each node in layer i represents a physical location in the memory or a CPU at the end of the last $(d - i + 1)$ -st time step.

Hence, we represent each node with (i, u) , where i is the layer number and u is either a CPU or a memory location.

- (b) **Edges.** Each edge is directed and points from a node in layer $i - 1$ to one in layer i for some $i \geq 1$. For each CPU or a memory location, there is a directed edge from its copy in layer $i - 1$ to one in layer i .

If a CPU c reads from some physical location ℓ in the last $(d - i)$ -th time step, then there is a directed edge from $(i - 1, \ell)$ to (i, c) . Since we allow concurrent read, the out-degree of a node corresponding to a physical location can be unbounded.

If a CPU c writes to some physical location ℓ in the last $(d - i)$ -th time step, then there is a directed edge from $(i - 1, c)$ to (i, ℓ) .

Observe that since we consider OPRAM with exclusive write, the in-degree of a node (either corresponding to a CPU or a memory location) is at most 2. In fact, the degree 2 bound holds even with concurrent write models as long as the write conflict resolution can be determined only by the access pattern.

The access pattern graph G captures the potential data flow of the last d time steps of the data access. Specifically, a path from $(0, \ell)$ to (d, c) means CPU c may learn the content of the memory location ℓ at the last d time step. If there is no such path, then CPU c cannot learn the content. This motivates the definition of compatible partitions.

Definition 2 (Compatible Partition). Let G be an access pattern graph and P_1, \dots, P_k be the partitions defined above. We say $P_i = \{S_i(a) : a \in [k]\}$ is compatible with G if there exist k distinct physical locations $\ell_{i,1}, \dots, \ell_{i,k}$ on the server such that for each $a \in [k]$, there are at least two CPUs c_1 and c_2 in $S_i(a)$ such that both nodes (d, c_1) and (d, c_2) are reachable from $(0, \ell_{i,a})$ in G .

Intuitively, compatibility is a necessary condition for the last d time steps of data access to “serve” an access request with induced partition P_i , assuming the requested data blocks are not stored in the CPU registers at the last d -th time step. Recall that each data block has at most s copies in the server, and each part $S_i(a)$ has size $s + 1$. By the Pigeonhole principle, for each part $S_i(a)$ in the induced partition, there must be at least two CPUs $c_1, c_2 \in S_i(a)$ obtaining the logical block from the same physical location ℓ_a on the server, which means the nodes (d, c_1) and (d, c_2) are reachable from $(0, \ell_a)$ in G . We note that verifying compatibility can be done in polynomial time.

Lemma 1 (Verifying Compatibility Takes Polynomial Time). *Given a CPU partition P and an access pattern graph G , it takes polynomial time to verify whether P is compatible with G .*

Proof. Given P and G as in the hypothesis of the lemma, we construct a bipartite graph H as follows. Each vertex in L is labeled with a memory location ℓ , and each vertex in R is labeled with a part S in P . There is an edge connecting a vertex ℓ in L to a vertex S in R iff there are at least two CPUs c_1 and c_2 in S such that both (d, c_1) and (d, c_2) are reachable from $(0, \ell)$ in G . This bipartite graph can be constructed in polynomial time.

Observe that P is compatible with G iff there is a matching in H such that all vertices in R are matched. Hence, a maximum matching algorithm can be applied to H to decide if P is compatible with G . \square

Now, the following key lemma states that an access pattern graph G with small depth d cannot be compatible with too many partitions. We will use the lemma to show two programs with efficiently distinguishable access patterns.

Lemma 2. *Let G be an access pattern graph with the depth parameter d , and P_1, \dots, P_k be the partitions defined above. Among P_1, \dots, P_k , there are at most $((s + 1) \cdot 2^d)^2$ partitions that are compatible with G .*

Proof. Recall that the in-degree of each node is at most 2. Thus, for each node (d, c) in layer d , there are at most 2^d nodes $(0, \ell)$ in layer 0 that can reach the node (d, c) . For the sake of contradiction, we show that if G is compatible with $u > ((s + 1) \cdot 2^d)^2$ partitions, then there exists a node (d, c) that is reachable by more than 2^d nodes in layer 0.

For convenience, we define a bipartite graph $H = (L, R, E)$ from G as follows. Each vertex in L is labeled with a CPU c , and each vertex in R is labeled with a physical location ℓ of the memory. There is an edge (c, ℓ) in H iff $(0, \ell)$ reaches (d, c) in G . Our goal can be restated as showing that if G is compatible with $u > ((s + 1) \cdot 2^d)^2$ partitions, then there exists $c \in L$ with degree $\deg(c) > 2^d$. We do so by lower bounding the number of edges $|E| > m \cdot 2^d$.

By definition, if P_i is compatible with G , then there exist k distinct physical locations $\ell_{i,1}, \dots, \ell_{i,k}$ on the server such that for each $a \in [k]$, there are at least two CPUs $c_{i,a}, c'_{i,a} \in S_i(a)$ such that $(d, c_{i,a})$ and $(d, c'_{i,a})$ are reachable from $(0, \ell_{i,a})$ in G , which means there are edges $(c_{i,a}, \ell_{i,a})$ and $(c'_{i,a}, \ell_{i,a})$ in H . Thus, a compatible partition certifies $2k$ edges in H , although two different partitions may certify the same edges.

Let P_{i_1}, \dots, P_{i_u} be the set of compatible partitions. While they may certify the same edges, the set of CPU pairs $\{(c_{i_j,a}, c'_{i_j,a}) : j \in [u], a \in [k]\}$ are distinct for the following reason: Recall that

the parts in partitions correspond to different line segments in \mathbb{Z}_k^2 . Since two points define a line, the fact that the parts correspond to different lines implies that all CPU pairs are distinct.

For each memory location ℓ , let w_ℓ denote the number of $\ell_{i_j, a} = \ell$. It means that ℓ is connected to w_ℓ distinct pairs of CPUs in H , which implies that $\deg(\ell) \geq \sqrt{2w_\ell}$ since there must be at least $\sqrt{2w_\ell}$ distinct CPUs. Also, note that $\sum_\ell w_\ell = u \cdot k$ and $w_\ell \leq u$ for every ℓ since ℓ can appear in each partition at most once. It is not hard to see that the above conditions imply a lower bound on $|E| = \sum_\ell \deg(\ell) \geq k \cdot \sqrt{2u} > m \cdot 2^d$. This in turn implies the existence of $c \in L$ with degree $\deg(c) > 2^d$, a contradiction. \square

Let us now consider a PRAM program P_1 that performs dummy access in the first $t - 1$ steps and a random access request at the t step with induced partition P_1 . Specifically, in the first $t - 1$ steps, all CPUs read the first logic data block. For the t -th step, let (b_1, \dots, b_k) be uniformly random k distinct logic data blocks. For $a \in [k]$, the CPUs in part $S_1(a)$ of P_1 read the block b_a at the t -th step. Let $d = (\alpha/3) \cdot \log m$ and $G(\Pi(P_1), d)$ denote the access pattern graph of $\Pi(P_1)$ with depth parameter d . The following lemma follows directly by Lemma 2 and an averaging argument.

Lemma 3. *There exists $i^* \in [k]$ such that*

$$\Pr[P_{i^*} \text{ is compatible with } G(\Pi(P_1), d)] \leq ((s+1) \cdot 2^d)^2/k \leq m^{-\alpha/3},$$

where the randomness is over Π and P_1 .

Now, consider a PRAM program P_2 that is identical to P_1 , except that the access request at the t -th step has induced partition P_{i^*} instead of P_1 . Namely, for $a \in [k]$, the CPUs in part $S_{i^*}(a)$ of P_{i^*} read the block b_a at the t -th step, where (b_1, \dots, b_k) are uniformly random k distinct logic data blocks.

Lemma 4. *Suppose that Π satisfies Eq. (1), then*

$$\Pr[P_{i^*} \text{ is compatible with } G(\Pi(P_2), d)] > m^{-\alpha/4}/4,$$

where the randomness is over Π and P_2 .

Proof. First note that since each CPU request a random data block at the t -th PRAM step, the probability that the requested data block is stored in the CPU register is at most r/n . By a union bound, with probability at least $1 - m \cdot (r/m) \geq 1 - m^{-\alpha/4}/4$, all data blocks requested at the t -th PRAM step are not in the corresponding CPU registers. In this case, the CPUs need to obtain the data blocks from the server. Furthermore, if the work-depth of the t -th PRAM step is $\leq d$, then the CPUs need to obtain the data blocks in the last d time steps of data access, which as argued above, implies compatibility. Therefore,

$$\Pr[P_{i^*} \text{ is compatible with } G(\Pi(P_2), d)] > m^{-\alpha/4} - m^{-\alpha/4}/4 - \epsilon_c > m^{-\alpha/4}/4.$$

\square

Recall by Lemma 1 that compatibility can be checked in polynomial time. The above two lemmas imply that assuming Eq. (1), $\Pi(P_1)$ and $\Pi(P_2)$ are efficiently distinguishable by a distinguisher D who checks the compatibility of P_{i^*} and the access pattern graph with depth parameter $d = (\alpha/3) \cdot \log m$. This is a contradiction and completes the proof of Theorem 3.

4 Background on Circuit OPRAM and Building Blocks

4.1 Preliminaries: Circuit OPRAM

As a warmup, we first briefly review the recent Circuit OPRAM algorithm [5] that we build on top of. For clarity, we make a few simplifying assumptions in this overview:

- We explain the non-recursive version of the algorithm where we assume that the CPU can store a position map for free that tracks the rough physical location of every block: this CPU-side position map is later removed using a standard recursion technique in Circuit OPRAM [5] — however, as we point out later, to obtain a small depth OPRAM in our paper, we must implement the recursion differently and thus in our paper we can no longer treat the recursion as blackbox technique.
- We assume that m is not too small and is at least polylogarithmic in N ; and
- A standard conflict resolution procedure proposed by Boyle et al. [2] has been executed such that the incoming batch of m requests are for distinct real blocks (or dummy requests).

Core data structure: a pool and $2m$ subtrees. Circuit OPRAM partitions the OPRAM data structure in memory into $2m$ *disjoint subtrees*. Given a batch of m memory requests (from m CPUs), each request will be served from a random subtree. On average, each subtree must serve $O(1)$ requests in a batch; and due to a simple balls and bins argument, except with negligible probability, even the worst-case subtree serves only $O(\alpha \log N)$ incoming requests for any super-constant function α .

In addition to the $2m$ subtree, Circuit OPRAM also maintains an overflow pool that stores overflowing data blocks that fail to be evicted back into the $2m$ subtrees at the end of each batch of m requests.

It will help the reader to equivalently think of the $2m$ subtrees and the pool in the following manner: First, think of a single big Circuit OPRAM [29] tree (similar to other tree-based OPRAMs [26]). Next, identify a height with $2m$ buckets, which naturally gives us $2m$ *disjoint subtrees*. All buckets from smaller heights as well as the Circuit OPRAM’s stash form the *pool*. As proven in the earlier work [5], at any time, the pool contains at most $O(m + \alpha \log N)$ blocks.

Fetch. Given a batch of m memory requests, henceforth without loss of generality, we assume that the m requests are for distinct addresses. This is because we can adopt the conflict resolution algorithm by Boyle et al. [2] to suppress duplicates, and after data has been fetched, rely on oblivious routing to send fetched data to all request CPUs. Now, look up the requested blocks in two places, both the pool and the subtrees:

- *Subtree lookup:* For a batch of m requests, each request comes with a position label — and all m position labels define m random paths in the $2m$ subtrees. We can now fetch from the m path in parallel. Since each path is $O(\log N)$ in length, each fetch can be completed in $O(\log \log N)$ parallel steps with the help of $\log N$ CPUs.

All fetched blocks are merged into the pool. Notice that at this moment, the pool size has grown by a constant factor, but later in a cleanup step, we will compress the pool back to its original size. Also, at this moment, we have not removed the requested blocks from the subtrees yet, and we will remove them later in the maintain phase.

- *Pool lookup:* At this moment, all requested blocks must be in the pool. Assuming that m is not too small, we can now rely on oblivious routing to route blocks back to each requesting CPU — and this can be completed in $O(\log m)$ parallel steps with m CPUs.

Maintain. In the maintain phase, perform the following: 1) remove all blocks fetched from the paths read; and 2) perform eviction on each subtree.

- *Efficient simultaneous removals.* After reading each subtree, we need to remove up to $\mu := O(\alpha \log N)$ blocks that are fetched. Such removal operations can lead to write contention when done in parallel: since the paths read by different CPUs overlap, up to $\mu := O(\alpha \log N)$ CPUs may try to write to the same location in the subtree. Circuit OPRAM employs a novel *simultaneous removal* algorithm to perform such removal in $O(\log N)$ parallel time with m CPUs. We refer the reader to the Circuit OPRAM paper for an exposition of the simultaneous removal algorithm. As noted in the Circuit OPRAM paper [5], simultaneous removal from m fetch paths can be accomplished in $O(\log m + \log \log N)$ parallel steps with $O(m \cdot \log N)$ total work.
- *Selection of eviction candidates and pool-to-subtree routing.* At this moment, we will select exactly one eviction candidate from the pool for each subtree. If there exists one or more blocks in the pool to be evicted to a certain subtree, then the *deepest* block (where deepest is precisely defined in Circuit ORAM [29]) with respect to the current eviction path will be chosen. Otherwise, a dummy block will be chosen for this subtree. Roughly speaking, using the above criterion as a preference rule, we can rely on oblivious routing to route the selected eviction candidate from the pool to each subtree. This can be accomplished in $O(\log m)$ parallel steps with m CPUs assuming that m is not too small.
- *Eviction.* Now, each subtree performs exactly 1 eviction. This can be accomplished in $O(\log N)$ runtime using the sequential procedure described in the original Circuit ORAM paper [29]. At the end of this step, each subtree will output an eviction leftover block: the leftover block is dummy if the chosen eviction candidate was successfully evicted into the subtree (or if the eviction candidate was dummy to start with); otherwise the leftover block is the original eviction candidate. All these eviction leftovers will be merged back into the central pool.
- *Pool cleanup.* Notice that in the process of serving a batch of requests, the pool size has grown — however, blocks that have entered the pool may be dummy. In particular, we shall prove that the pool’s occupancy will never exceed $c \cdot m + \alpha \log N$ for an appropriate constant c except with $\text{negl}(N)$ probability. Therefore, at the end of the maintain phase, we must compress the pool back to $c \cdot m + \alpha \log N$. Such compression can easily be achieved through oblivious sorting in $O(\log m)$ parallel steps with m CPUs, assuming that m is not too small.

Recursion. Thus far, we have assumed that the position map is stored on the CPU-side, such that the CPU knows where every block is in physical memory. To get rid of the position map, Circuit OPRAM employs a standard recursion technique that comes with the tree-based ORAM/OPRAM framework [26]. At a high level, the idea of the recursion framework is very simple: instead of storing the position map on the CPU side, we recurse and store the position map in a smaller OPRAM in physical memory; and then we recurse again and store the position map of this smaller OPRAM in a yet smaller OPRAM in physical memory, and so on. If each block can store $\gamma > 1$ number of position labels, then every time we recurse, the OPRAM’s size reduces by a factor of γ . Thus in at most $\log N$ recursion levels, the metadata size becomes at most $O(1)$ blocks — and at this moment, the CPU can store all the metadata locally in cache.

Although most prior tree-based ORAM/OPRAM papers typically treat this recursion as a standard, blackbox technique, in this paper we cannot — on the contrary, it turns out that the recursion becomes the most non-trivial part of our low-depth OPRAM algorithm. Thus, henceforth the reader will need to think of the recursion in an expanded form — we now explain what exactly happens in the recursion in an expanded form. Imagine that one of the memory requests among the batch of m requests asks for the logical address $(0101100)_2$ in binary format, and suppose that

each block can store 2 position labels. Henceforth we focus on what happens for fetching this logical address $(0101100)_2$ — but please keep in mind that there are m such addresses and thus the following process is repeated m times in parallel.

- First, the 0th recursion level (of constant size) will tell the 1st recursion level the position label for the address $(0*)_2$.
- Next, the 1st recursion level fetch the metadata block at level-1 address $(0*)_2$ and this fetched block contains the position labels for $(00*)_2$ and $(01*)_2$.
- Now, level-1 informs level-2 of the position label for $(01*)_2$; at this moment, level-2 fetches the metadata block for the level-2 address $(01*)_2$ and this fetched block contains the position labels for the addresses $(010*)_2$ and $(011*)_2$; and so on.
- This continues until the D -th recursion level (i.e., the final recursion level) — this final recursion level stores actual data blocks rather than metadata, and thus the desired data block will be fetched at the end.

As mentioned, the above steps are in fact replicated m times in parallel since there are m requests in a batch. This introduces a couple additional subtleties:

- First, notice that for obliviousness, conflict resolution must be performed upfront for each recursion level before the above procedure starts — this step can be parallelized across all recursion levels.
- Second, how do the m fetch CPUs at one recursion level *obliviously* route the fetched position labels to the m fetch CPUs waiting in the next recursion level? Circuit OPRAM relies on a standard oblivious routing procedure (initially described by Boyle et al. [2]) for this purpose, thus completely hiding which CPUs route to which.

Important observation. At this moment, we make an important observation. In the Circuit OPRAM algorithm, the fetch phase operations are inherently sequential across all recursion levels, and the maintain phase operations can be parallelized across all recursion levels. In particular, during the fetch phase, the m fetch CPUs at recursion level d must block waiting for recursion level $d - 1$ to pass down the fetched position labels before its own operations can begin. Due to the sequential nature of the fetch phase, Circuit OPRAM incurs at least $(\log m + \log \log N) \log N$ depth, where the $\log m$ stems from level-to-level oblivious routing, $\log \log N$ stems the depth needed to parallel-fetch from a path of length $\log N$ (and other operations), and the $\log N$ factor is due to the number of recursion levels. In comparison, the depth of the maintain phase is not the limiting factor due to the ability to perform the operations in parallel across recursion levels.

4.2 Other Important Building Blocks

Permutation-related building blocks. We will rely on the following building blocks related to generating and applying permutations.

To describe our ideas, we need a few basic building blocks as depicted in Figures 1, 2, and 3.

1. **Apply a pre-determined permutation to an array.** Figure 1 shows how to in parallel apply a pre-determined permutation to an array in a single parallel step.
2. **Permute an array by a secret random permutation.** Figure 2 shows how to generate a secret random permutation and apply it to an array obliviously, without revealing any information about the permutation. The formal abstraction for an oblivious random permutation and formal proofs for Figure 2 are deferred to the appendices.

Apply Permutation to Array

Inputs:

- An initial array $A[0..k]$ of size k .
- A permutation $\pi : [k] \rightarrow [k]$ written down as an array.

Outputs: A new array $\widehat{A}[0..k]$, where for each $i \in [k]$, $\widehat{A}[i] = A[\pi^{-1}(i)]$; in other words, each element $A[i]$ is copied to $\widehat{A}[\pi(i)]$.

Algorithm: Parallel runtime $O(1)$, total work $O(k)$

- Each CPU with index i reads $A[i]$ and $\pi(i)$, and then set $\widehat{A}[\pi(i)] := A[i]$.

Figure 1: Applying a pre-determined permutation to an array

3. **Obliviously construct a routing permutation that permutes a source to a destination array.** Figure 3 shows how to accomplish the following task: given a source array `snd` of length k containing distinct real elements and dummies (where each dummy element contains unique identifying information as well), and a destination array `rcv` also of length k containing distinct real elements and dummies, with the guarantee that the set of real elements in `snd` are the same as the set of real elements in `rcv`. Now, construct a routing permutation $\pi : [k] \rightarrow [k]$ (in an oblivious manner) such that for all $i \in [k]$, if `snd`[i] contains a real element, then `rcv`[$\pi[i]$] = `snd`[i].

Oblivious bin-packing. Oblivious bin-packing is the following primitive.

- *Inputs:* Let B denote the number of bins, and let Z denote the target bin capacity. We are given an input array denoted `ln`, where each element is either a dummy denoted \perp or a real element that is tagged with a bin number $g \in [B]$. It is guaranteed that there are at most Z elements destined for each bin.
- *Outputs:* An array `Out`[$1 : BZ$] of length $B \cdot Z$ containing real and dummy elements, such that `Out`[($g - 1$) $B + 1 : gB$] denotes contents of the g -th bin for $g \in [B]$. The output array `Out` must guarantee that the g -th bin contains all elements in the input array `ln` tagged with the bin number g ; and that all real elements in bin g must appear in the input array `ln` and are tagged with g .

There is an oblivious parallel algorithm that accomplishes oblivious bin packing in total work $O(\tilde{n} \log \tilde{n})$ and parallel runtime $O(\log \tilde{n})$ where $\tilde{n} = \max(|\text{ln}|, B \cdot Z)$. The algorithm works as follows:

1. For each group $g \in [B]$, append Z filler elements of the form (`filler`, g) to the resulting array — these filler elements ensure that every group will receive at least Z elements after the next step.
2. Obliviously sort the resulting array by the group number, placing all dummies at the end. When elements have the same group number, place filler elements after real elements.
3. By invoking an instance of the oblivious aggregation algorithm [2, 23] (see Section A.4 for the definition of oblivious aggregation), each element in the array finds the leftmost element in its own group. Now for each element in the array, if its offset within its own group is greater than Z , replace the element with a dummy \perp .
4. Obliviously sort the resulting array placing all dummies at the end. Truncate the resulting array and preserve only first $B \cdot Z$ blocks.
5. For every filler element in the resulting array, replace it with a dummy.

Oblivious Random Permutation

Input: An array A of size k . We assume that $\log^{0.5} N \leq k \leq N$, since if $k \leq \log^{0.5} N$, we can simply run a naïve quadratic oblivious random permutation algorithm (see Appendix A.3).

Outputs: An array \hat{A} which is obtained by applying a uniformly random permutation on array A in an oblivious manner.

Algorithm: Except with negligible in N probability, the parallel runtime is $O(\log k)$ and the total work is $O(k \log k)$.

1. Each CPU with index $i \in [k]$ copies $A[i]$ and samples $3 \log_2 N$ bits which form a key k_i in $[3N^3]$.
2. The k CPUs perform deterministic oblivious sort (e.g., Zigzag sort [15]) on the data tuple $(k_i, A[i])$ using k_i 's to determine the total order — henceforth this array is referred to as Y .
3. Each CPU contacts its neighbors to check if the element it holds has a key colliding with neighboring elements.
4. Perform parallel prefix sum (which can easily be achieved in $O(\log k)$ depth and $O(k)$ total work) such that each CPU learns how many CPUs before it have a collided element — this determines the position at which the CPU is going to write out its collided element.
If a CPU holds a collided element, write it out to a small colliding array at the position that was computed in the above step.
5. Now, perform naïve quadratic random permutation (see Appendix A.3) to permute the colliding array. At the end, the CPU that wrote the i -th element of the colliding array grabs back the i -th element of the permuted array and writes it back into the result array Y . Finally, output Y .

Figure 2: Randomly permuting an array obliviously, without revealing the secret permutation.

5 A Small-Depth OPRAM: Level-to-Level Routing Algorithm

5.1 Overview of Our OPRAM

We now show how we can improve the depth of OPRAM schemes [2] by a logarithmic factor, through employing the help of more CPUs; and importantly, we achieve this without incurring extra total work in comparison with the best known OPRAM scheme [5].

Challenges. As argued earlier in Section 4.1, for the case of general block sizes, the most sequential part of the Circuit OPRAM algorithm stems from the (up to) $\log N$ recursion levels. More specifically, (apart from the final data level), each recursion level's job is to fetch the metadata (referred to as position labels) necessary, and route this information to the next recursion level. In this way, the next recursion level will know where in physical memory to look for the metadata needed by its next recursion level, and so on (we refer the reader to Section 4.1 for a more detailed exposition of the recursion).

Thus, the fetch phase operations of Circuit OPRAM are inherently sequential among the D recursion levels, incurring $(D(\log m + \log \log N))$ in depth, where the $\log m$ term stems from the level-to-level oblivious routing of fetched metadata, and the $\log \log N$ term stems from fetching

Construct a Permutation that Maps a Source to a Destination

Inputs:

- A source array `snd` of length k containing distinct real elements and dummies;
- A destination array `rcv` also of length k containing distinct real elements and dummies;

Assume:

- The real elements in `snd` are guaranteed to be the same as the real elements in `rcv`.
- The dummy elements in `snd` each has a unique label that decides their relative order; similarly, the dummies in `rcv` each has a unique label too.

Outputs: The routing permutation π such that $\text{rcv}[\pi(i)] = \text{snd}[i]$ for all $i \in [k]$.

Algorithm: Parallel runtime $O(\log k)$, total work $O(k \log k)$

1. Tag each element in both arrays with an index within the array. More specifically, in parallel, write down the following two arrays:

$$[(1, \text{snd}[1]), (2, \text{snd}[2]), \dots, (k, \text{snd}[k])], \quad [(1, \text{rcv}[1]), (2, \text{rcv}[2]), \dots, (k, \text{rcv}[k])]$$

2. Obviously sort each of the above two arrays by the second value. Suppose the two output arrays are the following where “_” denotes a wildcard value that we do not care about.

$$[(s_1, -), (s_2, -), \dots, (s_k, -)], \quad [(t_1, -), (t_2, -), \dots, (t_k, -)]$$

Now in parallel write down $[(s_1 \rightarrow t_1), (s_2 \rightarrow t_2), \dots, (s_k \rightarrow t_k)]$

3. Now sort the above output by the s_i values, and let the output be

$$[(1 \rightarrow t'_1), (2 \rightarrow t'_2), \dots, (k \rightarrow t'_k)]$$

The output routing permutation is defined as $\pi(i) := t'_i$.

Figure 3: Obviously construct a routing permutation that maps a source to a destination.

metadata blocks from a path of length $\log N$. Ignoring the $\log \log N$ term, our goal therefore is to get rid of the $\log m$ depth that stems from level-to-level oblivious routing.

Our result. Our main contribution is to devise a low-depth algorithm to perform level-to-level routing of metadata. At first sight, this task seems unlikely to be successful — since each recursion level must *obviously* route its metadata to the next level, it would seem like we are inherently subject to the depth necessary for an oblivious routing algorithm [2]. Since oblivious routing in some sense implies oblivious sorting, it would seem like we have to devise an oblivious sorting algorithm of less than logarithmic depth to succeed in our goal.

Perhaps somewhat surprisingly, we show that this need not be the case. In particular, we show that by 1) allowing a negligible statistical failure probability; 2) exploiting special structures of our routing problem; and 3) introducing an offline/online paradigm for designing parallel oblivious algorithms, we can devise a special-purpose level-to-level oblivious routing algorithm such that

1. all work that is inherently $\log m$ in depth is pushed to an offline phase that can be parallelized across all recursion levels; and

2. during the online phase that is inherently sequential among all $\log N$ recursion levels, we can limit the work-depth of each recursion level to only $\log \log N$ rather than $\log m$ — note that for most interesting parameter regimes that we care about, $\log m \gg \log \log N$.

The details of this algorithm are rather involved and entail the novel usage of algorithmic building blocks such as oblivious random permutation in an offline/online paradigm. We defer the detailed introduction of this algorithm and its proofs to later in this section.

As a result, we obtain a new, *statistically* secure OPRAM algorithm (for general block sizes) that achieves $O(\log N \log \log N)$ depth blowup and $O(\log^2 N)$ total work blowup. In comparison, under our new performance metrics, the best known OPRAM algorithm [5] achieves $O(\log^2 N)$ total work blowup and $O(\log^2 N)$ depth blowup. Thus we achieve a logarithmic factor improvement in terms of depth.

Extensions. We consider several extensions. First, using a standard technique described by Fletcher et al. [10] and extended to the OPRAM setting by Chan et al. [5], we show how to obtain a *computationally* secure OPRAM scheme with $O(\log^2 N / \log \log N)$ total work blowup and $O(\log N)$ depth blowup, and supporting general block sizes. In light of our aforementioned OPRAM depth lower bound (which also applies to computationally secure OPRAMs), our OPRAM scheme is optimal for $m = N^\epsilon$ where $\epsilon > 0$ is an arbitrarily small constant.

Finally, we consider a setting with sufficiently large blocks, say, the block size is N^ϵ for any constant $\epsilon > 0$ — in this case, the recursion depth becomes $O(1)$. In this case, the limiting factor to an OPRAM’s work depth now is the eviction algorithm (rather than the level-to-level routing). We show how to leverage a non-trivial devise and conquer technique to devise a new, small-depth eviction algorithm, allowing us to perform eviction along a path of length $\log N$ in $\log \log N$ depth rather than $\log N$ — however, this is achieved at the cost of a small $\log \log N$ blowup in total work. As a result, we show that for sufficiently large blocks, there is an OPRAM scheme with depth as small as $O(\log \log N + \log m)$ where the $\log \log N$ part arises from our low-depth eviction algorithm (and other operations), and the $\log m$ part arises from the conflict resolution and oblivious routing of fetched data back to requesting CPUs — thus tightly matching our depth lower bound as long as m is at least logarithmic in N .

5.2 Small-Depth Routing of Position Identifiers: Intuition

Problem statement. As we explained earlier, in each recursion level, m fetch CPUs fetch the metadata (i.e., position labels) required for the next recursion level. The next recursion level contains m fetch CPUs waiting to receive these position labels, before its own operations can begin. Circuit OPRAM performs such level-to-level routing using a standard oblivious routing building block, thus incurring at least $D \log m$ depth where D is the number of recursion levels which can be as large as $\log N$, and $\log m$ is the depth of standard oblivious routing. How can we reduce the depth necessary for level-to-level routing?

We will first clarify some details of the problem setup. Recall that in each PRAM step, we receive a batch of m memory requests, i.e., m logical addresses. Given these m logical addresses, we immediately know which level- d addresses to fetch for each recursion level d (see Section 4.1 for details). We assume that conflict resolution has been performed for each recursion level d on all of the m level- d addresses, and thus, every real (i.e., non-dummy) level- d address is distinct. Now, note that from all these level- d addresses (and even without fetching the actual metadata in each recursion level), we can already determine the routing topology from level to level: as an example, a level-2 CPU that needs to fetch the level-2 address (010*) would like to receive position labels from the level-1 fetch CPU with the address (01*).

Our goal here is to improve the OPRAM’s depth to $O(\log N \log \log N)$ for general (worst-case) block sizes. We use the parameter Γ to denote the number of position labels that a block can store; we let $\gamma := \min\{\Gamma, m\}$ be an upper bound on the number of position labels in a block that is “useful” for the next recursion level. To achieve this, in the part of the algorithm that is sequential among all recursion levels (henceforth also referred to as the *online* part), we can only afford $O(\log \log N)$ depth rather than the $\log m$ necessary for oblivious routing. Indeed, for a general oblivious routing problem consisting of m senders and m receivers, it appears the best one can do is to rely on an oblivious routing network [2,6] that has $\log m$ depth — so how can we do better here? We rely on two crucial insights:

1. First, we observe that our routing problem has *small fan-in and fan-out*: each sender has at most γ recipients; and each recipient wants to receive from at most 1 sender. This is because that each fetched metadata block contains at most γ position labels, and obviously each fetch CPU in the next level only needs one position label to proceed.
2. Second, we will rely on an *offline-online paradigm* — in the offline phase, we are allowed to perform preparation work that indeed costs $\log m$ depth; however, in the online phase, the depth is kept to be small. Later when we employ this offline/online oblivious routing building block in our full OPRAM algorithm, we will show that the offline phase does not depend on any fetched data, and thus can be parallelized across all recursion levels, whereas the online phase must still be sequential — but recall that the online phase has much smaller depth.

First insight: localized routing. Our first idea is to rely on this observation to restrict oblivious routing to happen only within small groups — as we shall explain later, for this idea to work, it is essential that our routing problem has small fan-in and fan-out. More specifically, we would like that each small group of senders talk to a corresponding small group of receivers, say, sender group S_i talks only to receiver group R_i , where both S_i and R_i are $\mu := \alpha\gamma^2 \log N$ in size, where the choice of μ is due to Lemma 5. If we do this, then oblivious routing within each small group costs only $\log \mu$ depth.

How can we arrange senders and receivers into such small groups? For correctness we must guarantee that for every i , each receiver in R_i will be able to obtain its desired item from some sender in S_i .

To achieve this, we rely on a randomized load balancing approach. The idea is very simple. First, we pad the sender array with dummy senders to a size of $2m$ — recall that there are at most m real senders. Similarly, we pad the receiver array to a size of $2m$ as well. Henceforth if a receiver wants an item from a sender, we say that the sender and receiver are connected. Every dummy sender is obviously connected to 0 receivers.

Now, if we pick a random sender from the sender array, in expectation this sender will be connected to 0.5 receivers. Thus a random subset of μ senders will in expectation is connected to 0.5μ receivers — using measure concentration techniques, it is not difficult to show that a random subset of μ senders is connected to μ receivers except with negligible probability — note that this measure concentration result holds only when our routing problem has small fan-in and fan-out (see Lemma 5 for details).

Our idea is to randomly permute the source array, and have the first μ sender be group 1, the second μ senders be group 2, and so on. By relying on $O(1)$ number of oblivious sorts, we can now arrange the receiver array to be “loosely aligned” with the sender array, i.e., all receivers connected to sender group 1 are in the first size- μ bucket of the receiver array, all receivers connected to sender group 2 are in the second size- μ bucket of the receiver array, and so on.

Using the above idea, the good news is that oblivious routing is now constrained to μ -sized groups (each containing γ addresses), thus costing only $\log \mu$ depth. However, our above algorithm

still involves randomly permuting the sender array and oblivious routing to loosely align the receiver array with the sender array — these steps cost $\log m$ depth. Thus our idea is to perform these steps in an offline phase that can be parallelized across all recursion levels, and thus the depth does not blow up by the number of recursion levels. Nonetheless how to instantiate this offline/online idea is non-trivial as we explain below.

Second insight: online/offline paradigm. One challenge that arises is how to coordinate among all recursion levels. To help the reader understand the problem, let us first describe what would have happened if everything were performed online, sequentially level by level:

Imagine that each recursion has $2m$ fetch CPUs (among which at most m are real) first acting as receivers. Once these receivers have received the position labels, they will fetch data from the OPRAM’s tree data structure. At this point, they hold the position labels desired by the next recursion level, and thus the receivers now switch roles and become senders with respect to the next recursion level. Before the receivers become senders, it is important that they be randomly permuted for our earlier load balancing technique to work. Now, we can go ahead and prepare the next recursion level’s receivers to be loosely aligned with the permuted senders, and proceed with the localized oblivious routing.

Now let us consider how to divide this algorithm into a parallel offline phase and a subsequent low-depth online phase. Clearly, the oblivious routing necessary for loosely aligning each recursion level’s receivers with the last level’s senders must be performed in the offline phase — and we must parallelize this step among all recursion levels. Thus, our idea is the following:

- First, for each recursion level d in parallel, we randomly permute level d ’s fetch CPUs in an oblivious fashion (using a building block called oblivious random permutation), at the end of which we have specified the configuration of level d ’s sender array (that is, after level d ’s fetch CPUs switch roles and become senders).
- At this point, each recursion level d can prepare its receiver array based on the configuration of level $(d - 1)$ ’s sender array. This can be done in parallel too.
- During the online phase, after fetching metadata from the OPRAM tree, the receivers must permute themselves to switch role to senders — since the offline stage has already dictated the sender array’s configuration, this permutation step must respect the offline stage’s decision.

To achieve this in small online depth, our idea is that during the offline phase, each recursion level relies on an instance of oblivious routing to figure out exactly what permutation to apply (henceforth called the “routing permutation”) to switch the receiver array to the sender array’s configuration — and this can be done in parallel among all recursion levels once a level’s receiver and sender arrays have both been determined. Once the offline stage has written down this routing permutation, in the online stage, the receivers can simply apply the permutation, i.e., each receiver writes itself to some array location as specified by the permutation that offline stage has written down. Applying the permutation online takes a single parallel step.

One observation is that during the online stage, the routing permutation is revealed in the clear. To see why this does not leak information, it suffices to see that the result of this routing permutation, i.e., the sender array, was obliviously randomly permuted to start with (using a building block called oblivious random permutation). Thus, even conditioned on having observed the oblivious random permutation’s access patterns, each permutation is still equally likely — and thus the routing permutation that is revealed is indistinguishable from a random permutation (even when conditioned on having observed the oblivious random permutation’s access patterns).

5.3 Core Subroutine: Localized Routing

Notations and informal explanation. In the OPRAM’s execution, the instructions waiting to receive position labels at a recursion level d is denoted $\text{Instr}^{(d)}$. $\text{Instr}^{(d)}$ has been obliviously and randomly permuted in the offline phase. When these incomplete instructions have received position labels, they become complete and are now called $\text{CInstr}^{(d)}$ where $\text{CInstr}^{(d)}$ and $\text{Instr}^{(d)}$ are arranged in the same order. When data blocks are fetched in recursion level d , they are called $\text{Fetched}^{(d)}$, and $\text{Fetched}^{(d)}$ has the same order as $\text{CInstr}^{(d)}$. In the offline phase, $\text{Instr}^{(d)}$ is obliviously sorted to be loosely aligned with $\text{Fetched}^{(d-1)}$ resulting in $\overline{\text{Instr}}^{(d)}$, such that $\overline{\text{Instr}}^{(d)}$ can receive position labels from $\text{Fetched}^{(d-1)}$ through localized oblivious routing. The offline phase also prepares a routing permutation $\pi^{d \rightarrow d+1}$, that will permute $\overline{\text{Instr}}^{(d)}$ (after having received position labels) back to $\text{CInstr}^{(d)}$ — and the online phase will apply this routing permutation $\pi^{d \rightarrow d+1}$ in a single parallel step. We now describe our algorithms more formally.

We consider the following problem where there is a source array and a destination array, and the destination array wants to receive position identifiers from the source. Specifically, the source array is a set of fetched blocks in randomly permuted order, where each block may contain up to γ position labels corresponding to γ addresses in the next recursion level. The destination array is an incomplete instruction array where each element contains the address of the block to be read at the next recursion level — and each address must receive its corresponding position label before the fetch operations at the next recursion level can be invoked.

- *Inputs:* The inputs contain a randomly permuted source array $\text{Fetched}^{(d)}$ that represent the fetched position identifier blocks at recursion level d , and a randomly permuted destination array $\text{Instr}^{(d+1)}$ which represents the incomplete instruction array at recursion level $d + 1$.
 - The source array $\text{Fetched}^{(d)}$ contains $2m$ blocks, each of which contains up to γ (logical) pairs of the form $(\text{addr}, \text{pos})$ that are needed in the next recursion level. All the γ addresses in the same block comes from Γ contiguous addresses, and thus in reality the address storage is actually compressed — however, we think of each block in $\text{Fetched}^{(d)}$ as *logically* containing pairs of the form $(\text{addr}, \text{pos})$.
 - The destination array $\text{Instr}^{(d+1)}$ contains m elements each of which is of the form $(\text{addr}, _)$, where “ $_$ ” denotes a placeholder for receiving the position identifier for addr later. This array $\text{Instr}^{(d+1)}$ is also referred to as the incomplete instruction array.
 - We assume that
 - (1) all addresses in the destination array must occur in the source array;
 - (2) the γ addresses contained in the same block come from Γ contiguous addresses; and
 - (3) both the source array $\text{Fetched}^{(d)}$ and the destination array $\text{Instr}^{(d+1)}$ have been randomly permuted.
- *Outputs:* A complete instruction array denoted CInstr of length $2m$ where $\text{CInstr}^{(d+1)}[i]$ is of the form $(\text{addr}_i, \text{pos}_i)$ such that
 - $\text{Instr}^{(d+1)}[i] = (\text{addr}_i, _)$, i.e., the sequence of addresses contained in the output $\text{CInstr}^{(d+1)}$ agree with those contained in the input $\text{Instr}^{(d+1)}$; and
 - The tuple $(\text{addr}_i, \text{pos}_i)$ exists in some block in $\text{Fetched}^{(d)}$, i.e., the position identifier addr_i receives is correct (as defined by $\text{Fetched}^{(d)}$).

Offline phase. The inputs are the same as the above. In the offline phase, we aim to output the following arrays:

- a) A permuted destination array $\overline{\text{Instr}}^{(d+1)}$ that is a permutation of $\text{Instr}^{(d+1)}$ such that it is *somewhat aligned* with the source $\text{Fetched}^{(d)}$, where *somewhat aligned* means the following:

[Somewhat aligned:] Fix $\alpha := \omega(1)$ to be any super-constant function. For each consecutive $\mu := \alpha\gamma^2 \log N$ contiguous source blocks denoted $\text{Fetched}^{(d)}[k\mu+1 : (k+1)\mu]$, there is a segment of μ contiguous destination blocks $\overline{\text{Instr}}^{(d+1)}[k\mu+1 : (k+1)\mu]$ such that all addresses in $\text{Instr}^{(d+1)}$ that are contained in $\text{Fetched}^{(d)}[k\mu+1 : (k+1)\mu]$ appear in the range $\overline{\text{Instr}}^{(d+1)}[k\mu+1 : (k+1)\mu]$.

- b) A routing permutation $\pi^{d \rightarrow d+1} : [2m] \rightarrow [2m]$.

In other words, the goal of the offline phase is to prepare the source and the destination arrays such that in the online phase, we only perform oblivious routing from every $\mu := \alpha\gamma^2 \log N$ blocks (each containing at most γ labels) in the source to every μ tuples in the destination where $\alpha = \omega(1)$ is any super-constant function. This way, the online phase has $O(\log \mu)$ parallel runtime.

Before explaining how to accomplish the above, we first prove that if the source array, i.e., $\text{Fetched}^{(d)}$ has been randomly permuted, then every μ contiguous blocks contain at most μ position identifiers needed by the destination.

Lemma 5. *Let arr denote an array of $2m$ randomly permuted blocks, each of which contains γ items such that out of the $2m \cdot \gamma$ items, at most m are real and the rest are dummy.*

Then, for any consecutive n blocks in arr , with probability at least $1 - \exp(-\frac{n}{2\gamma^2})$, the number of real items contained in them is at most n .

Proof of Lemma 5. We use the Hoeffding Inequality for sampling without replacement [25]. Consider $2m$ integers a_1, a_2, \dots, a_{2m} , where the integer a_i denotes the number of real items contained in the i th pair in the array. Hence, we have $\sum_{i=1}^{2m} a_i \leq m$ and each $a_i \in [0, \gamma]$.

Since the pairs are randomly permuted in the initial array, each group of consecutive n blocks can be viewed as a sample of size n without replacement from the $2m$ integers. Let Z be the sum of the integers in the size- n sample. Then, $E[Z] = \frac{n}{2m} \sum_{i=1}^{2m} a_i \leq \frac{n}{2}$.

Hence, the Hoeffding inequality for sampling without replacement [25] gives the following as required:

$$\Pr[Z > n] \leq \Pr[Z - E[Z] > \frac{n}{2}] \leq \exp\left\{-\frac{2(\frac{n}{2})^2}{n \cdot \gamma^2}\right\} = \exp\left(-\frac{n}{2\gamma^2}\right)$$

□

We now explain the offline algorithm, i.e., permute the destination array to be somewhat aligned with the source array such that localized oblivious routing will be sufficient. We describe a parallel oblivious algorithm that completes in $O(m \log m)$ total work and $O(\log m)$ parallel runtime.

1. For each block in $\text{Fetched}^{(d)}$, write down a tuple $(\text{minaddr}, \text{maxaddr}, i)$ where minaddr is the minimum address contained in the block, maxaddr is the maximum address contained in the block, and i is the offset of the block within the $\text{Fetched}^{(d)}$ array.

Henceforth we refer to the resulting array as SrcMeta .

2. Imagine that the resulting array SrcMeta and the destination array $\text{Instr}^{(d+1)}$ are concatenated. Now, oblivious sort this concatenated array such that each metadata tuple $(\text{minaddr}, \text{maxaddr}, i) \in \text{SrcMeta}$ is immediately followed by all tuples from $\text{Instr}^{(d+1)}$ whose addresses are contained within the range $[\text{minaddr}, \text{maxaddr}]$.

3. Relying on a parallel oblivious aggregate operation [2, 23] (see Section A.4 for the definition), each element in the array (resulting from the above step) learns the first metadata tuple

$(\text{minaddr}, \text{maxaddr}, i)$ to its left. In this way, each address will learn which block (i.e., i) within $\text{Fetched}^{(d)}$ it will receive its position identifier from.

The result of this step is an array such that each metadata tuple of the $(\text{minaddr}, \text{maxaddr}, i)$ is replaced with a dummy entry \perp , and each address addr is replaced with (addr, i) , denoting that the address addr will receive its position identifier from the i -th block of $\text{Fetched}^{(d)}$.

4. For each non-dummy entry in the above array, tag the entry with a group number $\lfloor \frac{i}{\mu} \rfloor$. For each dummy entry, tag it with \perp .
5. Invoke an instance of the oblivious bin packing algorithm and pack the resulting array into $\lceil \frac{2m}{\mu} \rceil$ bins of capacity μ each. We refer to the resulting array as $\overline{\text{Instr}}^{(d+1)}$.
6. Obviously compute the routing permutation $\pi^{d \rightarrow d+1}$ that maps $\overline{\text{Instr}}^{(d+1)}$ to $\text{Instr}^{(d+1)}$.
7. Output $\overline{\text{Instr}}^{(d+1)}$ and $\pi^{d \rightarrow d+1}$.

Online phase. The online phase consists of the following steps:

1. For every k , fork an instance of the oblivious routing algorithm such that $\overline{\text{Instr}}^{(d+1)}[k\mu + 1 : (k+1)\mu]$ will receive its position identifiers from $\text{Fetched}^{(d)}[k\mu + 1 : (k+1)\mu]$.
This completes in $O(m \log \mu)$ total work and $O(\log \mu)$ parallel runtime.
2. Apply the routing permutation $\pi^{d \rightarrow d+1}$ to $\overline{\text{Instr}}^{(d+1)}$, and output the result as $\text{CInstr}^{(d+1)}$.

5.4 Level-to-Level Routing

Given our core localized routing building block, the full level-to-level position identifier routing algorithm is straightforward to state.

Offline phase. Upon receiving a batch of m memory requests, for each recursion level d in parallel:

- Truncate the addresses to the first d bits and perform conflict resolution. The result is an array of length m containing distinct addresses and dummies to read from recursion level d .
- Randomly permute the resulting array, and obtain an incomplete instruction array $\text{Instr}^{(d)}$. It is important for security that the random permutation is performed obliviously such that no information is leaked to the adversary about the permutation.

For $d = 0$, additionally fill in the position map identifiers and complete the instruction array to obtain $\text{CInstr}^{(0)}$.

- From the $\text{Instr}^{(d)}$ array, construct a corresponding incomplete $\text{Fetched}^{(d)}$ array where all position identifier fields are left blank as “_”. The blocks in $\text{Fetched}^{(d)}$ are ordered in the same way as $\text{Instr}^{(d)}$.
- If d is not the data level, fork an instance of the localized routing algorithm with input arrays $\text{Fetched}^{(d)}$ and $\text{Instr}^{(d+1)}$, and output a permuted version of $\text{Instr}^{(d+1)}$ denoted $\overline{\text{Instr}}^{(d+1)}$ a routing permutation $\pi^{d \rightarrow d+1}$.

Online phase. From each recursion level $d = 0, 1, \dots, D$ sequentially where $D = O(\frac{\log N}{\log \Gamma})$ is the total number of recursion levels:

- Based on the completed instruction $\text{CInstr}^{(d)}$, allocate an appropriate number of processors for each completed instruction and perform the fetch phase of the OPRAM algorithm. The result is a fetched array $\text{Fetched}^{(d)}$.
- Execute the online phase of the localized routing algorithm for recursion level d with the inputs $\text{Fetched}^{(d)}$, $\overline{\text{Instr}}^{(d+1)}$, and $\pi^{d \rightarrow d+1}$. The result is a completed instruction array $\text{CInstr}^{(d+1)}$ for the next recursion level.

6 Putting it Altogether: Detailed OPRAM Algorithm

We now describe our full OPRAM scheme in detail.

6.1 Our OPRAM’s Data Structure

We use the same data OPRAM structure as in earlier works [2, 6] but with slightly different parametrizations. We describe the data structures needed below.

Disjoint subtrees. In each of the D recursion levels (recursion will be described later), there are $2m$ disjoint subtrees — the subtrees can be viewed as truncating a big Circuit ORAM [29] tree at a height with $2m$ buckets, such that only the larger (i.e., closer to leaf) heights are preserved. (The reason we use $2m$ subtrees is that each requested block will lead to two path evictions later.) Just like in Circuit ORAM [29], each node in the subtree is a bucket of $O(1)$ capacity, storing $O(1)$ number of real or dummy blocks.

Overflowing pool. For the overflowing pool, we distinguish between two cases:

- When $m \geq \alpha \log \log N$: in this case, each recursion level has its own pool. Chan et al. [5] prove that the pool occupancy is upper bounded by $O(m + \alpha \log N)$ except with negligible probability.
- When $m < \alpha \log \log N$: in this case, all recursion levels share a single pool (henceforth referred to as the *shared pool*). Chan et al. [5] prove that except with negligible probability, this shared pool contains at most $O(mD + \alpha \log N)$ blocks, where D is the number of recursion levels.

Path invariant, position map, and recursion levels. As in all known tree-based ORAMs and OPRAMs [2, 6, 7, 26, 28, 29], the main invariant is that a block (in a specific recursion level) is assigned to a random path: this means that the block can reside anywhere along this tree path, or in the pool — in the case of small m , the block may reside in the shared pool.

If one did not care about CPU cache size, one can use a position map to keep track of the path identifier of each block. To achieve constant CPU cache, known tree-based ORAMs and OPRAMs [2, 6, 7, 26, 28, 29] rely on a standard recursion technique to recursively store the position map in a smaller and smaller ORAM/OPRAM.

6.2 Complete Description of Our OPRAM Algorithm

We next outline our small-depth OPRAM algorithm based on Circuit OPRAM [29]. At a very high level, our algorithm largely follows the blueprint established in the Circuit OPRAM paper. But as explained earlier, several non-trivial modifications are necessary to achieve asymptotically smaller

depth. Thus our presentation below also roughly follows the Circuit OPRAM framework (with parameters adjusted to our needs). We highlight the non-trivial steps (i.e., our contributions) in shaded boxes for clarity. The detailed algorithms for these shaded boxes are presented in Section 5.2 and Section 7.1 respectively.

Fetch phase. The fetch phase has an array of m addresses as input denoted $(\text{addr}_1, \dots, \text{addr}_m)$. Recall that there are $2m$ disjoint subtrees.

(i) *Preparation: all recursion levels in parallel.* For all recursion levels $d := 0, 1, \dots, D$ in parallel, perform the following:

- *Generate level- d prefix addresses.* Write down the level- d prefixes of all m requests addresses $(\text{addr}_1, \dots, \text{addr}_m)$. Clearly, this step can be accomplished in $O(1)$ parallel step with m CPUs.
- *Conflict resolution.* Given a list of m possibly dummy level- d addresses denoted $(\text{addr}_1^{(d)}, \dots, \text{addr}_m^{(d)})$, we run an instance of the oblivious conflict resolution algorithm to suppress duplicate requests (and pad the resulting array with dummies). This step can be accomplished in $O(\log m)$ parallel steps with m CPUs.
- *Choose fresh position labels for the next recursion level.* Let $\text{Addr}^{(d)} := \{\text{addr}_i^{(d)}\}_{i \in [m]}$ denote the list of level- d addresses after conflict resolution. By jointly examining $\text{Addr}^{(d)}$ and $\text{Addr}^{(d+1)}$, recursion level d learns for each non-dummy $\text{addr}_i^{(d)} \in \text{Addr}^{(d)}$, which of its children are needed for the next recursion level. For any child that is needed, recursion level d chooses a new position label for the next recursion level. For recursion level d , the result of this step is an instruction array

$$\{\text{addr}_i^{(d)}, (\text{npos}_j : j \in [\gamma])\}_{i \in [m]}$$

where npos_j is a fresh random label in level $d+1$ if $\text{addr}_i^{(d)} || j$ is needed in the next recursion level, otherwise $\text{npos}_j := \perp$. As described by Chan *et al.* [5], this can be accomplished in $O(\log m)$ parallel steps using m CPUs.

- *Pool lookup.* We have m CPUs each of which now seeks to fetch the level- d block at address $\text{addr}^{(d)}$. The m CPUs first tries to fetch the desired blocks inside the central pool; and at the end, the fetched blocks will be marked as dummy in the pool.
 - When $m \geq \alpha \log \log N$: we rely on an instance of the oblivious routing algorithm, such that each of these m CPUs will attempt to receive the desired block from the pool, and moreover, the received blocks will be removed from the pool. Since the pool size is bounded by $O(m + \alpha \log N)$ except with negligible probability as shown by Chan *et al.* [5], this step can be accomplished in $O(\log m + \log \log N)$ parallel steps consuming $(m \log N)$ total work.
 - When $m < \alpha \log \log N$: in this case, all recursion levels share a single pool of capacity $O(\alpha \log N + Dm)$ where D is the number of recursion levels. We rely on a single instance of oblivious routing to route answers to all requests over all recursion levels in $O(\log m + \log \log N) = O(\log \log N)$ parallel steps and $O((Dm + \alpha \log N)(\log m + \log \log N)) = O(mD \log \log N + \alpha \log N \log \log N)$ total work.
- *Offline preparation for small-depth routing.* Perform additional offline preparation that is necessary to route position identifiers across recursion levels in small online depth — this part of the algorithm was not there in the earlier Circuit OPRAM [5] since they did not aim to achieve small depth. Our new offline preparation completes in $O(\log m)$ parallel steps and $O(m \log m)$ total work per recursion level (see Section 5.2 for details).

- (ii) *Fetch: one recursion at a time sequentially.* Now, for each recursion level, m CPUs will each look for a block in one of the subtrees. This step must be performed sequentially one recursion level at a time since each recursion level must receive the position labels from the previous level before looking for blocks in the subtrees.

For each recursion level $d = 0, 1, \dots, D$ in sequential order, we perform the following:

- *Receive position labels from previous recursion level.* Unless $d = 0$ in which case the position labels can be fetched in $O(1)$ parallel step, each of the m level- d addresses will receive a pair of position labels from the previous recursion level denoted $(\mathbf{pos}, \mathbf{npos})$, where \mathbf{pos} represents the tree path to look for the desired block, and \mathbf{npos} denotes a freshly chosen label to be assigned to the block after the fetch is complete. In the Circuit OPRAM work [5], this step is accomplished through oblivious routing in $O(\log m)$ parallel steps with m CPUs — however, this approach will be too expensive for us, since over all levels of recursion, this would cost $O(\log m \log N)$ parallel steps. Therefore, in Section 5.2, we describe a new, small-online-depth algorithm for performing this level-to-level position identifier routing — this also turns out to be the most technically non-trivial part in terms of achieving a small-depth OPRAM.
 - *Subtree lookup.* We describe a small-depth variant of the subtree lookup algorithm described in [5]. At this moment, $O(m)$ fetch CPUs each receives an instruction of the form $(\mathbf{addr}^{(d)}, \mathbf{pos})$ that could be possibly dummy. Each fetch CPU will now recruit the help from $O(\log N)$ auxiliary CPUs. In parallel, each auxiliary CPU reads one physical slot on the tree path leading to the leaf node numbered \mathbf{pos} , in search of the block with logical address $\mathbf{addr}^{(d)}$ (but without removing the block). Then, the $O(\log N)$ CPUs for each fetch path invoke an instance of the oblivious select algorithm defined in Section A.4), such that at the end, the fetch CPU learns the block it is looking for, as well as the physical slot in which the block resides — and this information will later be used as input to the simultaneous removal algorithm that is part of the maintain phase. If a fetch CPU receives a dummy instruction, then the fetch CPU and its $O(\log N)$ auxiliary CPUs will simply scan through a random path in a random subtree. Clearly, the above can be accomplished in $O(\log \log N)$ parallel steps and $O(m \log N)$ total work over all $O(m)$ tree paths.
 - At this moment, each of the m CPUs has fetched the desired block either from the pool or the tree path (or the CPU has fetched dummy if it received a dummy instruction to start with). The fetched results (as well as the new position labels chosen for the next recursion level) are ready to be routed to the next recursion level.
- (iii) *Oblivious multicast: once per batch of requests.* Finally, when the data-OPRAM has fetched all requested blocks, we rely on a standard oblivious routing algorithm (see Section A.4) to route the resulting blocks to the request CPUs. This step takes $O(\log m)$ parallel time with m CPUs.

Maintain phase. All of the following steps are performed in parallel across all recursion levels $d = 0, 1, \dots, D$:

- (i) *Simultaneous removal of fetched blocks from subtrees.* After each of the m CPUs fetches its desired block from m tree paths, they perform a simultaneous removal procedure to remove the fetched blocks from the tree paths. Chan et al. [5] describes an algorithm that accomplishes this task in $O(\log N)$ parallel steps with m CPUs. It is not hard to observe that if we may have more

CPUs, the same algorithm can be completed in $O(\log m + \log \log N)$ parallel steps consuming $O(m \log N)$ total work.

- (ii) *Passing updated blocks to the pool.* Each CPU updates the contents of the fetched block — if the block belongs to a position map level, the block’s content should now store the new position labels (for the next recursion level) chosen earlier in the preparation phase. Further, each block will be tagged with a new position label that indicates where the block can now reside in the current recursion level — this position label was received earlier from the previous recursion level during the fetch phase (recall that each recursion level chooses position labels for the next recursion level).

The updated blocks are merged into the pool. The pool temporarily increases its capacity to hold these extra blocks, but the extra memory will be released at the end of the maintain phase during a cleanup operation.

This step can be accomplished in $O(1)$ parallel steps with m CPUs.

- (iii) *Selection of eviction candidates.* Following the deterministic, reverse-lexicographical order eviction strategy of Circuit ORAM [29], we choose the next $2m$ eviction paths (pretending that all subtrees are part of the same big ORAM tree). The $2m$ eviction paths will go through $2m$ subtrees henceforth referred to as *evicting subtrees*. If m has decreased (by a factor of 2) since the last PRAM step, then not all subtrees are evicting subtrees.

Our goal here is to output one (possibly dummy) block to evict for each evicting subtree, as well as the remainder of the pool (with these selected blocks removed). The block selected for each evicting subtree is based on the *deepest* criterion with respect to the current eviction path.

- When $m \geq \alpha \log \log N$: we rely on oblivious routing to accomplish this in $O(\log m + \log \log N)$ parallel steps and $O(m \log N)$ total work.
- When $m < \alpha \log \log N$: in this case, all recursion levels share a single pool containing at most $O(Dm + \alpha \log N)$ real blocks. We rely on a single instance of oblivious routing to route eviction candidates for all subtrees in all recursion levels $O(\log \log N)$ parallel steps and $O(mD \log \log N + \alpha \log N \log \log N)$ total work.

- (iv)

Eviction into subtrees. In parallel, for each evicting subtree, the eviction algorithm of Circuit ORAM [29] is performed for the candidate block the subtree has received. The straightforward strategy takes $O(\log N)$ parallel steps consuming m CPUs. In Section 7.1, we propose a new algorithm that accomplishes the eviction in small depth.

- (v) *Return eviction leftovers to pool.* After the eviction algorithm completes, if the candidate block fails to be evicted into the subtree, it will be returned to the pool; otherwise if the candidate block successfully evicts into the subtree, a dummy block is returned to the pool.

- (vi) *Cleanup.* Finally, since the pool size has grown in the above process, we perform a compression procedure to remove dummy blocks and compress the pool back to $c \cdot m + \alpha \log N$ size for an appropriate constant c . A standard measure concentration argument as used in [5] can be used to show that the pool occupancy is bounded by $c \cdot m + \alpha \log N$ except with $\text{negl}(N)$ probability, and thus ensures that no real blocks are lost during this reconstruction with all but negligible probability. Again, we handle the case of large m and small m separately:

- When $m \geq \alpha \log \log N$: in this case, we rely on oblivious sorting to accomplish pool compression in $O(\log m + \log \log N)$ parallel steps consuming $O(m \log N)$ total work.

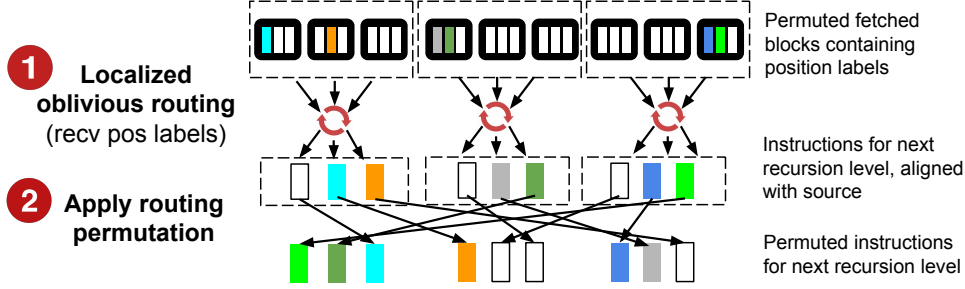


Figure 4: Online phase of the level-to-level routing algorithm

- When $m < \alpha \log \log N$: in this case, all recursion levels share a single pool containing at most $O(Dm + \alpha \log N)$ real blocks. We rely on a single instance of oblivious sorting (over all recursion levels) to perform pool compression in $O(\log \log N)$ parallel steps and $O(mD \log \log N + \alpha \log N \log \log N)$ total work.

Lemma 6 (Performance Analysis of Fetch Phase). *Suppose each block can store γ position labels. The number of recursion levels is $D = O(\frac{\log N}{\log \Gamma})$, and denote $\gamma := \min\{\Gamma, m\}$ and $\mu := \alpha \gamma^2 \log N$. Then, the fetch phase completes in $O(Dm \cdot (\log m + \log \mu))$ total work and $O(D \log \mu + \log m)$ parallel runtime.*

6.3 Obliviousness

We now argue why our scheme satisfies obliviousness. For simplicity, we may assume that whenever our OPRAM has any overflow (which happens with negligible probability), rather than aborting, the block that overflows is simply dropped — later if ever a fetch request did not find the desired block, the block’s value is treated as $\bar{0}$. In other words, henceforth we assume that the statistical failure is suffered in terms of correctness error — in this case, we will reason why the observable access patterns are identically distributed regardless of the input request sequences.

The security of our OPRAM scheme is based on that for all known tree-based ORAMs [26, 28, 29] and tree-based OPRAMs [2, 6], because whenever a block is requested, a new random path for the block is chosen without revealing the position identifier. Thus, in any recursion level, regardless of what next batch of m blocks are requested, the adversary will always observe m random paths being accessed. Hence, it suffices to show that the localized routing procedure in the fetch phase is also secure.

Lemma 7 (Security of Position Identifiers Routing). *The level-to-level routing procedure described in Section 5.4 produces a distribution of physical access pattern that is independent of the requested addresses.*

Proof. It suffices to check that in the core subroutine described in Section 5.3, the physical memory are accessed using the building blocks described in Section 4.2, which ensure that the access pattern is independent of the requested addresses. We next inspect each step more carefully.

In the offline phase, both the source array $\text{Fetched}^{(d)}$ and the destination array $\text{Instr}^{(d+1)}$ have been randomly permuted in an oblivious manner. Therefore, the routing permutation $\pi^{d \rightarrow d+1}$ (that can be observed by the adversary later in the online phase) is a uniformly random permutation, even when conditioned on having observed the access patterns of the oblivious random permutation in the offline phase — note that this is implied by our formal definition of oblivious random permutation

(see Definition 4). Other steps in the offline phase invoke oblivious building blocks such as oblivious aggregation, oblivious sort and oblivious bin packing, all of which produces deterministic access pattern independent of the requested addresses. The probability statement for the step involving oblivious bin packing refers to the probability that no real blocks are lost, but in any case, the access pattern produced is independent of the requested addresses.

In the online phase, (localized) oblivious routing is used and the access pattern produced is deterministic and independent of the requested addresses. As mentioned earlier, the routing permutation $\pi^{d \rightarrow d+1}$ is revealed, but it has an independent uniform distribution, because the destination array $\text{Instr}^{(d+1)}$ was permuted using a (secret) fresh random permutation. \square

With Lemma 7 and combining the standard security argument of tree-based ORAM and OPRAM schemes (where we can argue that for every batch of m requests, each recursion level accesses m random paths [5]), it is not difficult to see that our earlier small-depth OPRAM construction is indeed a statistically secure OPRAM, i.e., the following Theorem 4 holds.

Theorem 4 (Statistically secure, small-depth OPRAM). *There exists a statistically secure OPRAM scheme (for general block sizes) with $O(\log^2 N)$ total work blowup, and $O(\log N \log \log N)$ parallel runtime blowup, where the OPRAM consumes only $O(1)$ blocks of CPU private cache.*

6.4 Improving a $\log \log N$ Factor with Computational Security

Fletcher et al. [10] first proposed an elegant idea for transforming any *statistically secure*, tree-based ORAM into a *computationally secure* tree-based ORAM whose simulation overhead is a $\log \log N$ factor smaller than the original statistically secure scheme. Subsequently, Chan et al. [5] showed how to extend and improve their idea to the OPRAM context. Thus, given their *statistically secure* $O(\log^2 N)$ -overhead Circuit OPRAM construction, they show how to construct a *computationally secure* variant of the scheme with only $O(\log^2 N / \log \log N)$ simulation overhead — where simulation overhead is defined in the traditional way, i.e., the parallel runtime blowup of the OPRAM when the OPRAM is allowed to access only as many CPUs as the original PRAM.

At a high level, Fletcher et al.’s [10] and Chan et al.’s idea [5] is to compress the position map blocks using counters, and then generate the position labels on-the-fly pseudorandomly using a PRF whose secret key is known only to the CPU(s). In this way, when the block size is only $\Theta(\log N)$, we can pack $\Theta(\frac{\log N}{\log \log N})$ counters in each position map block, and therefore this reduces the depth of recursion by a $\log \log N$ factor. Although the high-level idea is simple, as Fletcher et al. and Chan et al. describe, several tricks are necessary to instantiate this idea. In particular, all counters in the same position map block must share the same outer counter, and each counter then has its own inner counter. When any inner counter overflows, the outer counter is incremented and all inner counters reset to 0 — at this moment, all corresponding blocks in the next recursion level must be relocated such that their physical locations are still consistent with the new counter values. Fletcher et al. and Chan et al. argue that the extra overhead incurred by such relocation (resulting from outer counter reset) can be amortized over all operations such that the relocation overhead gets absorbed asymptotically.

We refer the reader to the Circuit OPRAM [5] for a detailed explanation of this PRF+counter trick in the context of OPRAM. It is not difficult to see that our low-depth level-to-level routing trick applies nonetheless to the computationally secure variant of Circuit OPRAM [5]. Thus just like in the original Circuit OPRAM work, we gain an additional $\log \log N$ factor in both total work and depth assuming the existence of one-way functions.

Therefore we conclude with the following corollary which, due to the standard techniques explained earlier, arises as a straightforward extension to our main theorem earlier.

Corollary 1 (Computationally secure, small-depth OPRAM). *Assume that one-way functions exist. Then, there exists a computationally secure OPRAM scheme (for general block sizes) with $O(\log^2 N / \log \log N)$ total work blowup and $O(\log N)$ parallel runtime blowup, where the OPRAM consumes only $O(1)$ blocks of CPU private cache.*

7 The Depth of OPRAM Schemes for Larger Block Sizes

Thus far, we have focused on the case of general block sizes, and our results so far apply to any OPRAM whose memory block is at least large enough to store its own memory address. One cause of overhead for general block sizes is due to the up to $\log N$ levels of recursion. As all tree-based ORAM/OPRAM papers argue [5, 26, 29], when the block size is large, each block can store many position labels, and thus the depth of the recursion becomes asymptotically smaller.

In this section, we consider the depth of OPRAM schemes for larger block sizes. This question can be of interest because it can shed light on how optimal the non-recursive part of our OPRAM algorithm is. In particular, in the extreme case, when the block size is sufficiently large, i.e., N^ϵ bits per block for an arbitrarily small constant $\epsilon > 1$, the depth of recursion becomes constant. Now, the non-recursive part of the OPRAM algorithm becomes the limiting factor both in terms of total work and depth. In this case, it is not hard to see that our earlier OPRAM scheme would achieve $O(\log N)$ depth and $O(\alpha \log N)$ total work for any arbitrarily small super-constant factor α . In particular, the limiting factor for both the depth and the total work is the eviction algorithm.

In this section, our goal is the following:

- a) to obtain an OPRAM with $O(\log m + \log \log N)$ depth blowup and $O(\alpha \log N)$ total work blowup for sufficiently large block sizes (Section 7.1); and
- b) to generalize all techniques we have presented so far, and obtain a general theorem that states the best possible OPRAM depth parametrized by the block size (Section 7.2).

7.1 Small-Depth Path Eviction

As mentioned above, our first goal is to obtain an OPRAM with $O(\log m + \log \log N)$ depth blowup and $O(\alpha \log N)$ total work blowup for sufficiently large block sizes — in particular, assume that the block size is N^ϵ for an arbitrarily small constant ϵ , the depth of recursion becomes constant. At this moment, the limiting factor to Circuit OPRAM’s depth is the eviction algorithm.

If we used the naïve strategy to perform eviction on a path, each path would take $O(\log N)$ sequential steps (with a single CPU) to evict. For small block sizes, recall that since the evictions were done in parallel across all recursion levels, eviction was not the depth bottleneck — rather, the $O(\log N)$ recursion was the depth bottleneck. For sufficiently large blocks, the recursion can become as small as constant in depth, and thus the depth bottleneck is now the eviction. Thus in this section, we show how to perform eviction on each path in $O(\log \log N)$ parallel steps consuming $O(\log N \log \log N)$ total work. Henceforth in this section, let $L := O(\log N)$ denote the path length.

Intuition. Our key insight is to adopt a non-trivial divide and conquer strategy. To solve a problem instance of length n , we divide the problem into two sub-instances each of length $\frac{n}{2}$ and solve them in parallel. Then, in $O(1)$ parallel steps and $O(n)$ time, we reconstruct the solution to the length- n problem from the solutions of the two length- $\frac{n}{2}$ sub-problems.

The non-trivial challenge is that the two smaller instances are not independent. After solving each smaller instance, some information must be passed from one instance to another, and some correction must be performed on the receiving instance to make sure that the result is correct. Therefore in our algorithm design, we must consider what is the minimal amount of information

Algorithm 1 `EvictSlow(path)` *A slow, non-oblivious version of the eviction algorithm, only for illustration purpose [29]*

```

1:  $i := L$       /* start from leaf */
2: while  $i \geq 1$  do:
3:   if  $\text{path}[i]$  has empty slot then
4:      $(\text{block}, \ell) :=$  Deepest block in  $\text{path}[0..i-1]$  that can legally reside in  $\text{path}[i]$ , and its
       corresponding height in  $\text{path}$ .
       /*  $\text{block} := \perp$  if such a block does not exist. */
5:   if  $\text{block} \neq \perp$  then
6:     Move block from  $\text{path}[\ell]$  to  $\text{path}[i]$ .
7:      $i := \ell$     // skip to height  $\ell$ 
8:   else  $i := i - 1$ 

```

that must be passed from one completed instance to another, and how to perform the necessary a-posteriori correction fast, such that we can obtain a low-depth eviction algorithm while incurring little cost to total work.

Notation. We first revisit the eviction strategy used in Circuit ORAM, which is performed on a path of buckets $\text{path}[0..L]$, where $\text{path}[0]$ represents the (group) stash and $\text{path}[L]$ is the leaf. In our case, $\text{path}[1]$ corresponds to the root of one of the \hat{m} subtrees. Recall that each block has a leaf position, whose lowest common ancestor with $\text{path}[L]$ gives the deepest height at which the block can legally reside on path . We assume there is a total ordering imposed on the blocks determined by the maximum depths they can legally reside in path , where ties are consistently resolved (for instance by block addresses). The (non-oblivious) procedure in Algorithm 1 illustrates how the blocks are supposed to be evicted along the path.

Metadata Scan. As in Circuit ORAM [29], we first scan the metadata to collect the relevant information for moving the blocks in path . The difference here is that we give a parallel version for scanning the metadata. Algorithm 3 is supposed to produce an array $\text{target}[0..L]$, where the entry $\text{target}[i] = \text{dst}_i$ means that if $\text{dst}_i \neq \perp$, the deepest block in $\text{path}[i]$ will be moved to $\text{path}[\text{dst}_i]$.

Subroutine PrepareDeepest. A sequential version of this procedure was given in the Circuit ORAM paper [29].

After calling $\text{PrepareDeepest}(\text{path}[0..L])$, for $1 \leq i \leq L$, the array entry $\text{deepest}[i]$ stores the source height of the deepest block in $\text{path}[0..i-1]$ that can legally reside in $\text{path}[i]$.

We will use the following monotone property.

Fact 1 (Monotone property of deepest). *Suppose for some $1 \leq i \leq L$, $\text{deepest}[i] = j$. Then, for $j < k \leq i$, $\text{deepest}[k] = j$.*

Subroutine PrepareTarget. A sequential version was given in the Circuit ORAM paper [29] to process the metadata. We assume $\text{PrepareDeepest}(\text{path}[0..L])$ has already been called to have $\text{deepest}[0..L]$ ready. For the parallel version, $\text{PrepareTarget}(\text{path}[i..j], \text{dst})$ takes a path segment and a potential destination height $\text{dst} \geq j+1$. It prepares two arrays $\text{target1}[i..j]$ and $\text{target2}[i..j]$ and returns a pair $(\text{dst}_1, \text{dst}_2)$ with the following properties.

For $i \leq k \leq j$, $\text{target1}[k]$ indicates the destination height that the deepest block in $\text{path}[k]$ should be moved to, assuming that no block will be moved from $\text{path}[0..j]$ to $\text{path}[j+1..L]$; in this case, if $\text{dst}_1 \neq \perp$, then $i \leq \text{dst}_1 \leq j$, and there should be a block moving from $\text{path}[0..i-1]$ to

Algorithm 2 PrepareDeepest(path[i..j])

*/*Make parallel metadata scan to prepare the deepest array.*

After this algorithm, for each $i < k \leq j$, $\text{deepest}[k]$ stores the source height of the deepest block in $\text{path}[i..k-1]$ that can legally reside in $\text{path}[k]$, and $\text{goal}[k]$ stores the deepest height that the corresponding block can reside. Moreover, the procedure returns a pair (src, dst), where src is the source height the deepest block in $\text{path}[i..j]$ and dst is the deepest height the corresponding block can reside.

***Note:** The procedure can be implemented in an oblivious way. For instance, for the case when the condition of an “If” statement is false, dummy operations can access the corresponding variables without actually modifying them. */*

```
1: Initialize  $\text{deepest}[i..j] := (\perp, \perp, \dots, \perp)$  and  $\text{goal}[i..j] := (\perp, \perp, \dots, \perp)$ .
2: if ( $i == j$ ) then
3:   if  $\text{path}[i]$  is non-empty then
4:      $\text{dst} :=$  Deepest height that a block in  $\text{path}[i]$  can legally reside on  $\text{path}$ .
     return ( $i, \text{dst}$ )
5:   else return  $(\perp, \perp)$  ▷ Assume  $\perp$  is smaller than any integer.
▷  $i < j$ 
6:  $p := \lfloor \frac{i+j}{2} \rfloor$ 
7: Execute the following two statements in parallel:
   •  $(\text{src}_1, \text{dst}_1) := \text{PrepareDeepest}(\text{path}[i..p])$ 
   •  $(\text{src}_2, \text{dst}_2) := \text{PrepareDeepest}(\text{path}[p+1..j])$ 
8: for  $k = p+1$  to  $j$  in parallel do
9:   if  $\text{dst}_1 \geq \text{goal}[k]$  then
10:     $\text{deepest}[k] := \text{src}_1, \text{goal}[k] := \text{dst}_1$ 
11: if  $\text{dst}_1 \geq \text{dst}_2$  then return  $(\text{src}_1, \text{dst}_1)$ 
12: else return  $(\text{src}_2, \text{dst}_2)$ 
```

$\text{path}[\text{dst}_1]$. Similarly, $\text{target2}[k]$ indicates the corresponding destination height, assuming that if $\text{dst} \neq \perp$, then $\text{deepest}[j+1] = \text{deepest}[\text{dst}]$ and the deepest block in $\text{path}[\text{deepest}[j+1]]$ will be moved to $\text{path}[\text{dst}]$; similarly, in this case $\text{dst}_2 \neq \perp$ implies that there should be a block moving from $\text{path}[0..i-1]$ to $\text{path}[\text{dst}_2]$, where $\text{dst}_2 \geq i$ and $\text{deepest}[\text{dst}_2] < i$.

Hence, calling $\text{PrepareTarget}(\text{path}[0..L], \perp)$ will give the desired array $\text{target}[0..L] := \text{target1}[0..L]$ for moving the actual blocks.

Subroutine EvictFast. After the array $\text{target1}[0..L]$ is prepared by running $\text{PrepareTarget}(\text{path}[0..L], \perp)$, we can run EvictFast , which moves the blocks in parallel, as opposed to doing a linear scan from the root to the leaf. In particular, the effect of procedure $\text{EvictFast}(\text{path}[i..j])$ is to evict along the path starting from the i th height down to height j , following $\text{target}[i..j]$; the procedure returns a pair (B, dst) , where (B, dst) is any block that is supposed to be moved from $\text{path}[i..j]$ to $\text{path}[\text{dst}]$, where $\text{dst} > j$.

Hence, it suffices to call $\text{EvictFast}(\text{path}[0..L])$ to complete the eviction along the whole path.

Performance Analysis. All the algorithms are recursive and use a divide and conquer paradigm. Specifically, for an instance of size L , each algorithm first solves instances with sizes $\lfloor \frac{L}{2} \rfloor$ and $\lceil \frac{L}{2} \rceil$ in parallel. Then, the two solutions are combined to produce a solution to the original instance using $O(1)$ parallel runtime and $O(L)$ total work.

Algorithm 3 PrepareTarget(path[i..j], dst)*/* Parallel version to prepare the target arrays. */*

```
1: target1[i..j] := target2[i..j] := ( $\perp$ ,  $\perp$ , ...,  $\perp$ ), dst1 := dst2 :=  $\perp$ 
2: if (i == j) then
3:   if (path[i] has an empty slot) and (deepest[i]  $\neq$   $\perp$ ) then
4:     dst1 := dst2 := i
5:   if (deepest[dst] == i) then target2[i] := dst
6:     if deepest[i]  $\neq$   $\perp$  then dst2 := i
7:   else if deepest[dst] < i then dst2 := dst
8:   return (dst1, dst2)
▷ i < j
9: p :=  $\lfloor \frac{i+j}{2} \rfloor$ 
10: Execute the following two statements in parallel:
    • (dst1, dst2) := PrepareTarget(path[p+1..j], dst)
    • (dst'1, dst'2) := PrepareTarget(path[i..p], p+1)
11: if dst2  $\neq$   $\perp$  then ▷ deepest[dst2] = deepest[p+1]
12:   for k = i to p in parallel do
13:     if deepest[p+1] == k then target2[k] := dst2
14:   if dst1  $\neq$   $\perp$  then ▷ deepest[dst1] = deepest[p+1]
15:     for k = i to p in parallel do
16:       if (deepest[p+1] == k) then target1[k] := dst1
17:       else target1[k] := target2[k]
18:     if (dst'2 == p+1) then dst'1 := dst1
19:     else dst'1 := dst'2
20:   if (dst'2 == p+1) then dst'2 := dst2
▷ dst1 = dst2 =  $\perp$ 
21: if dst2 =  $\perp$  then
22:   for k = i to p in parallel do target2[k] := target1[k]
23:   dst'2 := dst'1
24: return (dst'1, dst'2)
```

Algorithm 4 EvictFast(path[i..j])

/ Evicts blocks along path[i..j] according to target[i..j] and return a pair (B, dst) for any block B to be moved from path[i..j] down to height dst. */*

```
1: if (i == j) then
2:   if target[i] ≠ ⊥ then
3:     B := Deepest block in path[i].
4:     Remove B from path[i].
5:     return (B, target[i])
6:   else return (⊥, ⊥)
7: for p := ⌊ $\frac{i+j}{2}$ ⌋ to j in parallel do
8:   Execute the following two statements in parallel:
9:     • (B1, dst1) := EvictFast(path[i..p])
10:    • (B2, dst2) := EvictFast(path[p+1..j])
11: for k = p+1 to j in parallel do
12:   if (dst1 == k) then put block B1 in path[k].
13: if dst1 > j then return (B1, dst1)
14: else return (B2, dst2)
```

▷ $i < j$

Hence, the recursion for parallel runtime is $T(L) = T(\frac{L}{2}) + O(1)$, which gives $T(L) = O(\log L)$; the recursion for total work is $W(L) = 2W(\frac{L}{2}) + O(L)$, which gives $W(L) = O(L \log L)$.

Correctness. We next show that EvictFast (Algorithm 4) achieves the same effect on the locations of the blocks as EvictSlow (Algorithm 1). The intuition is that the sequential version of the path eviction consists of three linear scans along the path. Hence, using the divide-and-conquer strategy in the parallel version, it suffices to keep track of what computation each sub-instance can perform on its own, and what information needs to be passed from one sub-instance to the other in order to complete the computation.

Lemma 8 (Correctness of EvictFast). *EvictFast has the same effect on the locations of the blocks in the eviction path as EvictSlow.*

Proof. We show that each of the algorithms PrepareDeepest, PrepareTarget and EvictFast achieve the effects in the above description. In particular, as seen from [29], in the sequential versions of these algorithms, PrepareDeepest and EvictFast each performs a root-to-leaf scan, and PrepareTarget performs a leaf-to-root scan. We explain how the solutions to the two sub-instances in each algorithm can be combined to give a solution to the larger instance.

1. **PrepareDeepest.** Since the effect of the algorithm can be achieved by a linear scan from the root to the leaf in path, after the sub-instances [i..p] and [p+1..j] have been handled in parallel, we just need to consider what information needs to be passed from the first instance to the second one to get a solution for the larger instance [i..j].

Observe that for each $k \in [p+1..j]$, $\text{deepest}[k]$ is supposed to store the source height of the deepest block from $\text{path}[i..k-1]$ that can legally reside in $\text{path}[k]$.

Starting from a solution for the sub-instance [p+1..j], we just need to know if where the deepest block in $\text{path}[i..p]$ can reside in $\text{path}[p+1..j]$, i.e., the source level src_1 of the deepest block in $\text{path}[i..p]$ and the deepest height dst_1 that it can reside. Equipped this information, the solution for [p+1..j] can be updated accordingly as shown in Algorithm 2.

2. **PrepareTarget.** Assuming that **PrepareDeepest** has already been called to construct the array **deepest**, Algorithm 3 performs a leaf-to-root scan. In the sequential version, we can imagine that when the scan passes from height $i + 1$ to height i , the algorithm just needs to remember a destination level $\text{dst} \geq i + 1$ that has an empty slot and is supposed to take a block from $\text{deepest}[\text{dst}]$. However, the monotone property from Fact 1 implies that $\text{deepest}[\text{dst}] = \text{deepest}[i + 1]$, which means the algorithm only needs to remember **dst**.

Now, suppose we break the larger instance $[i..j]$ into two smaller sub-instances $[i..p]$ and $[p+1..j]$. The difficulty is that if we try to solve the sub-instance $[i..p]$, we do not know whether a block is supposed to go from $\text{path}[i..p]$ to $\text{path}[p + 1..j]$, without solving the sub-instance $[p + 1..j]$ first. The important observation is that if a block is supposed to go from $\text{path}[i..p]$ to some destination height **dst** in $\text{path}[p + 1..j]$, even if we do not know what **dst** is, we know the source level of this block is given by $\text{deepest}[p + 1] = \text{deepest}[\text{dst}]$.

This intuition suggest that we need to prepare for two scenarios. The first scenario is that no block is supposed to go from $\text{path}[i..p]$ to $\text{path}[p + 1 :]$, and this corresponds to the solution in $\text{target1}[i..p]$. The second scenario is that a block is supposed to go from $\text{path}[i..p]$ to $\text{path}[p + 1 :]$, and so we pretend that the destination height is $p + 1$ for the time being to produce $\text{target1}[i..p]$.

Once the solutions **target1** and **target2** have been produced for the sub-instance $[p + 1..j]$, we know whether a block is supposed to go from $\text{path}[i..p]$ to $\text{path}[p + 1 :]$ in each case, and we also know the true destination of that block. Hence, we can use this information to correct the solutions for the sub-instance $[i..p]$ accordingly.

3. **EvictFast.** After **PrepareTarget** has been called, the sequential version just needs to perform a root-to-leaf scan to move the blocks according to $\text{target1}[0..L - 1]$.

After the sub-instances $[i..p]$ and $[p + 1..j]$ are solved, we just need to know if any block is supposed to be moved from $\text{path}[i..p]$ to $\text{path}[p + 1 :]$. If yes, the block can either be moved to $\text{path}[p + 1..j]$ or passed to $\text{path}[j + 1 :]$, as shown in Algorithm 4.

□

7.2 Theorem Statement for Larger Block Sizes

The next lemma summarizes the performance analysis when the small-depth path eviction algorithm is used. Observe that there is an extra $\log \log N$ factor in the work for eviction. Hence, this algorithm is used for moderate block sizes, i.e., when the number of recursion levels is $o(\log N)$. The lemma is parameterized with Γ , which is the number of position labels that a block can store.

Lemma 9. *Suppose each block can store Γ position labels. The number of recursion levels is $D = O(\frac{\log N}{\log \Gamma})$, and denote $\gamma := \min\{\Gamma, m\}$. Then, each step of the PRAM with m CPUs can be simulated with $O(Dm(\log m + \log N \log \log N))$ total work and $O(D(\log \gamma + \log \log N) + \log m)$ parallel runtime.*

Proof. By Lemma 6, the fetch phase can be performed with $O(Dm \cdot (\log m + \log \mu))$ total work and $O(D \log \mu + \log m)$ parallel runtime, where $\mu := \alpha \gamma^2 \log N$.

With the small-depth eviction algorithm, the maintain phase takes total work $O(Dm \log N \log \log N)$ and $O(\log m + \log \log N)$ parallel runtime.

Combining the performance analysis of both phases gives the result.

□

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Appendix

A Additional Preliminaries

A.1 Notations

Statistical and computational indistinguishability. Given two ensembles of distributions $\{X_N\}$ and $\{Y_N\}$ (parameterized with N), we use the notation $\{X_N\} \stackrel{\epsilon(N)}{\equiv} \{Y_N\}$ to mean that for any (possibly computationally unbounded) adversary \mathcal{A} ,

$$\left| \Pr[\mathcal{A}(x) = 1 \mid x \stackrel{\$}{\leftarrow} X_N] - \Pr[\mathcal{A}(y) = 1 \mid y \stackrel{\$}{\leftarrow} Y_N] \right| \leq \epsilon(N).$$

We use the notation $\{X_N\} \stackrel{\epsilon(N)}{\equiv_c} \{Y_N\}$ to mean that for any non-uniform p.p.t. adversary \mathcal{A} ,

$$\left| \Pr[\mathcal{A}(1^N, x) = 1 \mid x \stackrel{\$}{\leftarrow} X_N] - \Pr[\mathcal{A}(1^N, y) = 1 \mid y \stackrel{\$}{\leftarrow} Y_N] \right| \leq \epsilon(N).$$

A.2 Adaptive, Composable Notion of OPRAM

As mentioned earlier, although using a static notion of security allows us to prove a stronger lower bound, our upper bound in fact satisfies a stronger, adaptive and composable notion of security as we formally specify below. This notion of security was also adopted in several recent works [1]. For convenience, below we first define the computationally secure version and then we remark how to extend it to the statistically secure notion.

Definition 3 (Adaptively secure OPRAM). We say that a stateful parallel algorithm `opram` is a strongly-oblivious OPRAM algorithm with computational security iff there exists a non-uniform probabilistic polynomial-time simulator `Sim`, such that for any non-uniform probabilistic polynomial-time adversary \mathcal{A} , \mathcal{A} 's view in the following two experiments, $\text{Expt}_{\mathcal{A}}^{\text{real,opram}}$ and $\text{Expt}_{\mathcal{A}, \text{Sim}}^{\text{ideal}}$ are computationally indistinguishable:

$\text{Expt}_{\mathcal{A}}^{\text{real,opram}}(N, m):$ $\text{out}_0 = \text{addresses}_0 = \perp$ For $t = 1, 2, \dots \text{poly}(N)$: $\vec{I}^{(t)} \leftarrow \mathcal{A}(N, m, \text{out}_{t-1}, \text{addresses}_{t-1})$ $\text{out}_t, \text{addresses}_t \leftarrow \text{opram}(\vec{I}^{(t)})$	$\text{Expt}_{\mathcal{A}, \text{Sim}}^{\text{ideal}}(N, m):$ $\text{out}_0 = \text{addresses}_0 = \perp$ For $t = 1, 2, \dots \text{poly}(N)$: $\vec{I}^{(t)} \leftarrow \mathcal{A}(N, m, \text{out}_{t-1}, \text{addresses}_{t-1})$ $\text{out}_t \leftarrow \mathcal{F}_{\text{mem}}(\vec{I}^{(t)}), \text{addresses}_t \leftarrow \text{Sim}(N, m)$
--	---

In the above, \mathcal{F}_{mem} is the “ideal logical memory functionality” defined in the most natural way, and addresses_t denotes the physical memory locations accessed for the t -th batch of requests, and out_t denotes the outcome of the computation at the end of the t -th batch of requests.

The above definition is for computational security, but we can easily extend it to statistical security in the usual manner. Specifically, if we replace computational indistinguishability with statistical indistinguishability and remove the requirement for the adversary to be polynomially bounded, then we then say that `opram` is an OPRAM algorithm with statistical security.

We remark that although most existing works on ORAM [13, 14, 19, 26, 28, 29] and OPRAM [2, 6] often adopt a weaker, statically secure notion that is not composable (equivalent to our earlier

definition in Section 2.2) it is not difficult to observe that almost all known ORAM and OPRAM constructions [2,6,13,14,19,26,28,29] also satisfy the stronger notion, i.e., our Definition 3. Similarly, it is not difficult to see that our small-depth OPRAM construction also satisfies this stronger notion of security.

A.3 Details on Oblivious Random Permutation

For completeness, we formally define what an oblivious random permutation is. We also prove that the algorithm in Figure 2 realizes this formal abstraction, and we analyze its performance.

Definition 4 (Oblivious random permutation). We say that a PRAM algorithm Alg realizes a perfectly secure oblivious random permutation if there exists a simulator Sim such that for any input array I , where n is the length of I , we have

$$\text{REAL}^{\text{Alg}}(I) \equiv (\mathcal{F}_{\text{perm}}(I), \text{Sim}(n)),$$

where the notation \equiv denotes identically distributed, and the notation $\text{REAL}^{\text{Alg}}(I)$ denotes a pair of random variables containing: 1) the output of the running Alg on the input I , and 2) the ordered sequence of addresses observed in all steps of the PRAM algorithm Alg upon input I . Here, $\mathcal{F}_{\text{perm}}(I)$ is the ideal “random permutation functionality” which permutes the input array I at random and outputs the permuted array.

In the above definition, although not explicitly denoted, we assume that the memory word size and the number m of CPUs are public knowledge hard-coded in both Alg and Sim .

Realizing a perfectly secure oblivious random permutation. We now prove that the algorithm described in Figure 2 realizes an oblivious random permutation; we also bound its parallel runtime and total work.

Before we do that, we fill in the missing details regarding the naïve quadratic random permutation algorithm that Figure 2 used as a subroutine. Given a (small) array containing s elements, a naïve quadratic random permutation algorithm proceeds sequentially as the following. Initially, every element in the array is *unconsumed*. For $i = 1$ to s : 1) sample a random number j between 1 and $s - i + 1$; and 2) in one linear scan of the array, find the j -th unconsumed element, mark it as *consumed*, and write it into the i -th position of the result array. It is straightforward to see that this naïve quadratic algorithm realizes an oblivious random permutation. Its total work and runtime are both quadratic in the array length s .

We proceed to bound the parallel runtime and total work of the algorithm described in Figure 2.

Lemma 10. *Assume that the length k of the input array satisfies $\log^{0.5} N \leq k \leq N$. Except with probability negligible in N , the algorithm in Figure 2 completes with $O(\log k)$ parallel runtime and $O(k \log k)$ total work.*

Proof. We first prove that for the super-constant function $\alpha(N) := \log \log N$, after one iteration of the algorithm, except with negligible in N probability, at most $\alpha(N)$ elements receive the same random key.

Since $3 \log_2 N$ random bits are assigned to each element of A as the key, we can model the process as throwing $k \leq N$ balls into N^3 bins. Without loss of generality, henceforth we simply assume $k = N$. For any fixed set S of $\alpha(N)$ elements, the probability that every element in S collides with other elements is upper bounded by $\left(\frac{N}{N^3}\right)^\alpha = \frac{1}{N^{2\alpha}}$. Thus, the probability that there exist $\alpha(N)$ elements that collide with other elements is upper bounded by:

$$\binom{N}{\alpha} \cdot \frac{1}{N^{2\alpha}}.$$

For $\alpha(N) = \omega(1)$, the above expression is bounded by a negligible function in N .

Hence, we next assume that at most $\alpha(N)$ elements receive the same key. Recall that we run the naïve quadratic random permutation for each subset of elements receiving the same key (in parallel). Observe that in this part, the parallel runtime is $O(\alpha(N)^2)$ and the total work is $O(k \cdot \alpha(N)) = O(k \log k)$. \square

Lemma 11. *The algorithm in Figure 2 realizes a perfectly secure oblivious random permutation as defined in Definition 4.*

Proof. We describe a simulator that simulates the access pattern. Since the oblivious sort and the naïve quadratic permutation parts of the algorithm have deterministic access patterns, without loss of generality, henceforth we ignore simulating this part of the access patterns. Similarly, some other parts of the algorithm (e.g., sequential scan to assign random numbers to each coordinate) are deterministic and we ignore those as well in describing the simulator. The remaining non-trivial access patterns to simulate include which coordinates have colliding elements — the access patterns of the oblivious aggregation step and the step of copying out the collided elements are solely determined by this information and thus we ignore simulating these parts of the access patterns too.

In summary, we imagine a simulator denoted Sim that chooses k random keys of $3 \log_2 N$ bits long, and sorts them in increasing order. Now, the simulator outputs which coordinates in the sorted array have collisions. By definition, we need to show that the following joint distributions are identical:

1. the joint distribution of the real-world algorithm’s resulting permutation and which elements collide in the real world henceforth denoted as

$$(\pi_{\text{real}}, C_{\text{real}}),$$

2. the joint distribution containing 1) a completely random permutation, and 2) the simulated output of which elements collide, henceforth denoted as

$$(\mathcal{F}_{\text{perm}}(k), \text{Sim}).$$

We now show that the above joint distributions are identical. First, it can be shown that the marginal distributions $\pi_{\text{real}} \equiv \mathcal{F}_{\text{perm}}(k)$ and $C_{\text{real}} \equiv \text{Sim}$, where \equiv denotes identically distributed. It thus suffices to prove that even when conditioned on observing C_{real} , the real-world resulting permutation π_{real} is still distributed as a completely random permutation. This can be proven using a simple coupling argument as follows. Consider all sample paths where the collision pattern is fixed to some pattern, and the random string chosen in the second phase (i.e., the quadratic permutation phase) is fixed to some π — note that π can be interpreted as a permutation. Given any two final permutations π_0 and π_1 , if the first-phase random string χ_0 results in π_0 (conditioned on collision pattern and π), then observe that $\chi_1 = \chi_0 \pi_0^{-1} \pi_1$ would have the same collision pattern at the end of the first phase, and would result in the final permutation π_1 . Further, to result in the final permutations π_0 and π_1 , χ_0 and χ_1 are obviously unique in the conditional sample space we consider (by fixing the access pattern and π) and they have equal probability mass in the conditional sample space we consider. The remainder of the argument follows in a straightforward manner by summing over all possible choices of collision pattern and π . \square

A.4 Additional Building Blocks

We define some additional building blocks. All of these building blocks are oblivious in the sense that they have deterministic access patterns that do not depend on the inputs.

Oblivious aggregation for a sorted array. Oblivious aggregation is the following primitive where given a somewhat sorted array of (key, value) pairs, each representative element for a key will learn some aggregation function computed over all pairs with the same key.

- *Input:* An array $\text{Inp} := \{k_i, v_i\}_{i \in [n]}$ of possibly dummy (key, value) pairs, where all pairs with the same key appear in consecutive locations. Henceforth we refer to all elements with the same key as the same *group*. We say that index $i \in [n]$ is a *representative* for its group if i is the leftmost element of the group.
- *Output:* Let Aggr be a publicly known, commutative and associative aggregation function and we assume that its output range can be described by $O(1)$ number of blocks. The goal of oblivious aggregation is to output the following array:

$$\text{Outp}_i := \begin{cases} \text{Aggr}(\{(k, v) \mid (k, v) \in \text{Inp} \text{ and } k = k_i\}) & \text{if } i \text{ is a representative} \\ \perp & \text{o.w.} \end{cases}$$

Oblivious routing. Oblivious routing is the following primitive where n source CPUs wish to route data to n' destination CPUs based on the key.

- *Inputs:* The inputs contain two arrays: 1) a source array $\text{src} := \{(k_i, v_i)\}_{i \in [n]}$ where each element is a (key, value) pair or a dummy element denoted (\perp, \perp) ; and 2) a destination array $\text{dst} := \{k'_i\}_{i \in [n']}$ containing a list of (possibly dummy) keys.

We assume that each (non-dummy) key appears only once in the src array, however, each (non-dummy) key can appear multiple times in dst .

- *Outputs:* We would like to output two arrays: 1) an array $\text{Out} := \{v'_i\}_{i \in [n']}$ such that for each $i \in [n]$, it holds that either $(k'_i, v'_i) \in \text{src}$, or $k'_i \notin \text{src}$ in which case $v'_i := \perp$, or $k'_i = v'_i = \perp$; and 2) an additional array Remain of length n that holds all remaining elements of src that have not been routed to the destination array (padded with dummies).

Oblivious select. Parallel oblivious select solves the following problem: given an array containing n values, v_1, v_2, \dots, v_n , compute the minimum value $\min(v_1, \dots, v_n)$. Parallel oblivious select can be achieved through an aggregation tree: imagine a binary tree where the leaves represent the initial values. We assign one CPU to each node in the tree. Now for each height ℓ going from the leaves to the root, in parallel, each internal node at height ℓ computes the minimum of its two children. Clearly, oblivious select can be attained in $O(n)$ total work and $O(\log n)$ parallel steps.

In a straightforward fashion, oblivious select can also be applied to the following variant problem: suppose that each of n CPUs holds a value, and that only one CPU is holding a real value s (but which one is not known), and all others hold the dummy value \perp . At the end of the algorithm, we would like a designated fetch CPU to obtain the real value s .