Empirical Evaluation of the Parallel Distribution Sweeping Framework on Multicore Architectures

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Abstract

In this paper, we perform an empirical evaluation of the Parallel External Memory (PEM) model in the context of geometric problems. In particular, we implement the parallel distribution sweeping framework of Ajwani, Sitchinava and Zeh to solve batched 1-dimensional stabbing max problem. While modern processors consist of sophisticated memory systems (multiple levels of caches, set associativity, TLB, prefetching), we empirically show that algorithms designed in simple models, that focus on minimizing the I/O transfers between shared memory and single level cache, can lead to efficient software on current multicore architectures. Our implementation exhibits significantly fewer accesses to slow DRAM and, therefore, outperforms traditional approaches based on plane sweep and two-way divide and conquer.

1 Introduction

Modern multicore architectures have complex memory systems involving multiple levels of private and/or shared caches, set associativity, TLBs, and prefetching effects. It is considered challenging to design and even engineer algorithms to directly optimize the running time on such architectures [16]. Furthermore, algorithms optimized for one architecture may not be optimal for another. To address these issues, various computational models [4, 8, 9, 13, 14] have been proposed in recent years. These computational models are simple (usually assuming only two levels of memory hierarchy, out of which one is shared) as they abstract away the messy architectural details. Also, the performance metric of these models involve a single objective function such as minimizing shared memory accesses. The simplicity of these models allows the design of practical algorithms that are expected to work well on various multicore architectures. It also allows us to compare the relative performance of algorithms theoretically.

The success of a computational model crucially depends on how well the theoretical prediction of an algorithm in that model matches the actual running time on real systems. Unfortunately so far, there has been little empirical work (such as [21]) to evaluate the predictions of algorithmic performance using these models on real multicore architectures. It is not even clear if these models can lead to the design of algorithms that are faster on current multicore systems (with 2 - 48 cores) than those designed in the traditional RAM model, external memory model and the PRAM model. In fact, many of the algorithms designed in these models for multicore seem quite sophisticated and are likely to have high constant factors that can pay off only for architectures with hundreds
of cores. This state of affairs is in sharp contrast with the sequential cache-efficient models, where a considerable empirical work (e.g., [3, 10]) evaluating the algorithms on real systems exists.

At the core of the debate for the computational model is the choice of the performance metric that an algorithm designer should optimize for the current multicore systems. In the traditional RAM (and PRAM) model of computation, the algorithms are designed to minimize the number of instructions (and parallel instructions) executed by the algorithm. The external memory (EM) model [1] when applied to cached memories (e.g., see [17]) aims at minimizing the cache misses, ignoring the number of instructions. The parallel external memory (PEM) model [4] aims at minimizing the number of parallel cache misses.

In this work, we demonstrate that algorithms designed in simple models, that focus on minimizing the parallel I/O transfers between shared memory and a single level cache, can lead to a software performing great in practice on real multicore systems. For this purpose, we consider the algorithms to solve the problem of answering batched planar orthogonal stabbing-max queries. This problem is a fundamental geometric primitive and together with its variants is used as subroutines in solutions of many popular geometric problems such as point location in an orthogonal subdivision of the plane, orthogonal ray shooting, batched (offline) dynamic predecessor queries in 1-dimensional array and batched union-find. Also, this problem has been well-studied in various computational models and many different optimal solutions for it are known in these models. Thus, it provides a test-bed for evaluating the efficacy of theoretical analysis in various models on real multicore architectures. Another reason for selecting this non-HPC application is that the ratio of memory accesses to computation in the solutions of this problem is similar to that of many data-intensive geometric applications. For instance, our engineered PEM solution for this problem is based on the parallel distribution sweeping framework and this framework has been used for designing a wide range of other geometric algorithms in the PEM model [2, 3] and a basis for PEM data structures [20].

We empirically compare the different solutions and show that a carefully engineered solution based on an algorithm in the PEM model gives the best performance on various multicore systems, outperforming traditional approaches based on plane sweep, sequential distribution sweeping and two-way divide-and-conquer. Using hardware profilers, we show that this solution exhibits significantly fewer number of accesses to slow DRAM which is correlated with the improved running time.

Since the cache line on modern systems is typically 64 bytes, I/O-efficient solutions also need to be work-efficient to compete with RAM algorithms. In other words, the total number of instructions of a cache-efficient algorithm should asymptotically match that of the best RAM solution. Therefore, we design an algorithm that is both I/O-optimal and work-efficient. To the best of our knowledge, this is the first work-efficient I/O-optimal algorithm for this problem.

2 Computational Models

External Memory Model. The widely used external memory model or the I/O model by Aggarwal and Vitter [1] assumes a two level memory hierarchy. The internal memory has a limited size and can hold at most \( M \) objects (points/line-segments) and the external memory has a conceptually unlimited size. The computation can only use the data in the internal memory, while the input and the output are stored in the external memory. The data transfer between the two memories happens in blocks of \( B \) objects. The measure of performance of an algorithm
is the number of I/Os (cache misses) it performs. The number of I/Os needed to read $n$ contiguous items from the external memory is $\text{scan}(n) = \Theta(n/B)$. The number of I/Os required to sort $n$ items is $\text{sort}(n) = \Theta(\frac{n}{B} \log \frac{M}{B}(n/B))$. For all realistic values of $n$, $B$, and $M$, $\text{scan}(n) < \text{sort}(n) < n \log_2 n$.

**Parallel External Memory (PEM) Model.** The parallel external memory (PEM) model \[4\] is a simple parallelization of the EM model. It consists of $P$ processors, each with a private cache of size $M$ (see Figure 1). Processors communicate with each other through access to a shared memory of conceptually unlimited size. Each processor can use only data in its private cache for computation.

The caches and the shared memory are divided into blocks of size $B$. Data is transferred between the caches and shared memory using parallel input-output (I/O) operations. During each such operation, each processor can transfer one block between shared memory and its private cache. The cost of an algorithm is the number of I/Os it performs. Concurrent reading of the same block by multiple processors is allowed but concurrent block writes are disallowed (similar to a CREW PRAM). The cost of sorting in this model is $\text{sort}_P(n) = O(\frac{n}{PB} \log \frac{M}{B}(\frac{n}{B}))$ parallel I/Os, provided $P \leq \frac{n}{B^2}$ and $M = B^{O(1)}$ \[4\].

The PEM model provides the simplest possible abstraction of current multicore chips, focusing on the fundamental I/O issues that need to be addressed when designing algorithms for these architectures, similar to the I/O model \[1\] in the sequential setting.

### 3 1-D Stabbing Max Algorithms

In this section, we describe various algorithms that we implemented and used for our experimental study. We begin with formally describing the problem.

**Definition 1 (Batched 1-D Stabbing-Max Problem)** Given a set of $n$ horizontal line segments and points on the plane, report for each point the closest segment that lies directly below it.

**RAM algorithm.** In the classical RAM model, this problem is solved using the sweep line paradigm \[18, 6\]. We sweep a hypothetical vertical line across the plane in increasing $x$-coordinate and perform some computation at each segment endpoint or query point. We maintain an ordered set $A$ of active segments — all segments which intersect the sweep line, ordered by the $y$-coordinates. A segment is inserted into $A$ when the sweep line encounters its left endpoint and removed when it encounters the right endpoint. An answer to a query point $q$ is the segment in $A$ with the largest $y$-coordinate that is smaller than the $y$-coordinate of $q$, i.e., the predecessor of $q$ in $A$ according to the $y$-ordering.

For $n$ line segments and query points, there are $O(n)$ insertions, deletions and predecessor searches in $A$. Since each of these operations can be performed in $O(\log n)$ time by maintaining $A$ as a balanced binary search tree, the total complexity of this algorithm is $O(n \log n)$ instructions.

**Sequential I/O-optimal solution.** The sequential I/O-efficient solution for this problem proceeds using the distribution sweeping framework of Goodrich et al. \[15\] as follows.
Let \( r_q \) be a variable associated with each query point \( q \) which we will use to store the answer. Initially \( r_q \) is initialized to a virtual horizontal line \( y = -\infty \).

We partition the space into \( K = \min\{M/B, n/M\} \) vertical slabs \( \sigma_1, \ldots, \sigma_K \), so that each slab contains equal number of points (endpoints of horizontal segments or query points) and perform a sweep of the input by increasing \( y \)-coordinate. During the sweep we maintain for each slab \( \sigma_i \) a segment \( s_{\sigma_i} \) which is the highest segment that spans \( \sigma_i \) encountered by the sweep. When the sweep line encounters the query point \( q \in \sigma_i \), we update \( r_q \) with \( s_{\sigma_i} \) iff \( y(s_{\sigma_i}) > y(r_q) \). During the sweep we also generate slab lists \( \sigma_i \). A copy of a query \( q \) (resp., segment \( s \)) is added to \( \sigma_i \) if \( q \) (resp., at least one of the endpoints of \( s \)) lies in slab \( \sigma_i \). The sweep is followed by a recursive processing of each slab, using \( \sigma_i \) as input for the recursive call. The recursion terminates when each slab contains \( O(M) \) points and the problem can be solved in internal memory, for example, by using the plane sweep algorithm.

Note, that if the initial objects are sorted by \( y \)-coordinates, we can generate the inputs \( \sigma_i \) for the recursive calls sorted by \( y \)-coordinate during the sweep. Thus, the sweep at each of \( O(1 + \log_K(n/M)) \) recursive levels takes \( O(n/B) \) I/Os and the total I/O complexity of distribution sweeping is \( O \left( \frac{n}{P} \right) \left( 1 + \log_K(n/M) \right) = \text{sort}(n) \) I/Os.

**Work-optimal solution.** Note that a naive implementation of the sweep in internal memory might potentially result in updating \( K \) different variables \( s_{\sigma_i} \) whenever a segment is encountered during the sweep. This could lead to \( O(Kn) \) instructions at each recursive level, resulting in total \( O(Kn \log_K n) \) instructions, which is larger than \( O(n \log_2 n) \) instructions of the plane sweep algorithm. At the same time, the plane sweep algorithm could result in up to \( O(n \log_2 n) \) I/Os, which is larger than \( \text{sort}(n) \) I/Os of the above algorithm.

To achieve optimal internal computation time while maintaining the optimal \( \text{sort}(n) \) I/O complexity we store segments \( s_{\sigma_i} \) in a segment tree \( T \) over \( K \) intervals defined by the slabs \( \sigma_i \). Since, we are interested only in segments that fully span the slabs, each segment is stored only in one node. Also, at each node we store only the highest segment encountered up to that point in the sweep. Thus, \( |T| = O(K) \), i.e. \( T \) fits in internal memory. Consider the nodes on the root to leaf path which correspond to the intervals containing \( q \). We update \( r_q \) to the highest segment stored at these nodes. Thus, maintaining \( T \) and updating \( r_q \) takes \( O(\log_K K) \) instructions per update/query, and over \( O(1 + \log_K N/M) \) recursive levels of distribution sweeping adds up to at most \( O(n \log_2 n) \) instructions, which is optimal.

**Parallel External Memory Solution.** The PEM solution is based on the parallel distribution sweeping framework introduced by Ajwani et al. [2]. It differs from the sequential distribution sweeping by recursively dividing the plane into \( K := \max\{2, \min\{\sqrt{n/P}, M/B, P\}\} \) vertical slabs\(^1\) and performing the sweep in parallel using all \( P \) processors. During recursion, the slabs are processed concurrently using sets of \( \Theta(P/K) \) distinct processors per slab. The parallel recursion proceeds for \( O(\log_K P) \) rounds, until there are \( \Theta(P) \) slabs remaining, at which point, each slab is processed concurrently using a single processor running the sequential I/O-efficient solution.

To perform the sweep of a single recursive level in parallel using multiple processors, each processor performs distribution sweeping on an equal fraction of the input. Note, that such a sweep sets the values of \( r_q \) correctly only if both the query \( q \) and the spanning segment \( s_{\sigma_i} \) below it are processed by the same processor. To correct the values \( r_q \) across the boundaries of the parallel sweeps we perform a round of parallel reduction on segments and queries using MAX associative operator [7]. Finally, we compact the portions of slab lists \( \sigma_i \) generated by different processors.

\(^1\)The explanation for this choice of \( K \) can be found in [4].
into contiguous slab lists to be used as input for recursive calls. The details of the algorithm follow directly from [2] but are also presented in Appendix A for completeness.

The parallel I/O complexity of the above algorithm is \( O(\text{sort}_P(n)) \) I/Os.

**Work-optimal solution.** Similar to the sequential I/O model, we can achieve work optimality in the PEM model algorithm by maintaining a segment tree \( T \) on the \( K \) child slabs. In this case, all processors keep their own copy of \( T \) and the parallel reduction (using MAX operator) is performed over not only the \( K \) leaves, but also the \( K - 1 \) internal nodes of \( T \). This does not affect the asymptotic number of parallel I/Os, but makes the scheme work-optimal, i.e. \( O(\frac{n}{P} \log n) \) instructions per processor.

**2-way Distribution Sweeping.** As a PRAM solution, we consider a recursive 2-way distribution sweeping algorithm. This framework is akin to divide-and-conquer paradigm, that is archetype for many PRAM algorithms. The 2-way distribution is continued recursively till the slab size is smaller than a fixed constant and at that stage, plane sweep algorithm is used as a base case. The distribution step is a simplified version of the corresponding step in the PEM algorithm, as the considerations of work-optimality no longer apply.

### 4 Implementation Details

We implemented our algorithms in C++, using OpenMP for parallelization. The engineered implementation uses some simple techniques to improve the running time of the theoretical algorithm, while trying to preserve its worst-case asymptotic guarantee on the number of shared cache accesses.

The parallel distribution sweeping calls for setting the branching parameter at \( K = \max\{2, \min\{M/B, \sqrt{n/P}, P\}\} \). The parameter \( M \) also defines the size of the recursive base case. We experimentally determine the best choice of \( M \). In particular we found that setting \( M \) to be a large fraction (e.g., 1/3 or 1/4) of the L3 cache results in best running times.

Having determined \( M \), we observe that for computing \( K \), in our compute systems the number of processors (up to 12) is far below the other two terms. Thus, the first recursive level is always a single \( P \)-way parallel distribution sweeping round, which results in \( P \) vertical slabs each of which can be processed independently of others in the consequent phases. Thus, after the parallel distribution, each of \( P \) resulting vertical slabs is assigned to a separate thread which processes it using a sequential distribution sweeping algorithm.

To perform the parallel sweep, we divide the input based on the \( y \)-coordinate among the \( P \) threads, conceptually, assigning a horizontal slab of objects to each thread. The thread with the smaller ID gets the lower \( y \) values. This can be viewed as a \( P \times P \) matrix where the columns correspond to the different slabs and the rows correspond to the different threads.

We perform the prefix sum on the \( P \times P \) array sequentially as the overheads associated with the synchronization barrier of OpenMP are too high to justify this operation in parallel.

We combine the second scan of the data (due to reduction) with the step of compacting child slab lists into contiguous vectors. During the compaction, each processor \( p_j \) copies all partial chunks of child slab \( \sigma_j \) into the contiguous space. Note, the propagation of the results of the prefix sums simply needs to update the result of each query point that had been assigned the sentinel line \( y = -\infty \) with the result of the prefix sums value. Thus, the propagation of the prefix sums values can be performed during this copying process.

\(^2\)In our experiments, performing this step sequentially takes less than a millisecond, while the overall running time is in dozens or hundreds of seconds.
Next, we process the $P$ child slabs in parallel using sequential distribution sweeping. This recursively subdivides the slabs till the pre-specified threshold $M$ is reached. When generating the input lists for the child slabs, we also store the total number of segments and query points for the child slabs. If for any slab, either the number of segments or query points is zero, we do not process it or its child invocations any further.

**Space efficiency.** We carefully engineered our algorithms to reduce the space requirement of our implementations considerably. This is done while ensuring that the running time of our implementations is not affected by the space reduction. We provide more details of this in Appendix B.

**Randomized vs. deterministic computation of slab boundaries.** Deterministic identification of slab boundaries such that all the child slabs at each level of recursion contain the same number of objects, requires sorting the input based on the $x$-coordinate and storing $O(n/M)$ equally spaced entries of the sorted input in a separate array. We avoid the extra sort by instead determining the slab boundaries by partitioning the space into uniform vertical slabs. This optimization works well for random input, but in the worst case can result in the recursion depth as large as $O(\log_K \delta)$, where $\delta$ is the spread of the point set – the ratio between the largest and the smallest (horizontal) distance between a pair of points. In case of a large base case of the recursion and randomized input, this is not an issue. But in the case of double precision coordinates, the worse case analysis dictates that the depth of the recursion can be very large.

**Constant factors vs. EM implementation.** The I/O complexity of the sequential distribution sweeping framework is $O(n/B(1 + \log_K n/M))$, where $K = \min\{M/B, n/M\}$. Since in our experimental settings $K = n/M$, there are only 2 recursive levels: one for distribution sweeping and one for the sweep line at the base case. Thus, the implementation performs two sequential scans of the input.

In the parallel version, we have to perform two additional scans. Specifically, we perform one extra recursive step – the parallel distribution. During this step, each processor scans $n/P$ items and writes them out into its private child slabs. After the prefix sums, which takes negligible amount of time, we must (a) propagate the result of the prefix sums to the queries that contain only sentinel values as the result and (b) construct each child slab in contiguous space. As described earlier, we combine these two tasks into a single scan.

Thus, combined with the two scans of the parallel recursive invocation of the sequential distribution sweeping, the parallel implementation performs a total of four scans of the input, i.e., twice as many as the sequential version. Since all scans are performed in parallel and in expectation each child slab contains equal number of items, the total I/Os performed by each processor is $2/P$ times the number of sequential I/Os, and (ignoring the speedup due to faster parallel internal computation) we should expect the speed up of $P/2$ on $P$ processors.

**Sorting.** To perform the initial sorting of the input by the $y$-coordinate, we used the sorting implementation from the C++ Multicore Standard Template Library (MCSTL) [19] that is now part of the GNU libstdc++ library. For the base case of plane sweep algorithm, we use the C++ Standard Template Library (STL) sorting implementation.

**Choice of $P$.** While in theory, $P$ denotes the number of cores, there are many considerations involved in picking the correct value of the parameter $P$ in practise. These considerations are discussed in Appendix C.
5 Experiments

We performed extensive experimentation studying the performance of these algorithms on various input types and on many different multicore architectures. In addition to measuring the running time of these algorithms, we used papi library and the Linux perfctr kernel module to read the hardware performance counters and measure cache misses, DRAM accesses, TLB misses, branch mispredictions, number of instructions etc. This section summarizes the key findings of our experiments.

Our query points were generated uniformly at random inside the grid of size Grid Size \times Grid Size. To elicit the asymptotic worst case performance of point location algorithms, we focus on long segments, whose length is chosen uniformly at random between Grid Size/4 and 3 \cdot Grid Size/4 and are at a random y-coordinate.

Configuration. We ran our implementation on the following multicore systems:

1. A system with a single 4-core 2.66 GHz Intel Core i7-920 processor and a total of 12.3GB RAM. Each core can run 2 threads due to hyperthreading. The processor has an L3 cache of size 8192 KB that is shared among all 4 cores. The L2 cache of 256 KB is only shared among pairs of cores.

2. A system with 4 \times 12-core 1.9 GHz AMD Opteron 6168 processors and total of 264 GB of RAM. Each core contains a private L2 cache of 512 KB and groups of 6 cores share an L3 cache of 5118KB. Thus, each processor contains two L3 caches of combined size of just over 10MB.

3. A system with 2 \times 16-core 2.6 GHz AMD Opteron 6282 SE processors and total of 96 GB RAM. Each core has its private L2 cache while the L3 cache is shared between 16 cores. The L2 cache size is 2 MB and L3 cache size is 16 MB.

All configurations run Linux kernels and the codebase was compiled using g++-2.4 compiler and -O3 flag.

Spatio-temporal locality in our setting. The cache line size for all cache levels on all 3 systems is 64 bytes. Since our objects take 32 bytes of space, it appears that each cache line can hold only two objects. Therefore, at a first glance it is not clear if I/O efficient algorithm can utilize the spatial locality for any improvement in runtime. However, we observed that given an array that is too large to fit in cache and which contains our 32-byte objects, it takes 4-5 times faster to access the objects sequentially rather than performing access in random locations. This observation can be explained by the fact that the memory system prefetches 2-3 cache lines when performing a sequential scan. Thus, during sequential scan the prefetcher amplifies the size of the cache line by the number of lines being prefetched\(^3\).

Another benefit of performing K-way distribution sweeping is that it allows us to utilize temporal locality by reducing the number of recursive calls. In particular, \(K = \min\{n/M, M/B\}\) and the number of recursive levels is \((1 + \log_K(n/M))\). Given limit of RAM size on our systems and the large size of L3 cache, it appears from our experiments that \(K\) is set to \(n/M\) on configuration 1 and 2, resulting in a single recursive level dedicated to (sequential) distribution (with the recursive base case performing plane sweep on chunks that fit in L3 cache).

\(^3\)For this experiment, the array must contain the actual objects and not just pointers to the objects, which could be allocated anywhere in memory.
On configuration 3, it requires two recursive calls. The various trade-offs involved in selecting the correct values of parameters $K$ and $M$ and the effect of these parameters on the actual run-time of our PEM implementation are described in Appendix D.

**Random access vs. I/O-efficient algorithms.** Figure 2 shows the absolute running times for the plane sweep and (parallel) distribution sweeping algorithms. One can see improvements in runtimes with the increase in the number of processors used. Also note the difference in the slopes in the graphs of the plane sweep algorithm compared to distribution sweeping algorithms. This is due to larger asymptotic number of cache misses of the plane sweep algorithm.

Figure 3 demonstrates this difference better. It shows the speedup of the sequential and parallel distribution sweeping algorithms relative to the plane sweep algorithm for long segments. In this figure one can see the effects of cache-efficiency on runtimes. It clearly shows that the I/O-efficient algorithms outperform the plane sweep algorithm as the input sizes increase. Recall our discussion that for the parameters of our systems $K = n/M$ and the I/O complexity of the distribution sweeping algorithm is $O((n/B)(1+\log_K n/M)) = O(n/B)$. This explains the non-linear asymptotic speedup over plane sweep algorithm (with I/O complexity of $O((n/B)\log n/M)$) as a function of the input size.
Figure 4: Speedup of the parallel distribution sweeping algorithms relative to the sequential distribution sweeping algorithm on configuration 2 (left) and configuration 1 (right) systems. The plots exclude the times to perform initial sorting of inputs by the y-coordinate.

Figure 4 shows the speedup that parallel distribution sweeping algorithm achieves relative to the sequential distribution sweeping algorithm.

Figure 5: Comparison of PEM and PRAM algorithms on 16 cores of configuration 3 is shown in the left figure. Running time and DRAM traffic for long segments on 12 cores of configuration 2 in the right.

**PRAM vs. PEM performance.** Figure 5 (left) shows the comparative performance of the various algorithms on configuration 3. We observe that the PRAM implementation is significantly slower than the PEM algorithm. For instance, with 51.2 million segments and the same number of queries, PRAM implementation takes 96 seconds with 16 cores, while the PEM implementation only requires 30 seconds with the same number of cores (excluding the time for loading the input and sorting it, which is 18 seconds for both implementations). This is largely accounted for by the fact that the PRAM implementation makes poor use of temporal locality and thus, has larger number of recursive levels. In each recursive level, it scans all the segments and query points, increasing the DRAM accesses significantly.

**DRAM Accesses and Cache Misses.** We could not find a reliable way to measure only L3 cache misses: the papi library does not support measurement of shared cache events, while the hardware counters for LLC (Last Level Cache) counters returned suspiciously similar results to L2 cache misses. Instead we measured the total traffic to DRAM using perf tool. Figure 5 (right) shows a clear correlation between the total DRAM traffic and running times. It is interesting to note that although our algorithms are designed in simple 2-level cache model, they minimize the
total traffic to DRAM, in spite of complex nature of modern memory systems.

**Random, short, medium and long segments.** We refer the reader to Appendix E for the relative behavior of the different point location algorithms on different segment types.

## 6 Conclusions And Future Work

In this work, we explored the effects of caches on actual run-times observed on various multicore architectures in the context of the geometric stabbing-max query problem. This is used to understand how accurately the PEM model predicts the running time of combinatorial algorithms on current multicore architectures. On single-socket multicore architectures, our results show a direct correlation between traffic on DRAM memory controller and running times of implementations. Thus, the algorithms designed I/O-efficiently via the (parallel) distribution sweeping framework outperform the plane sweep algorithms which do not address the I/O-efficiency.

We chose to perform our experiments on single-socket architectures, because the PEM model assumes uniform access latencies to shared memory. We conjecture that NUMA effects of DRAM access on multi-socket architectures might be better modeled by distributed computational models, where each processor copies/moves data into “local” memory — address space associated with its socket — before processing it. Once the data is in its “local” memory, one can use the PEM model to design algorithms to process the data. The experimental evaluation and modeling NUMA effects of multi-socket architectures is left for future investigations.

While we chose to implement an algorithm which was designed in the PEM model, it would be interesting to see how the implementations in other cache-conscious parallel models (for example, [8]) will fare in practice in similar setting.

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A Details of the PEM solution

In this section we present the details of the PEM solution.

As mentioned earlier, our algorithm is based on the parallel distribution sweeping framework introduced by Ajwani et al. [2]. Parallel distribution sweeping recursively divides the plane into vertical slabs, starting with the entire plane as one slab and in each recursive step dividing a given slab into \( K := \max\{2, \min\{\sqrt{n/P}, M/B, P\}\} \) child slabs. This division is chosen so that each slab at a given level of recursion contains roughly the same number of objects (segment endpoints and query points). The first level of recursion divides the plane into \( P \) slabs, each containing \( \Theta(n/P) \) input elements. Viewing the recursion as a rooted tree defines leaf invocations and children of a non-leaf invocation. An invocation on slab \( \sigma \) at the \( k \)th recursive level is denoted as \( I_k^{\sigma} \).

Each invocation \( I_k^{\sigma} \) receives as input a \( y \)-sorted list \( Y_k^{\sigma} \) containing segments and query points. The root invocation \( I_0^{\emptyset} \) contains all segments and query points of the input and the input \( Y_0^{\emptyset} \) is generated by sorting the horizontal segments and query points by the \( y \)-coordinate. For a non-leaf invocation \( I_k^{\sigma} \), let \( I_{\sigma_1}^{k+1}, I_{\sigma_2}^{k+1}, \ldots, I_{\sigma_K}^{k+1} \) denote its child invocations. The input \( Y_{\sigma_j}^{k+1} \) for a child invocation \( I_{\sigma_j}^{k+1} \) consists of the \( y \)-sorted list of segments in \( Y_k^{\sigma} \) with an endpoint in \( \sigma_j \) and the query points in \( \sigma_j \). In processing \( I_k^{\sigma} \), we consider all the children slabs of \( \sigma \): \( \sigma_1, \sigma_2, \ldots, \sigma_K \) and compute a segment \( l_{\sigma_j}^q \in Y_k^{\sigma} \) for all query points \( q \in \sigma_j \), that is the highest segment lower than \( q \) and spans \( \sigma_j \). If \( l_{\sigma_j}^q \) is higher than \( r_q \), we let \( r_q := l_{\sigma_j}^q \). Thus, \( r_q \) always stores the highest segment lower than \( q \) that spans \( \sigma_j \) (and all its parent slabs).

At every leaf invocation \( I_k^{\sigma} \), the highest segment below the query points are found using sequential I/O-efficient distribution sweeping technique as described in the previous section. This is then compared to the current stored value \( r_q \) and the maximum of the two values is stored in \( r_q \).

To compute the values of \( l_{\sigma_j}^q \) for all child slabs, we process them using the \( P \) processors as follows: We partition the input sets \( Y = \cup_i Y_i^{\sigma_i} \) into \( P \) equal chunks \( Y_1, Y_2, \ldots, Y_P \) based on the \( y \)-coordinate, each one of size \( \Theta(n/P) \). Then, processor \( p_j \) processes \( Y_j \) using the sequential I/O-efficient algorithm independently of others. Note, that this process sets the initial values of \( s_{\sigma_i} \) to \( l_{-\infty} = \emptyset \) – a virtual horizontal line \( y = -\infty \). Thus, if at the end of the process \( l_{\sigma_i}^q = l_{-\infty} \) for some query \( q \in Y_j \) and there is a segment \( s \in Y_k \), \( k < j \) such that \( s \) lies below \( q \), then \( l_{\sigma_i}^q \) is still not set correctly. To fix this, at the end of sequential sweep by each processor \( p_j \), \( p_j \) saves the values of \( s_{\sigma_i} \) for each of its slabs. These values are then processed by all processors in a way similar to segmented prefix sums to propagate the values of \( s_{\sigma_i} \) to the appropriate processors as follows. Assume, the segment identifiers of segments increase with the increase in segments’ \( y \)-coordinates. Now consider \( K \) independent prefix sums with max as the associative operator applied on the \( K \cdot P \) values of \( s_{\sigma_i} \) (one prefix sums on the values within a single slab). Finally, we initialize the value of \( s_{\sigma_i} \) at each processor \( p_j \) to the final value (after the prefix sums) of \( s_{\sigma_i} \) at processor \( p_{j-1} \) (in case of \( p_0 \), \( s_{\sigma_i} = l_{-\infty} \) for all \( \sigma_i \)) and repeat the sequential sweep by each processor. Note, the purpose of the prefix sums is to propagate the values of last seen segment across multiple processors and after the second sweep, all \( l_{\sigma_i}^q \) are set correctly.

The parallel I/O complexity of the above algorithm is \( O(\text{sort}_P(n)) \) I/Os.
B Data representation and space efficiency of our implementations

We implement the segments and queries as a single vector of objects. To achieve this, we represent each line segment and query point with a single 32-byte structure as follows. Using double precision for the coordinates, we need 16 bytes to represent a point on a plane. Additional 8 bytes are used for each segment to represent the x-coordinate of the second endpoint or for each query to record the y-coordinate of the segment below it as an answer. Additional 4 bytes are used for the identifier of each segment (we do not assume that segments have unique y coordinates, and therefore, they cannot be identified with this ID field). The same field entry is used for recording the segment ID of the output for a query. Finally, we need at least one boolean value for distinguishing a segment from a query. However, since memory allocation in C++ is aligned at 4 byte memory intervals, the last field of the structure takes up at least 4 bytes. While keeping segments and points as a single object type might cost extra 4 bytes of memory for each item, this approach has the advantage of simplifying computations and internal data structures for the slabs as we no longer need to keep separate lists for query points and line-segments, but can keep a single sorted list.

Let $s$ be the number of segments in the input and $q$ be the number of query points in the input, i.e., $n = s + q$. The sequential distribution sweeping can be implemented to use space taken up by up to $3s + 2q$ objects. This number arises from the fact that during distribution we might create up to two copies of each segment (one for each end point) to be placed into child slabs. In addition the distribution at each recursive level cannot be performed in place and, hence, we need to allocate additional $2s + q$ memory during the process. In practice to achieve this bound, one needs to know exactly how many objects will be distributed to each child slab. This would require an additional pass over the data to count the the sizes of each child slabs, increasing the running time by a factor of 2. Instead, we use dynamic arrays (e.g., vectors in C++ STL) which grow automatically when the data exceeds the preallocated capacity. During the resizing, the contents of vectors are copied over into the newly allocated vector, seemingly resulting in the same double the running time. However, copying a vector is performed using low level memory copying routines which are more efficient than traversing the input twice. In addition, we utilize the fact that we run our experiments on uniformly distributed data. By allocating 10% more space than $1/K$-th fraction of the total data to be distributed to $K$ children, with high probability no child exceeds the preallocated space.

C On the right choice of $P$

The PEM model prescribes algorithms for an architecture which contains a private cache per processor connected with an independent channel to the shared memory. The PEM algorithms measure the number of parallel I/Os performed. If the bandwidth of channels is fully utilized, increasing the number of processors without increasing the number of memory channels would not result in reduced parallel I/Os. In reality, modern multicores have much fewer memory channels than there are cores so an interesting question is whether it is beneficial to increase the number of processors beyond the number of memory channels.

The Intel i7 system contains 3 memory channels and implements hardware counters which record overall DRAM accesses per each channel. Interestingly, throughout the computation, one of the channels recorded no DRAM accesses, while the other two channels shared the traffic to DRAM
unevenly. The discrepancy in traffic load between the two channels decreased with higher number of cores used. We cannot explain the reason why the system did not use all memory channels and this is worth further investigations.

We could not measure the bandwidth utilization of the memory channels. However, we did measure the IPC – an average number of instructions executed per clock cycle. The results showed IPC on the Intel i7 system being close to 1 regardless of number of cores used for the distribution sweeping implementation. However, for the planesweep implementation the IPC dropped down to .7 for large inputs. This leads us to believe that the random access of planesweep results in inefficient use of the memory bandwidth. At the same time full cache line transfers of the distribution sweeping implementation barely saturates the memory channels even with all 4 cores running 2 threads each.

D Effect of parameters $K$ and $M$ on runtime.

Figure 6 shows how much each task takes as a fraction of total runtime as we vary the threshold parameter $M$. One can see that by reducing $M$ down to the L2 cache size results in faster combined execution time of all invocations of the plane sweep algorithm at the base case of the recursion. This is due to the fact that the binary search tree $T$ used for the plane sweep algorithm fits in the faster L2 cache.

However, this decrease in runtime of the base case is offset by the increase in the runtime of the distribution sweeping phase due to the following reason. Recall that $K = \min\{n/M, M/B\}$. If $K = n/M$, larger value of $M$ results in increase of $K$ – the number of slabs to distribute the objects at each recursive level. While the number of slabs is still small enough to fit in cache, and, therefore, there is no increase in cache misses, the time it takes to identify the slab where to distribute each item takes $O(\log K)$ internal computation time, i.e. it grows with $K$. If, on the other hand, $M$ decreases so much that $K = M/B$, the number of recursive levels grows as a function of $\log_K(n/M)$. Thus, decrease in $K$ results in more recursive levels which in turn results in more scans of the input and, therefore, more (capacity) cache misses.

We also observed that setting $M$ equal to exactly the size of a cache does not result in the best
run times. This can be explained by the more complex nature of caches, such as set associativity and the replacement policy: the external memory model assumes fully associative cache with optimal replacement policy, while modern architectures implement set-associative caches with (most likely\(^4\)) the Least Recently Used (LRU) replacement policy. To achieve the best results we set \(M\) to a quarter of the L3 cache size for the Intel i7 architecture and a third of the L3 cache size for the AMD Opteron architectures.

Since the caches are shared among subsets of processors, in case of the parallel execution, our initial intuition was to reduce \(M\) by the number of processors sharing the cache. However, our experiments showed that this is unnecessary and the same \(M\) as for the sequential implementation works just as well in the parallel implementation.

This can be explained by the fact that in our experiments, \(K = n/M\) which constitutes a much smaller portion than the L3 cache. Therefore, during the parallel distribution sweeping, maintaining one block for each child slab in cache and maintaining for each processor the tree \(T'\) of size \(2 \cdot K\) for work-optimal distribution sweeping does not interfere with other processors’ cache data.

E Effect of segment type on the performance of point location algorithms

To analyze the behavior of our algorithms on segments of varying lengths, we first generate the line-segments in different ways:

**Random segments:** Our first input set is a set of random line-segments in the grid. The random lines are generated by selecting a random y-coordinate and two random x-coordinates in the grid. Thus, the expected length of the line segments is \(O((\text{Grid Size}))\).

**Short segments:** Here, \(n\) line segments are generated with length chosen uniformly at random between Grid Size/\(n\) and \(4 \cdot \text{Grid Size}/n\).

**Medium segments:** Here, \(n\) line segments are generated with length chosen uniformly at random between Grid Size/\(\sqrt{n}\) and \(4 \cdot \text{Grid Size}/\sqrt{n}\).

**Long segments:** Here, we generate line segments with lengths chosen uniformly at random between Grid Size/4 and \(3 \cdot \text{Grid Size}/4\).

In Figures 7 through 9 one can see how different sizes of segments affect the running times of our different implementations. We observe that the plane sweep (Figure 7) performs much worse on the long segments than on the short segments. This is because the expected number of short segments intersecting any vertical line is expected to be constant and the set of active segments \(A\) fits in cache at all times. Thus, the updates and predecessor queries on \(A\) do not incur any additional cache misses. On the other hand, the number of long segments intersecting any vertical line is expected to be linear with the input size and the traversal of \(T\) will incur a lot of cache misses. Contrast this with the runtimes for sequential (Figure 8) and parallel (Figure 9) distribution sweeping algorithms for the different segment sizes, which show much smaller variance of runtimes as a function of different segment lengths. This confirms that the results that Chiang [11, 12] observed 17 years ago still hold on modern architectures.

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\(^4\)It is hard to determine the true replacement policy because processor manufacturers keep this information confidential.
Figure 7: Runtimes for small, medium, long and randomly sized segments for the plane sweep on AMD Opteron 6168 (left) and Intel i7 (right) systems as a function of input size.

Figure 8: Runtimes for small, medium, long and randomly sized segments for the sequential distribution sweeping on AMD Opteron 6168 (left) and Intel i7 (right) systems as a function of input size.

Figure 9: Runtimes for small, medium, long and randomly sized segments for the parallel distribution sweeping on AMD Opteron 6168 (left) and Intel i7 (right) systems as a function of input size. The results are for the maximum number of threads for each system.