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To cite this version:
Ram A Raghunathan, Stefan A Muller, Umut A Acar, Guy A Blelloch. Hierarchical Memory Management for Parallel Programs. Proceedings of the 21st ACM SIGPLAN International Conference on Functional Programming, Sep 2016, Nara, Japan. 10.1145/3022670.2951935. hal-01416237

HAL Id: hal-01416237
https://inria.hal.science/hal-01416237
Submitted on 14 Dec 2016

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Hierarchical Memory Management for Parallel Programs

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Abstract
An important feature of functional programs is that they are parallel by default. Implementing an efficient parallel functional language, however, is a major challenge, in part because the high rate of allocation and freeing associated with functional programs requires an efficient and scalable memory manager.

In this paper, we present a technique for parallel memory management for strict functional languages with nested parallelism. At the highest level of abstraction, the approach consists of a technique to organize memory as a hierarchy of heaps, and an algorithm for performing automatic memory reclamation by taking advantage of a disentanglement property of parallel functional programs. More specifically, the idea is to assign to each parallel task its own heap in memory and organize the heaps in a hierarchy/tree that mirrors the hierarchy of tasks.

We present a nested-parallel calculus that specifies hierarchical heaps and prove in this calculus a disentanglement property, which prohibits a task from accessing objects allocated by another task that might execute in parallel. Leveraging the disentanglement property, we present a garbage collection technique that can operate on any subtree in the memory hierarchy concurrently as other tasks (and/or other collections) proceed in parallel. We prove the safety of this collector by formalizing it in the context of our parallel calculus. In addition, we describe how the proposed techniques can be implemented on modern shared-memory machines and present a prototype implementation as an extension to MLton, a high-performance compiler for the Standard ML language. Finally, we evaluate the performance of this implementation on a number of parallel benchmarks.

Categories and Subject Descriptors  D.3.4 [Programming Languages]: Processors — Memory management (garbage collection); D.1.3 [Programming Techniques]: Concurrent Programming — Parallel Programming; D.4.1 [Operating Systems]: Process Management — Scheduling

General Terms  Languages, Performance

Keywords  Languages, Memory Management, Parallelism, Scheduling

1. Introduction
In the past two decades, there has been a large body of work on both parallel garbage collection and parallel scheduling. In this paper, we propose an approach for coupling the two together, which is particularly useful in functional languages. The basic idea is to organize memory in a way that mirrors the structure of parallelism in the computation and takes advantage of the independence of parallel tasks to perform garbage collection. While this general idea may be applicable to the broader array of parallel functional programming languages, we focus here on strict languages, such as the ML family, extended with support for nested parallelism.

More specifically, instead of using a single shared heap [6, 9, 21, 26, 32] or a two-level heap consisting of local processor heaps and a global heap [5, 8, 23, 25, 31, 37], we suggest a hierarchy of heaps (tree of heaps) that is tied to the nesting of tasks. In nested parallel (fork-join) parallelism, the tasks form a natural nesting, and memory that is allocated and referenced has a structure that is related to this nesting, which we would like to leverage. In strict functional computations, pointers only point up the hierarchy, creating what we refer to as a disentangled heap. Furthermore, what is often understood as temporal locality in sequential programs becomes nested task locality in nested parallel computations. For example, in divide-and-conquer algorithms, each node in the recursion tree has some locality within its subtree [14, 45].

The motivation for organizing the heap into a hierarchy is three-fold. Firstly, it allows for simpler parallel collection since the nodes in the tree can be collected individually and separately, especially if they can be kept disentangled. Secondly, the hierarchical organization of heaps is better suited for supporting memory locality at multiple levels of a modern cache hierarchy, for which some of the higher levels are shared among cores on multiprocessors. By having the size of heaps grow going up the levels, and moving values up the heap as they are collected, the approach can be seen as the natural extension of a multi-level generational collector. Finally, the organization allows us to naturally tie memory management decisions to scheduler decisions. For example, a good time to collect is probably when a task finishes, or a good time to create a heap in the hierarchy is likely when stealing a task.

We propose a specific memory manager and garbage collection algorithm based on the hierarchical heaps and disentanglement. Rather than going directly to a low-level description of the algorithms, we formalize a functional, ML-like language with nested parallelism and a memory manager by presenting a reasonably high-level operational semantics that accounts for all the key aspects of the memory management and collection, while abstracting away from the details. The semantics models parallelism using non-deterministic interleaving of parallel steps. It captures a parallel fork by “activating” a parallel tuple. This is meant to capture the lower-level idea of scheduling a parallel task (e.g. stealing the task in a work-stealing scheduler), and allows for lazy task creation (i.e. if
we present a high-level, informal overview of the ideas proposed in this paper. The rest of the paper makes these ideas precise.

2. Overview

We present a high-level, informal overview of the ideas proposed in this paper. The rest of the paper makes these ideas precise.

We use the quicksort algorithm as an example and assume an ML-like, strict, purely functional language with nested parallelism, where the primitive par creates a tuple by allowing its components to be evaluated in parallel. Figure 1 shows the code for quicksort in such a language. For simplicity, we assume that the input contains no duplicates. The function qsort uses the first item as a pivot to partition the input list into less-than and greater-than parts, written \( l1 \) and \( l2 \), sorts them recursively in parallel, and concatenates the sorted lists to compute the final sorted output.

**Task Tree.** We can represent an evaluation of the quicksort algorithm with a task tree consisting of parallel tasks where each task corresponds to a parallel sub-computation as indicated by the par primitive. Figure 2 shows such a tree representation of the evaluation of the qsort function on the input list \([5; 6; 3; 4; 1; 9]\). The nodes of the task tree represent parallel tasks and the edges represent the control dependencies between them. For example, the root task, labeled \( T_A \), which corresponds to the first call to the function, forks two new tasks \( T_B \) and \( T_C \).

```
fun qsort l =
  case l of
    nil ⇒ nil
  | h::nil ⇒ h:
  | h::t ⇒
    let
      \( (l1,l2) = \text{partition } h t \)
      \( \text{in} \)
      append sll (h::slg)
    in
  end
```

**Figure 1.** The code for quicksort.

**Figure 2.** An example run of quicksort. The nodes represent (parallel) tasks and the edges represent the control dependencies between them.

**Memory as a Heap Tree and Disentanglement.** Classically, memory is viewed in programming languages as a flat, unstructured pool of objects, or a fixed hierarchy of several heaps, referred to as generations. In this paper, we propose an approach where we structure memory as a dynamic hierarchy by partitioning memory into heaps, each of which corresponds to a task in the task tree. To create this hierarchy, we dedicate a fresh heap to each forked task and allocate all memory requested by the task in its heap, which the task owns. When the task completes, we join its heap with that of its parent. As a result, we can organize the memory at any point during execution as a hierarchy, represented as a tree of heaps as follows. Let \( H_A \) and \( H_B \) be two heaps. If there is an edge between their owner tasks, \( T_A \) and \( T_B \), respectively, then insert an edge between \( H_A \) and \( H_B \). Otherwise, don’t insert any edge. We refer to the resulting tree on the heaps as a heap tree. The motivation behind our approach is to organize memory in a way that reflects the structure—specifically the dependencies—of the computation.

Figure 3 illustrates the task and the heap trees for our example. Each “qsort” node represents a task. The gray box around the node represents its heap. Tasks and the edges between them define the task tree. Heaps and the edges between them define the heap tree. The figure also shows the data stored in each heap. The root heap, \( H_3 \), contains the input list \( l0 \) and the two lists obtained via partitioning, \( l1 \) and \( l2 \). The task \( T_B \) takes the input \( l1 \) from the root heap \( H_A \) and partitions it into two new lists \( l3 \) and \( l4 \). The task \( T_C \) takes the input \( l2 \) from the root heap \( H_A \) and produces the lists \( l5 \) and \( l6 \).

Having organized memory as a tree of heaps, we observe the following key disentanglement property. Let \( T_A \) and \( T_B \) be any two tasks in the task tree and let \( H_A \) and \( H_B \) be their heaps in the heap tree. If some object in \( H_A \) references (or points to) an object in \( H_B \) at any time during execution, then \( H_A \) is an ancestor of \( H_B \) in the heap tree. Intuitively speaking, this property holds because, in a purely functional language, a parallel task has access only to the objects that have been allocated before its execution or by the task itself, excluding the objects that are allocated by tasks that might have been executed in parallel (even if they are not actually executed in parallel). Such objects are allocated precisely by the tasks that are on the path from the task to the root of the task tree. Since all memory allocated by a task is placed in the heap of that task, and since there
is an isomorphism between the heap tree and the task tree as defined by the ownership relation, the disentanglement property holds. As a corollary, we can establish the following: if \( T_b \) is neither an ancestor nor a descendant of \( T_b \), then objects in \( H_b \) do not reference objects in \( H_b \) and objects in \( H_b \) do not reference objects in \( H_b \). In other words, if \( T_b \) and \( T_b \) are concurrent tasks that may be executed in parallel, then \( H_b \) and \( H_b \) are guaranteed to be “disentangled”. As suggested by disentanglement, in the quicksort example shown in Figure 3, all pointers in the memory point “up” in the heap tree: if an object in heap \( H_b \) points to another object in another heap \( H_b \), then \( H_b \) is an ancestor of \( H_b \).

In Sections 3.1 and 3.2, we make precise this high-level description of hierarchical heaps and disentanglement by presenting a semantics that restricts a task’s memory access to its own heap and ancestor heaps. We prove that a program evaluated with hierarchical heaps and restricted memory access has the same behavior as it would have under a standard operational semantics, thus establishing the disentanglement property.

Garbage Collection. When considered in the context of the heap tree, the disentanglement property leads us to observe that any leaf heap \( H \) in the heap tree can be collected by the task \( T \) that owns that heap independently of all the other heaps and tasks. This holds because, via disentanglement, \( H \) does not have any “roots” or pointers into it. Thus, garbage collection of \( H \) can proceed as all the other tasks are executing. A moving (copying) or a non-moving collection can be used. We can generalize this observation to any subtree in the heap tree as follows: the heaps in any subtree of the heap tree can be collected independently of all other heaps. As with a leaf of the heap tree, a subtree of heaps has no incoming pointers from outside the subtree, and thus the collection can proceed concurrently with all other tasks operating on the other heaps. This means that we can garbage collect a subtree of heaps by stopping only those tasks that own them as we allow other tasks to continue executing. Furthermore, the subtree can be garbage collected in parallel by taking advantage of disentanglement: all the leaves can be collected in parallel and, after their collections are completed, the parents of the leaves may be collected in parallel, and so on until the root is reached.

For example, in Figure 3, we can garbage collect heaps \( H_b, H_b, \) and \( H_b \), which form a subtree rooted at \( H_b \), by stopping only the tasks \( T_b, T_b, \) and \( T_b, \) which own the heaps. To garbage collect the subtree, we can collect \( H_b, H_b, \) and \( H_b \) in parallel, and then \( H_b \).

In Section 3.3, we make precise this informal description of garbage collection and establish its correctness by showing that collection does not free rechable locations (memory safety) or alter the behavior of the program (meaning preservation).

### 3. The Language

We present an ML-style functional calculus called \( \Lambda^{HP} \) and an operational semantics that accounts for memory allocation and the scheduling of tasks, and ensures the property that tasks access memory only in their own heap and ancestor heaps. We then show that any \( \Lambda^{HP} \) program is disentangled in that it may be run safely and that it is both deterministic and equivalent to a standard “flat” semantics where memory is treated conventionally. We then formalize a hierarchical memory manager on top of \( \Lambda^{HP} \) and prove its correctness.

#### 3.1 A Core Calculus for Disentanglement

**Abstract syntax.** Figure 4 shows the syntax of the language \( \Lambda^{HP} \). The types include natural numbers (as the sole base type), functions and products. To present a precise accounting of memory operations, we distinguish between large values \( v \), which are always allocated in the heap, and small values consisting only of memory locations, written \( \ell \). Large values include natural numbers, named recursive functions and pairs of heap locations. Large values are not irreducible, but rather step to heap locations \( \ell \), which are the only
irreducible values of the language. Since small values consist only of locations, we use the term “value” to refer to a large value and “location” to refer to a small value.

In \( \lambda^{HP} \), we organize memory as a tree of heaps and restrict expressions to access only part of the memory (along a path going up the hierarchy). A heap, written \( H \) (and variants), is simply a mapping from locations to (large) values. A heap path or simply a path, written as \( P \) (and variants), is a list or stack of heaps. A path represents a leaf-to-root path in the heap tree; it is the collection of heaps to which an expression has access.

Expressions of \( \lambda^{HP} \) include locations and (large) values, introduction and elimination forms for the standard types such as function application, projection and pair creation, and parallel tuples \( \langle e_1, e_2 \rangle \). Parallel tuples are evaluated sequentially by default, but may be activated for parallel evaluation. An active parallel tuple, \( \langle T_1, T_2 \rangle \), consists of two parallel tasks, \( T_1 \) and \( T_2 \), which correspond to the components of the tuple.

A task, \( H \cdot e \), consists of an expression \( e \) and a heap \( H \). The heap \( H \) serves as the portion of the memory that is private to the expression \( e \), i.e., \( H \) is visible only to the expression \( e \) (including subexpressions of \( e \)). The reader may wonder why we introduce tasks as a level of indirection in the syntax since they can take only one form \( (H \cdot e) \). When we formalize the memory manager, we will introduce another form of task, and the separation will make the presentation cleaner, so we adopt it now.

Statistics. Figure 5 illustrates the static semantics of the calculus. Apart from the treatment of memory, the statics semantics of \( \lambda^{HP} \) is relatively standard. We first describe the typing of heaps and paths.

We ascribe type signatures to heaps and paths. A signature \( \Sigma \) is a mapping from locations to types. The heap typing judgment 

\[ H \vdash \ell \vdash \ell_{\text{fresh}} \]


\[ H \vdash \ell \vdash \ell_{\text{fresh}} \quad \text{D-TypE} \]

\[ H \vdash e \vdash e' \vdash e'_{\text{fresh}} \quad \text{D-TypS} \]

\[ H \vdash e \vdash e' \vdash e'_{\text{fresh}} \quad \text{D-TypA} \]

\[ H \vdash e \vdash e' \vdash e'_{\text{fresh}} \quad \text{D-TypS} \]

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\[ H \vdash e \vdash e' \vdash e'_{\text{fresh}} \quad \text{D-TypA} \]

\[ H \vdash e \vdash e' \vdash e'_{\text{fresh}} \quad \text{D-TypS} \]

\[ H \vdash e \vdash e' \vdash e'_{\text{fresh}} \quad \text{D-TypA} \]
rule D-Alloc allocates a fresh location \( \ell \) in the heap \( H \), binds the location to \( v \), and returns \( \ell \).

The rules for function and product types are standard for a call-by-value calculus with left-to-right evaluation. Function application and projection evaluate the subexpressions down to heap locations, look up the locations in the path, and perform the operation. For technical reasons, the function application rule requires that the location \( \ell \) be bound, but doesn’t immediately use the value.

Sequential and inactive parallel tuples evaluate the first component to a location, followed by the second. If both components of a parallel tuple are evaluated fully before the tuple is activated, it changes into a pair \( (\ell_1, \ell_2) \) (rule D-ParE). A parallel tuple may also be activated at any time, non-deterministically, by rule D-Activate. This models the possibility that another processor may begin evaluating one of the components in parallel. The activation rule changes the tuple into an active parallel tuple with two new tasks (each with a new, initially empty, private heap). Once a parallel tuple is activated, either component task may evaluate at any time using rules D-ParAS1 and D-ParAS2. (The rule D-ParAS2 does not require \( e_1 \) to be fully reduced to a location.)

The transition judgment for tasks, \( T \Rightarrow P T' \) indicates that, under path \( P \), task \( T \) steps to \( T' \). It doesn’t include a private heap since that is internal to the task. Only the internal private heap may be modified, so \( P \) is read-only. Rule D-TaskStep (Figure 7) allows \( e \) to step using the expression step judgment.

When a component of a parallel tuple is stepped (the premises of D-ParAS1 and D-ParAS2), the heap for the tuple, \( H \), is consed onto the current path and becomes part of the shared memory. This means that a heap becomes read-only when it is not at a leaf of the heap tree. When both components of an active parallel tuple are evaluated down to locations, the tuple steps to an ordinary pair using rule D-ParAE and merges the two local heaps into the parent heap.

### 3.2 Disentanglement

We state and prove one of this paper’s key contributions: that the instance of this paper’s key contributions: that

\[
\ell \rightarrow_v \tau \quad \text{then there exists a value } v \text{ such that } P; \ell \rightarrow_v v \quad \text{and } \Gamma \vdash_{\text{local}} \ell ; v : \tau.
\]

**Proof.** By construction.

The second lemma is an important property of heaps, which states that a merged heap is well-typed.

\[
\text{Lemma 2.} \quad H : \Sigma \text{ and } H' : \Sigma', \text{ then } (H \uplus H') : \Sigma, \Sigma'.
\]

**Proof.** By induction on the derivation of \( H' : \Sigma' \). If \( H' = \emptyset \), then the result is trivial. Otherwise, \( H' = H'[\ell \rightarrow v] \) and \( \Sigma' = \Sigma'', \ell : \tau \). By induction, \((H \uplus H'') : \Sigma, \Sigma''\). By the heap typing rules and weakening, \( H \uplus H' = (H \uplus H'')[\ell \rightarrow v] : \Sigma, \Sigma'' \), \( \ell : \tau = \Sigma, \Sigma' \).

We now state and prove the preservation and progress lemmas.

\[
\text{Lemma 3 (Preservation).} \quad \Gamma \vdash_{\Sigma} e : \tau \text{ and } H : \Sigma_1 \text{ and } P : \Sigma_2 \text{ and } H ; e \rightarrow_{P} e' ; H'; \text{ then } H' : \Sigma_1' \text{ where } \Sigma_1' \text{ is an extension of } \Sigma_1, \text{ and } \Gamma \vdash_{\Sigma} e : \tau.
\]

**Proof.** By induction on the derivation of \( H ; e \rightarrow_{P} e' ; H' \).

**Lemma 4 (Progress).** \( \vdash_{\Sigma} e : \tau \text{ and } H : \Sigma_1 \text{ and } P : \Sigma_2 \text{ then either } e \text{ is a location or there exist } e' \text{ and } H' \text{ such that } H ; e \rightarrow_{P} e' ; H' \).

**Proof.** By induction on the derivation of \( \vdash_{\Sigma} e : \tau \).

### Correspondence with flattened semantics

Figure 8 defines the flattening transformation, which converts a \( \lambda^{HP} \) expression into a flat expression with no nested tasks and no free heap locations (but which is still a valid expression in the \( \lambda^{HP} \) syntax). The judgment \( \llbracket e \rrbracket_p \rightarrow \hat{e} \) indicates that expression \( e \) flattens to \( \hat{e} \) under path \( P \). The transformation is defined inductively. There are two important rules to note. Locations are looked up in the path and the resulting value is recursively flattened (this will terminate since heaps and heap paths have no cycles). The flattened value is substituted for the location. The rule for active parallel tuples recursively flattens the two subexpressions, adding the local heaps to the path, and then converts the active tuple into an inactive tuple, thus flattening the hierarchy.

We make three observations about the flattening transformation:

- If \( e \) contains no active parallel tuples or heap locations (it may be a source program that has not yet begun executing), then \( \llbracket e \rrbracket_p \rightarrow_1 e \).
- If \( e \in \text{dom}(P) \) and \( \llbracket e \rrbracket_p \rightarrow_1 \hat{v} \), then \( \hat{v} \) is simply the value at \( e \) “lifted” so that it is closed with respect to the heap.
- Flattening is deterministic: if \( \llbracket e \rrbracket_p \rightarrow_1 \hat{e}_1 \) and \( \llbracket e \rrbracket_p \rightarrow_1 \hat{e}_2 \), then \( \hat{e}_1 = \hat{e}_2 \).

Flattened expressions may be run using the simple operational semantics of Figure 9. No path or heap is required, since these expressions have no heap locations. There is now no notion of “active” parallel tuples, and instead (formerly “inactive”) parallel tuples may evaluate in parallel using rules F-ParS1 and F-ParS2. While these parallel evaluation rules allow for nondeterministic interleavings, evaluation is nevertheless deterministic. As usual, we show this using an intermediate result, the diamond lemma, which states that if an expression steps to two different expressions in one step, they can be “brought back together” in one step. We do not immediately prove confluence, the generalization of the diamond property to multi-step evaluation. This will be part of the disentanglement theorem.

**Lemma 5 (Diamond Lemma).** \( e \rightarrow e_1 \text{ and } e \rightarrow e_2 \text{, then there exists } e' \text{ such that } e_1 \rightarrow e' \text{ and } e_2 \rightarrow e'. \)
Lemma 9. Suppose that $\hat{e} \rightarrow e'$ and $||e||_{H, F-P} \rightarrow \hat{e}$. Then there exist $H'$ and $P$ such that $H; e \rightarrow P; e'; H'$ and $||e'||_{H', P} \rightarrow e'$.

Proof. By induction on the derivation of $e \rightarrow e'$. □

Disentanglement. Finally, we state and prove the disentanglement theorem, which formalizes the two properties (type safety and correspondence with flattened semantics) outlined at the beginning of this section. If a source program $e$ is well-typed and it evaluates to $e'$, then (1) $e'$ is not stuck and (2) for any flattened evaluation of $e$ to $\hat{e}'$ ($e'$ and $\hat{e}'$ need not be related), $e'$ and $\hat{e}'$ may be brought back together in the style of confluence.

Theorem 1 (Disentanglement). Suppose $e$ is a source program with no active parallel tuples. If $\vdash e; \tau$ and $e \rightarrow_{s} e'; H'$, then

1. Either $e'$ is a location or there exists $e''$ such that

$$H'; e' \rightarrow_{s} e''; H''$$

and $\hat{e}' \rightarrow e''$ and $||e''||_{H''} \rightarrow \hat{e}'$.

Proof. 1. We show by induction on the derivation of $e; \rightarrow_{s} e'; H'$. Suppose that $H'; \rightarrow_{s} e'; H'$ for some $\Sigma$ and that $\vdash_{s} e'; \tau$. By this, the result follows from Lemma 4. If $e' = e$ and $H' = \emptyset$, then $H'; \rightarrow_{s}$ and the typing result on $e'$ was assumed. Otherwise, $e \rightarrow_{s} e'$ and $H_{0}, e'_{0}; \rightarrow_{s} e'; H'$. By induction, there exists $\Sigma_{0}$ such that $\vdash_{s} H_{0}; \Sigma_{0}$ and $\vdash_{s} e'_{0}; H'$. By typing, the result for $e'$ follows from Lemma 3. 2. We have $||e||_{\emptyset} \rightarrow e$. By inductive application of Lemma 8, we have $e \rightarrow_{s} e''$ and $||e''||_{H''} \rightarrow e''$. By a standard inductive application of Lemma 5, there exists $e'''$ such that $e'' \rightarrow_{s} e'''$ and $e''' \rightarrow e''$. By inductive application of Lemma 9 to the latter derivation, there exist $e'''$ and $H''''$ such that $H'; e'' \rightarrow_{s} e'''; H''''$ and $||e''''||_{H''''} \rightarrow e''''$. □

Part (2) of Theorem 1, as well as its proof, is illustrated graphically in Figure 11. This statement is surprisingly powerful. In particular, it immediately implies that if $e$ terminates with flattened semantics, it terminates with hierarchical semantics, and vice versa, that both semantics are deterministic and that they both produce the same result up to flattening. Each of these statements is simply an instantiation of Theorem 1.

Corollary 1. If $e \rightarrow_{s} \hat{e}$, then there exist $e; H$ such that $\vdash e; \rightarrow_{s} e; H$ and $||e||_{H} \rightarrow \hat{e}$. 2. If $\emptyset; e \rightarrow_{s} e; H$, then there exists $\hat{e}$ such that $e \rightarrow_{s} \hat{e}$ and $||\hat{e}||_{H} \rightarrow \hat{e}$.

3.3 Hierarchical Garbage Collection

We describe a hierarchical garbage collector, dubbed HGC, that takes advantage of disentanglement to perform concurrent garbage collection at the granularity of tasks. The collector HGC can collect any heap in the heap hierarchy by stopping the task that owns the heap along with its descendant tasks, while all other (non-descendant) tasks continue executing. Any number of independent heaps, which are not descendants and ancestors of each other, can be collected in parallel. The collector HGC is thus able to perform garbage collection without incurring additional synchronization.
The collector we formalize is a semispace copying collector, though whose
state of collection, in which some locations have already been copied
collection. The tuple represents a heap which is in an intermediate
to dynamic rules to perform garbage collection on tasks.
other garbage collection algorithms could be used as well. Since the
subsection is the proof of correctness for HGC.

\[
\begin{array}{c}
\text{F-ArrE} & e_1 \to e'_1 \\
\text{F-PasE} & e_1 \to e'_1 \\
\text{F-PasS1} & e_2 \to e'_2 \\
\text{D-FstE} & e \to e' \\
\end{array}
\]

Figure 9. The dynamics for the flattened semantics.

\[
\begin{array}{c}
\text{F-ArrS1} & e_2 \to e'_2 \\
\text{F-PasS1} & e_2 \to e'_2 \\
\text{D-FstS} & e \to e' \\
\text{D-SndE} & e \to e' \\
\end{array}
\]

Figure 10. An illustration of the relationship between hierarchical and flattened evaluation. Straight lines are (hierarchical or flattened) evaluation and curved lines indicate the flattening transformation.

Between the mutator and the collector. The main result of this
subsection is the proof of correctness for HGC.

We formalize the outlined hierarchical garbage collector for \(\lambda^{HP}\).
The collector we formalize is a semispace copying collector, though other garbage collection algorithms could be used as well. Since the structure of a \(\lambda^{HP}\) program already contains hierarchical tasks and heaps, formalizing the collection algorithm requires few changes to
the semantics. In fact, we simply need to introduce a new form for
tasks to represent a task which is locked for collection, and several
dynamic rules to perform garbage collection on tasks.

Our approach in formalizing the collection algorithm is inspired
by Morrisett et al. [40]. While they analyze only stop-the-world collectors for sequential languages and their formalisms do not capture some of the details that ours does, we borrow their language-based approach for high-level descriptions of collection algorithms as well as much of their notation.

Syntax. The first change is to introduce a new form of task, as shown at the top of Figure 12. The new form, \(\langle H_j; S; H; F \rangle [e]\), represents an expression \(e\) which is locked to perform garbage collection. The tuple represents a heap which is in an intermediate state of collection, in which some locations have already been copied to the to-heap \(H_j\) and the rest remain in the from-heap \(H\). The second component, \(S\), is the scan set or frontier, the set of locations that have been seen by the collector but not yet copied. The final component, \(F\), is a forwarding map. It is a finite map whose domain
and codomain both consist of heap locations, and it maps locations formerly bound in \(H_j\) to the locations in \(H\) to which they have been copied. This allows us to make sense of values in \(H_j\), which may refer to locations which have already been renamed and copied to \(H\). The forwarding map models forwarding pointers, which are commonly used to perform this function in practical copying garbage collectors. For notational convenience, we implicitly extend the domain of forwarding maps by mapping locations that are not in the domain to themselves:

\[
F(\ell) := \begin{cases} 
F(\ell) & \ell \in \text{dom}(F) \\
\ell & \ell \notin \text{dom}(F).
\end{cases}
\]

We will also use the notation \(F(e)\) to denote an expression in which every free location \(\ell\) has been replaced with \(F(\ell)\). We extend \(F\) to operate on heaps in the following way:

\[
F(H) := \{ F(\ell) \mapsto F(v) \mid \ell \mapsto v \in H \}
\]

and on signatures in the following way:

\[
F(\Sigma) := \{ F(\ell) \mapsto F(v) \mid \ell \mapsto v \in \Sigma \}
\]

A particularly common usage of the forwarding map is the notation \(\ell \mapsto \ell'\) (and similar for heaps and signatures), in which \(e\) is forwarded with the singleton forwarding map forwarding \(\ell\) to \(\ell'\). In addition, we use \(F[\ell \mapsto \ell']\) to indicate the extension of the map \(F\) with a mapping from \(\ell\) to \(\ell'\).
Statics. The typing rule for the new form of task is given in Figure 12. The rule uses the rules for typing heaps to assign a signature to the concatenation of the from- and to-heaps. Since values in \( H \) might refer to locations that have already been copied, we forward the values of \( H \) before concatenating with \( H' \). The expression must be well-typed under the concatenation of the global signature \( \Sigma \) and this new signature \( \Sigma \). In addition, the typing rule also enforces a key invariant of garbage collection: that the scan set \( S \) is equal to the set \( FL(H_1, e) \) of locations that appear free in either \( e \) or a binding of \( H_1 \) and are not bound in \( H_1 \). Formally,

\[
FL(H_1, e) \equiv FL(e) \cup \bigcup_{(\ell \in dom(H_1))} FL(H_1(\ell)) \cap dom(H_1)
\]

Dynamics. Note that if a collection is in progress and the task is of the form \( H \cdot e \), no rule allows \( e \) to stop, so no changes are required to the semantics to ensure that \( e \) is in fact locked during collection. Instead, we introduce a new transition rule for tasks, D-GCStep, which allows the task to proceed with collection. The auxiliary judgment \( T \longrightarrow_{gc} T' \) indicates that a step of garbage collection transforms \( T \) into \( T' \).

Our garbage collection algorithm uses a small-step semantics, where one step (atomically) copies a single location. A full garbage collection will therefore involve many individual steps, which can be freely interleaved with evaluation or collection on other processors. Figure 12 gives the transition rules for collection. Rule HGC-S stacks GC locks for collection and sets up the tuple by using the existing heap, \( H \), as the from-heap, initializing the to-heap and the forwarding map to be empty and the scan set to be \( FL(e) \), the roots of the expression. Rule HGC-Cory performs one step of collection. The rule takes a location from \( S \) which is present in \( H \), copies its value \( v \) and adds to \( S \) the locations of \( v \) which have not already been copied. Rule HGC-EnvGC requires that \( dom(H_1) \cap S = \emptyset \), i.e., that there are no remaining locations to be collected. If this is the case, \( e \) is unlocked and the to-heap becomes the new heap.

Rule HGC-S stacksGC can apply to a task nondeterministically at any time, and so our semantics does not specify when collections can or should occur. Also note that a collection in the semantics can collect any heap in the hierarchy at any time. An entire subtree is locked, but only the heap at the root of that subtree is collected. However, this property also means that, because our safety proofs are parameterized over any possible evaluation, our theorems will allow for a variety of possible policies for both when to trigger a collection and what subtrees to collect at what times. The policy for collecting a subtree described in Section 2 may be reproduced in the semantics by locking and collecting all leaves of a subtree, then immediately locking and collecting their parents, and so on until the root of the subtree is collected. Because any collection occurs over a number of single steps, unrelated tasks can execute or perform garbage collection while a collection is taking place. Thus, we effectively model parallel collection.

3.3.1 Safety and Correctness Proofs

We will now proceed to prove that our collection algorithm is memory safe, that is, it never frees a location that will be dereferenced later. Because our calculus is typed and dereferencing a freed location is a stuck state, memory safety is an immediate corollary of type safety. Thus, we only need to update the progress and preservation proofs to handle the new rules. Several lemmas are required. One important lemma is that garbage collection preserves typing, in that if \( e \) is well-typed under heap \( H \) and \( H' \) and \( H \) is discarded by garbage collection, \( e \) is also well-typed under heap \( H' \). The signature of the higher levels of the hierarchy, notated \( \Sigma_1 \) in the lemma, does not change. Both \( H \) and \( H' \) are typed under \( \Sigma_1 \). We also allow the expression to refer to additional locations in another signature, \( \Sigma_2 \), this contains locations at levels lower than the level being collected. In using this lemma to prove preservation, we will only be interested in the case where \( \Sigma_2 \) is empty because we want to show type preservation of the expression whose heap is being collected. However, \( \Sigma_2 \) is necessary for the proof of the lemma, when inducing on expressions containing nested heaps.

**Lemma 10.** If \( \cdot, H \bowtie H' \bowtie \Sigma \) and \( \cdot, H' \bowtie \Sigma' \) and \( \Gamma \bowtie_1, \bowtie_2 \Sigma_1 \cdot e : \tau \) and \( FL(H' \cdot e) \cap dom(H) = \emptyset \), then \( \Gamma \bowtie_1, \bowtie_2 \Sigma(e) : \tau \).

**Proof.** By induction on the derivation of \( \Gamma \bowtie_1, \bowtie_2 \Sigma(e) : \tau \).

Lemma 11 states that typing of an expression is preserved by performing renamings on both the expression and signature.

**Lemma 11.** If \( \Gamma \bowtie e : \tau \), then \( \Gamma '\bowtie e' \bowtie \Sigma' \bowtie e : \tau \).

**Proof.** By induction on the derivation of \( \Gamma \bowtie e : \tau \).

We now restate the preservation and correctness theorems and prove them for the full language, including garbage collection.

**Theorem 2** (Preservation). If \( \Gamma \bowtie_2, \bowtie_2 e : \tau \) and \( \cdot \bowtie_1 \bowtie_2 \cdot, \Sigma_1 \bowtie P : \Sigma_2 \bowtie H ; e \rightarrow_\rho e' ; H' \bowtie \Sigma_1' \bowtie e \), then \( \cdot \bowtie_1 \bowtie_2 \cdot, \Sigma_1 \bowtie P : \Sigma_2 \bowtie H ; e \rightarrow_\rho e' ; H' \bowtie \Sigma_1' \bowtie e \).

**Proof.** We consider the cases for garbage collection \( T \longrightarrow_{gc} T' \) where, without loss of generality, \( e = \bullet T_1, T_2 \bullet \) and \( T_1 \times T_2 \).

- HGC-SavesGC. Then \( T = H \cdot e_1 \). By typing inversion, \( H : \Sigma \) and \( \Gamma \bowtie_2, \bowtie_2 e_1 : \tau_1 \). We have that \( (H) \bowtie H = \emptyset = \Sigma \), so by task typing, \( \Gamma \bowtie_2, \bowtie_2 (H ; FL(e_1) ; \emptyset, e_1) : \tau_1 \). By definition, \( FL(e_1) = FL(\emptyset, e_1) \).

- HGC-Cory. Then \( T = (H(\ell \mapsto v), S \bowtie \{\ell\}, H; F \cdot e_1) \). By typing inversion, \( F(H(\ell \mapsto v)) \bowtie H = \emptyset = \Sigma \) and \( \Gamma \bowtie_2, \bowtie_2 e_1 : \tau_1 \). We have \( F(H(\ell \mapsto v)) \bowtie H') = \emptyset = \emptyset = \emptyset \) and \( (H; F(e_1)) = \emptyset \). Since \( FL(e_1) \bowtie \Sigma = \Sigma_1, \Sigma_2 \bowtie e_1 : \tau_1 \), Lemma 11 gives \( \Gamma \bowtie_2, \bowtie_2 (H ; FL(e_1) ; \emptyset, e_1) : \tau_1 \). It is a straightforward exercise to show that \( S_1 = FL(H(\ell \mapsto v)) \).

- HGC-EnvGC. Then \( T = (H ; S, H ; F \cdot e_1) \). By typing inversion, \( F(H) \bowtie H = \emptyset = \Sigma \) and \( \Gamma \bowtie_2, \bowtie_2 e_1 : \tau_1 \) and \( S = FL(H_1) \bowtie e_1 \). By the task typing rules, we have \( \Gamma \bowtie_2, \bowtie_2 H_1 ; e_1 : \tau_1 \).

**Theorem 3** (Progress). If \( \cdot, \bowtie_2 e : \tau \) and \( H \bowtie \Sigma_1 \bowtie P : \Sigma_2 \bowtie e, \cdot \bowtie_2 \cdot, \Sigma_1 \bowtie e \), then \( e \) is a location or there exist \( e' \) and \( H' \) such that \( H ; e \rightarrow_\rho e' ; H' \).

**Proof.** We consider the case where \( e = \bullet T_1, T_2 \bullet \), where either \( T_1 = (H_1 ; S, H_1 ; F \cdot e_1) \) or \( T_2 = (H_1 ; S, H_2 ; F \cdot e_2) \) or both. Without loss of generality, suppose \( T_1 = (H_1 ; S, H_1 ; F \cdot e_1) \). If \( dom(H_1) \cap S = \emptyset \), then apply HGC-EnvGC. If \( e \in dom(H_1) \cap S \), then apply HGC-Cory. In either case, \( T_1 \) steps with D-GCStep.

Together, the progress and preservation theorems imply the standard type safety property: a well-typed term will not become stuck. Formally, if \( e : \tau \), then, for some \( H \) and \( H \), we have \( \emptyset, e \rightarrow_\rho \emptyset, H \). In particular, evaluation of \( e \) will never dereference a freed memory location. This completes the proof of memory safety. This is only half of the correctness of the garbage collector; we have not yet shown that collection is meaning-preserving, i.e., that the behavior of a garbage-collected program is identical to an execution without collection. This property is straightforward to prove using the same flattening machinery that was introduced in the previous section. The proof simply requires extending the definitions and proofs related to flattening to account for garbage collection. We
first extend the definition of flattening to cover tasks which are in the process of collection:

\[
\|\ell\|_{H;F;\mathit{HFL};P} \leadsto \hat{\ell}_1 \quad \|\ell\|_{F';\mathit{HFL};\mathit{H};P} \leadsto \hat{\ell}_2
\]

\[
\mathit{v}(H_f;S;H;F) \cdot \ell_1, (H_f';S';H';F') \cdot \ell_2 \quad \leadsto \mathit{v}(\hat{\ell}_1, \hat{\ell}_2)
\]

The two other rules for parallel tuples in which one task is of the form \(H \cdot e\) and the other is of the form \((H_f';S;H';F') [e']\) are admitted but are defined in the natural way. We also require two more lemmas regarding flattening, which will be used in the extension of the simulation proof.

**Lemma 12.** If \(\mathit{dom}(H) \cap \mathit{dom}(F) = \emptyset\) and \(\mathit{dom}(H) \cap \mathit{dom}(F) \subseteq \mathit{dom}(H)\) and \(\|\ell\|_{H;F;\mathit{HFL};P} \leadsto \hat{\ell}\) then \(\|F(e)\|_{H;F;\mathit{HFL};P} \leadsto \hat{\ell}\).

**Proof.** By induction on the derivation of \(\|\ell\|_{H;F;\mathit{HFL};P} \leadsto \hat{\ell}\). The interesting case is the case for a location \(\ell\). □

**Lemma 13.** If \(\mathit{dom}(H_1) \cap \mathit{FL}(e) = \emptyset\) and \(\|\ell\|_{H_1;H_2;P} \leadsto \hat{\ell}\) then \(\|\ell\|_{H_1;H_2;P} \leadsto \hat{\ell}\).

**Proof.** By induction on the derivation of \(\|\ell\|_{H_1;H_2;P} \leadsto \hat{\ell}\). The interesting case is the case for a location \(\ell\). □

Lemma 14 is simply a restatement of Lemma 8, which showed that a step of the hierarchical semantics can be simulated by zero or one steps of the flattened semantics. The lemma now includes garbage collection steps.

**Lemma 14.** Suppose that \(H; e \rightarrow e'; H'\) and \(H :: P : \Sigma\) and \(F \mapsto e : \tau \|\ell\|_{H;F;\mathit{HFL};P} \leadsto \hat{\ell}\). Then either \(\|\ell'\|_{H';F;\mathit{HFL};P} \leadsto \hat{\ell}'\) or \(\|e'\|_{H;F;\mathit{HFL};P} \leadsto \hat{\ell}'\) and \(\hat{\ell} \rightarrow \hat{\ell}'\).

**Proof.**

The new cases are those that instantiate D-ParAS1 or D-ParAS2 with one of the GC rules. We show that if \(\mathit{T} \rightarrow_{\mathit{GC}} \mathit{T}'\), then for any \(T_2\), if \(\|\mathit{T}, T_2 \|_{H;F;\mathit{HFL};P} \leadsto \hat{\ell}\) then \(\|T', T_2 \|_{H;F;\mathit{HFL};P} \leadsto \hat{\ell}\).

- **HGC-Copy.** Then \(\mathit{T} = (H_f[\ell \mapsto v];S \cup (\ell_1;H;F) [e])\). Let \(H_f = F(H_f[\ell \mapsto v] \cup H)\) and \(H_f' = F(H_f[\ell \mapsto v])\) and \(H_f' \rightarrow_0 H_f\). We have \(\|\ell_1\|_{H_f;H;F;\mathit{HFL};P} \leadsto \hat{\ell}_1\) and wish to show \(\|\ell_1\|_{H_f;H;F;\mathit{HFL};P} \leadsto \hat{\ell}_1\).

This follows from Lemma 12 since

\[
\begin{align*}
\|\ell \mapsto v\|(H_f) &= (\ell \mapsto v')(F(H_f[\ell \mapsto v] \cup H), H) \\
&= (\ell \mapsto v')(F(H_f) \cup H, F(v')) \\
&= F(H_f') \cup H \rightarrow_0 F(v') \\
&= H',
\end{align*}
\]

- **HGC-StarRG.** Then \(\mathit{T} = (H_1 \cdot e_1)\) and \(\|\ell_1\|_{H_1;H;F;\mathit{HFL};P} \leadsto \hat{\ell}_1\). If \(F = \emptyset\), then \(F(H_1) \cup H = H_1, \therefore \|\mathit{(H_1, F(e_1), \emptyset, e_1, T_2)}\|_{H;F;\mathit{HFL};P} \leadsto \hat{\ell}\).

- **HGC-EmptyGC.** Then \(\mathit{T} = (H_f;S;H;F) [e]\) and \(\|\ell_1\|_{H_f;H;F;\mathit{HFL};P} \leadsto \hat{\ell}_1\).

By typing, \(\|\ell_1\|_{H_f;H;F;\mathit{HFL};P} \leadsto \hat{\ell}_1\). By Lemma 13 gives \(\|\ell_1\|_{H_1;H;F;\mathit{HFL};P} \leadsto \hat{\ell}_1\) and \(\|H_f, e_1, T_2\|_{H;F;\mathit{HFL};P} \leadsto \hat{\ell}\).

Finally, Theorem 4 states the correctness of HGC: the hierarchical semantics produces a value from an expression if and only if the flattened semantics produces a value from the same expression, and the two values agree up to flattening. It was earlier shown that the flattened semantics produces a consistent with the non-collected hierarchical semantics, so this is sufficient to show that garbage collection does not change the meaning of a hierarchical program.

**Theorem 4.** Let \(e\) and \(\tau\) be such that \(\cdot \cdot \cdot e : \tau\).

- If \(e \rightarrow \hat{\ell}, \text{ then there exist } \ell \text{ and } H \text{ such that } e \rightarrow \hat{\ell}\) and \(\|\ell\|_{H;F} \leadsto \hat{\ell}\).
- If \(\hat{\ell} \rightarrow \hat{\ell}'\) and \(\|\ell\|_{H;F} \leadsto \hat{\ell}\), then there exists \(\hat{\ell}'\) such that \(e \rightarrow \hat{\ell}'\) and \(\|\ell\|_{H;F} \leadsto \hat{\ell}'\).

**Proof.** These are simply items 1 and 2 of Corollary 1, and are instances of Theorem 1, which is easily proven for the full language including garbage collection by using Lemma 14 instead of Lemma 8 in the proof. □

4. The Design

There are three primary challenges to the design of a practical implementation of the hierarchical semantics. These challenges primarily stem from the fact that a practical design must handle concurrency to realize parallelism.

- **Scheduling:** the semantics models scheduling using a nondeterministic interleaving of threads; a practical design must use a specific scheduler.
- **Algorithms for the heap hierarchy:** a practical design must use efficient and correct data structures and algorithms to represent the heap hierarchy.
- **Hierarchical garbage collection:** a practical design must ensure correct and performant garbage collection.

4.1 Scheduling

For this design, we use a work-stealing scheduler as described by Blumofe et al. [15]. This scheduler assigns every processor a double-ended queue, or “deque”, containing tasks associated with that processor. As a computation spawns new tasks, they are pushed onto the bottom end of the processor’s deque. Once a task is completed, the processor fetches the next task by popping the task from the bottom end of the deque. If the processor finds the deque to be empty, the processor steals a task from a randomly chosen victim processor by popping from the top of the victim’s deque.

The semantics presented in Section 3.1 accounts for some of the scheduler actions, such as the creation of parallel tasks, by distinguishing between active and inactive parallel tuples. There is a subtle but a crucial difference, however, between the management of tasks in the semantics and in work stealing. The semantics allows a parallel pair to be activated at any time by creating a task for each component and by pairing each task with its own heap. In work stealing, however, task creation is a lazy and asynchronous operation: a task is created only when one of the components of a parallel tuple is stolen, as the other component continues to execute serially. This subtle difference creates a challenge for identifying the heap for a task.

For example, consider a processor \(Q\) executing a parallel tuple \(\langle e_1, e_2 \rangle\). To execute the tuple, the processor starts executing the left component \(e_1\). Sometime after \(e_1\) is reduced to \(e'_1, e_2\) may be stolen and converted to a task \(T_2\) so that it can be executed by the thief processor. Since the algorithm has not converted \(e_2\) to a task before \(e_2\) was stolen, we don’t have a heap for it, and thus don’t know the objects allocated during the execution of \(e_1\) thus far.

To solve this problem, we assign to each parallel pair of the form \(\langle e_1, e_2 \rangle\) a level that corresponds to its nesting depth. For example, if the expression \(e = \langle e_1, e_2 \rangle\) has level \(m\), then \(e_1\) and \(e_2\) have level \(m + 1\). We then use levels to identify the heap of a task by tracking the levels at which steals take place.

4.2 Hierarchical Heaps

In our design, the abstract heaps of the semantics are implemented using pages and chunks, described below. These smaller units of memory are collected in data structures which we call superheaps,
which implement the heap hierarchy. The structure of the pages, chunks and superheaps allow efficient allocation, lookup, and other required operations on heaps and the heap hierarchy.

Pages, Chunks, and Chunk Pool. To reduce contention at memory allocation, we partition memory into chunks, each of which is a contiguous block of memory. Each chunk is further divided into one or more contiguous pages. A page can coincide with a system-level page, typically 4KB, but can be set to any power of two bytes.

Processors allocate and free chunks from and to a shared chunk pool. Once a chunk is allocated, it is locally divided into individual objects without further communication with other processors. Prior work has proposed design techniques for such a data structure [21, 27, 38]. Typical designs maintain a number of free chunk lists, each of which may, for example, contain chunks of a particular size. A number of different allocation policies can be employed to maximize efficiency and minimize fragmentation of the global memory [10, 18].

Distinguishing between chunks and pages might seem redundant, but it is not. Our design uses pages to find the chunk and the heap to which a given object belongs to by storing this information in the metadata for each page. Since a page size is a fixed power of two, the start of the page containing a given object can be located by truncating the address of the object. Once the page is found, a simple look up suffices to find the relevant metadata.

Superheaps. In our design, we use superheaps to identify the heap of a task by using levels. A superheap is a collection of chunks, where each chunk is tagged with the level at which it is allocated. The computation starts with a root superheap. During the execution, each steal starts a superheap for the stolen task. Thus at any point in the execution, a processor works on one superheap. When a steal happens, we remember the level of the stolen task at the superheap of the victim processor by marking it as “activated.” A heap in the semantics corresponds to all chunks in one superheap belonging to an activated level and possibly a contiguous range of unactivated levels. Figure 13 illustrates the structure of a superheap, where chunks at each level are drawn together at the same “level.” The levels 5, 6, and 7 are activated, since the scheduler has stolen the corresponding right tasks at these levels. Level 8 is not yet activated since its corresponding right task has not yet been activated and is still on the processor’s deque. Levels 5 and 6 each correspond to a heap of the semantics, and levels 7 and 8 together correspond to the local heap of the currently running task. We store various pointers in each chunk, such as a pointer to its level, and a pointer to the next chunk in its level to aid in the implementation of various operations.

Heaps operations To realize the semantics, the implementation of heaps needs to support four key operations: creating heaps, merging two heaps, allocating a new location in a heap, and looking up the value of a location in a heap. The design allows creating a heap by marking the level for the stolen task in the victim processor as activated and creating a new superheap for the stolen task. Merging two heaps corresponds to simply unioning their chunks. The allocation operation is simply a bump allocation within the current chunk. When the chunk is full, a new one is fetched from the chunk pool and assigned a level. Looking up an object is simply a pointer dereference.

4.3 Garbage collection

The semantics allow for collection of any heap in the memory hierarchy. The heaps that are leaves of the heap hierarchy, however, are special in the sense that they can be collected without communicating with any other processor, since they are disentangled. In our design, we therefore distinguish between two forms of collection. Local collection performs garbage collection on a leaf in the heap hierarchy; non-local collection collects all of the heaps in a subtree of the hierarchy.

Local Collection. A processor starts a local collection by locking its deque, ensuring that no tasks can be stolen during collection. A steal of a task from the deque corresponds in the semantics to the activation of a parallel tuple contained in the current expression, so locking the expression in rule HGC-StartGC corresponds to locking the deque.

After locking its deque, the processor proceeds to run a standard algorithm such as Cheney (copying) collection [20] adapted for hierarchical heaps. Such a collection would scan the root set (the collecting processor’s stack) for all pointers within the scope of collection. A pointer is considered to be within scope if it points to an object residing in one of the levels being collected. All roots in scope of the collection are copied to a new collection of chunks. The in-scope check also appears in the semantics, since rule HGC-Copy only copies locations that appear in both the scan set and the heap being collected.

After all roots are copied, copy collection continues normally, level-by-level in descending order. Any pointers to objects in scope of the collection but in lower levels are copied and scanned once collection gets to that level. Note that by the disentanglement property, no pointers from lower levels to higher levels can exist, so once a level is completely scanned, it does not need to be revisited. Once copy collection completes, all chunks in the levels being collected are released to the chunk pool, the processor’s deque is unlocked and execution resumes (corresponding to transition rule HGC-EndGC).

Non-Local Collection. As specified in Section 3.3, the GC semantics allows non-deterministic collection of any heap in the hierarchy. Here, we describe a specific non-local GC algorithm that collects all heaps in a given subtree at each non-local collection. Other algorithms are also possible.

A non-local collection starts by first choosing a subtree of the heap hierarchy to collect and identifying the processors participating in the collection. All participating processors then synchronize; each locks its deque and starts collection. All other, non-participating, processors continue to work on their execution as their accessible data will not be affected by the collection.

Participating processors then, in parallel, scan their root sets for all pointers within the scope of collection, just like for local collection. These pointers are all copied to the corresponding to-space. After all the roots are copied, collection proceeds level-by-level in parallel across all superheaps involved in the collection. However, each superheap is collected sequentially on its own.
Therefore, once collection on all of a parent superheap’s children is completed, one processor will continue collection on the parent superheap while the other processors remain idle. In this way, collection proceeds up the hierarchy of heaps until complete, at which point all the processors involved in the collection unlock theirdequeus and resume execution.

**Open issues.** Our design leaves open several important questions: 1) when to garbage collect, 2) whether to perform a local or non-local collection and 3) if non-local, which heaps, or subheap, to collect. The answers to these questions are not as straightforward as in traditional collection algorithms, where relatively simple amortization arguments can guide them. We leave the investigation of these questions to future work. In the implementation described below, we choose a simple technique adopted from traditional garbage collection literature.

5. **Implementation and Evaluation**

We describe a prototype implementation and preliminary evaluation of the proposed hierarchical memory manager.

5.1 **Implementation**

We implement our hierarchical memory manager on top of Daniel Spoonhower’s parallel extension of the MLton compiler [47]. As a whole-program optimizing, high-performance compiler for Standard ML, MLton [39] offers an excellent starting point for parallel computing. Spoonhower extended MLton to support a number of parallelism primitives including nested (fork-join) parallelism and futures. Spoonhower implemented many different schedulers; we use his implementation of a standard work-stealing scheduler.

Our implementation, which we call mlton-parmem, closely follows the design described in Section 4 but of course has to consider many more details. One such detail is the treatment of stacks. MLton executes all code with a traditional call stack, which is allocated on the heap. We allocate computation threads’ initial stacks in their associated heap. When the stack runs out of space, a new stack is allocated in the same heap at the current level of the computation. Upon first inspection, this model seems to blatantly violate disentanglement as a stack allocated in a lower level may have pointers in a higher level of computation. However, since the stack cannot be pointed to from any object, there is no issue of having parallel reads from another processor. In addition, since the stack is explicitly scanned as a root set for collection, we keep track of all the entangled pointers at collection time.

Currently, our implementation only supports local collection, which collects unactivated levels in descending order until the first activated level. For example, in Figure 13, local collection will only collect level 8 as all lower levels (drawn higher in the figure) have been activated. Our algorithm for deciding when to collect is a straightforward adaptation of the corresponding algorithm from traditional GC literature. We initiate a local collection when the local heap becomes half full and we resize the heap to a small multiple (8 in our experiments) of the size of the live set.

5.2 **Benchmarks**

We evaluate performance under five benchmarks. Each benchmark is written in a purely functional style and in three different dialects of ML: Standard ML, our parallel extensions to it, dubbed “mlton-parmem”, and Manticore.

All of our benchmarks use sequences as the core data structure. Sequences are implemented as weight-balanced trees with data elements at the leaves. This data structure allows for naturally expressing parallel computation over the sequences. The benchmarks use a manual approach to granularity control: any potentially parallel operation first checks if the length of the input sequence is less than or equal to a specified grain size. If so, the operation is carried out sequentially without creating parallel tasks. If the length of the input sequence is greater than the grain, the operation will create the parallel tasks. The grain size can be supplied at run-time.

The MemStress benchmark is a synthetic benchmark designed to heavily stress the memory manager by allocating a large amount of data and releasing it almost immediately. This is achieved by building, in parallel, an integer sequence (using the `tabulate` operation on sequences) where, at each element, a large list is sequentially allocated. The list’s tabulation function is simply the identity function and the sequence’s element is then either the first or second element of the list depending on whether the element index is even or odd. Note that the amount of computation done per allocation is very small, ensuring that the bulk of the work of the program lies in the allocation and collection of the data. The benchmark tabulates a 100,000 element sequence with granularity 100 and allocates a 10,000 element list per sequence element.

The Tabulate benchmark is designed to evaluate the performance of an embarrassingly parallel sequence tabulation operation. In this benchmark, a sequence is tabulated and every element is set to a Fibonacci number, calculated each time. In our benchmark, we tabulate a 100,000 element sequence with granularity 100. For each element, the 20th Fibonacci number is calculated sequentially using the classical recursive algorithm and stored.

The Raytracer benchmark is adapted from the raytracer benchmark written for the Manticore language [8]. It renders a 512px × 512px scene in parallel. The original program was written in ID [42] and does not use any special data structures to improve performance.

The SMVM benchmark performs a sparse matrix vector multiplication. The sparse matrix is represented in a typical fashion as a sequence of rows, where each row only contains the non-zero entries as an index-value pair. Multiplication is then done in parallel over the rows. Our benchmark performed the test on a 10,000 × 10,000 matrix with granularity 100.

The DMM benchmark performs a dense matrix multiplication. The matrices are represented as a sequence of sequences. Multiplication is then done in parallel over the rows of the first matrix. Our benchmark performed the test on 500 × 500 matrices at the granularity of individual dot-products.

5.3 **Measurements**

For our experiments, we use a 64-core AMD machine with 128 gigabytes of memory, but due to technical limitations of the underlying MLton codebase, we are not able to utilize more than 32 cores. We use interleaved NUMA allocation in all of our experiments. For the grain sizes we use, our benchmarks create anywhere from several thousand to hundreds of thousands of tasks.

In addition to our mlton-parmem compiler, we run our benchmarks on the manticore compiler, the mlton compiler for Standard ML and Spoonhower’s parallel extension of MLton, which we call mlton-spoonhower. The memory manager of mlton-spoonhower allows for parallel allocation, but performs stop-the-world sequential collection. We choose this compiler as a naive baseline for comparing parallel garbage collectors.

The Manticore compiler for Parallel ML [28] offers support for an ML-like language extended with several paradigms for parallelism, including fork-join parallelism and speculation. Manticore uses an advanced NUMA-aware memory manager based on the Doligez-Leroy-Gonthier and Appel semi-generational collectors. It is a reasonably well-developed system with particular innovations in memory management and has been shown to deliver good performance [8, 11, 13, 28, 44]. Since it is based on ML, Manticore is a good basis for comparison for the approach suggested in this paper.

We use the mlton compiler timings as the sequential baseline for all our speedup calculations. We derive the sequential versions of our
benchmarks by replacing the par calls by sequential composition. In performance evaluations, it is sometimes preferable to use highly optimized serial code as a baseline. Since, however, we wish to study the performance and the scalability of our memory manager, doing so would obscure our results.

We run each benchmark five times and report the median of the results in the tables and graphs. We measure two quantities across these benchmarks. The first quantity is the single-processor slowdown, or overhead, compared to a sequential run. As the name suggests, the overhead indicates the overhead incurred by the parallel runtime and memory manager and helps us establish the practical viability of the compiler for parallel computations against a sequential compiler. The second quantity is the speedup of the mlton-parmem, mlton-parmem and Manticore compilers against the mlton baseline.

5.4 Performance and Scalability

Figure 14 shows some of the key measurements across our benchmarks using the following quantities.

- Serial run time, $T_s$, is the time in seconds for a serial run with mlton.
- Uniprocessor run time, $T_1$, is the uniprocessor run-time of the parallel code reported in seconds;
- The overhead $O$ is the overhead calculated as $T_1/T_s$.
- 32-processor run time, $T_{32}$, is the 32-processor run-time of the parallel code reported in seconds;
- The speedup $S$ is the 32-core speedup calculated as $T_s/T_{32}$.

Across all the benchmarks, we notice that the overhead compared to mlton is comparable for both mlton-spoonhower and mlton-parmem and range between $-15\%$ and $51\%$. These measurements suggest that there is some overhead to parallelism but using a hierarchical memory manager is not significantly more expensive than the naïve parallel memory manager employed in mlton-spoonhower. In fact, overheads of mlton-parmem are slightly better in 3 benchmarks than those of mlton-spoonhower. The manticore compiler incurs significantly larger overheads compared to mlton. This is consistent with earlier observations with the Manticore compiler [12], which focuses on performant parallel execution but not fast sequential execution. Our observed overhead of between $1.19$ and $2.95$ is in line with expectations [12].

In terms of scalability, which we can observe by comparing the Speedup “$S$” columns across compilers, our mlton-parmem compiler delivers the best speedups for all benchmarks except Raytracer. Figure 15 shows the speedup curves for individual benchmarks. In all benchmarks, we observe that mlton-parmem scales nicely up to 32 processors. While mlton-spoonhower also scales reasonably in some benchmarks, it consistently fails to scale as well as mlton-parmem. Indeed, in both SMM and DMM, we see the speedup of mlton-spoonhower starting to plateau towards 32 processors. In MemStress, mlton-spoonhower fails to achieve much more than $3\times$ speedup regardless of the number of processors it utilizes. The manticore compiler scales very well, but also trails mlton-parmem, except in Raytracer. As programs compiled with manticore typically suffer from large overhead over the baseline, mlton, they require more processors in order to overcome the
overhead. Indeed, manticore requires about 3 processors just to match the performance of mlton on the MemStress benchmark. In Raytracer, the overhead of manticore against mlton poses less of a problem and allows manticore to exhibit high scalability.

These results suggest that the proposed parallel hierarchical memory manager can be implemented in the context of a high-performance compiler such as MLton and can deliver good performance. The overheads and the speedups that our mlton-parmem compiler delivers suggest that the basic idea behind the approach of coupling parallel scheduling and memory management is sound. We note, however, that our comparative study does not carefully account for many important factors that can have significant effects on performance including various internal “hardwired” constants such as block and page sizes and amortization constants that vary between compilers. We leave a more careful implementation and a more detailed experimental study for future work.

6. Related Work

There has been significant work on parallel garbage collection and on scheduling for parallel programs, which we discuss below.

Garbage Collection. The work on parallel garbage collection dates back at least to the multilisp [31] parallel collector. The multilisp collector kept separate heaps for each processor and had no shared heap. A variety of parallel collectors use a shared heap, sometimes with a local block-allocated pool for allocation [6, 9, 21, 26, 32]. This requires significant synchronization among the processors, and does not give the locality advantages of a generational collector.

The Doligez-Leroy-Gonthier (DLG) collector [23, 24] introduced the idea of having local nurseries (heaps) for each processor in addition to a global heap. Memory is allocated in the local heaps and promoted to the global heap on a local collection, or in various other situations, such as when writing into state that is in the global heap. The approach has a similar advantage as a single-level generational sequential collector—by the time a local heap is promoted to the global heap, much of what has been allocated is already garbage. Also if the local nursery fits in cache, its locations can be kept “warm” in cache so that allocation is cheaper. Finally, the local collections can be done independently. Several other parallel collectors have refined the idea of local nurseries and a global heap [5, 8, 25, 37, 51]. These collectors differ in precisely how the two levels are divided and what collection approach is used at each level. For example, the original DLG collector allows the global heap to be collected concurrently using Dijkstra’s concurrent mark-sweep algorithm [22]. The ABFR collector [8], on the other hand, uses a form of stop-the-world copying collection on the global heap.

To handle mutable data, these two-level collectors ensure a property similar to disentanglement by requiring that most or all (depending on the strategy used by the individual collector) mutable data be stored in the global heap. Thus, mutation-heavy programs will require a great deal of global collection. In all of these collectors, the decision to collect is based purely on when memory is full, or perhaps when mutations happen across boundaries, but is not related to scheduling decisions or other aspects of the tasks themselves. In fact, they are agnostic to the tasks running on the processors.

Marlow et al. use a generational block-structured parallel collector [38] with support for multiple generations. However, the hierarchy does not form a tree, but rather is a sequence of global heaps of varying size.

Pizlo et al. describe a hierarchical heap structure [43], which has some similarities to ours. It is, however, designed for realtime collectors, and in particular as a replacement for the Real-time Specification for Java (RTSJ). The key difference is that the hierarchical structure is meant to be (mostly) static and user-defined, while ours is meant to be highly dynamic and invisible to the user. The idea in their work is that real-time threads can use heaplets at the leaf, and that each heaplet can support its own style of GC.

Bocchino et al. suggest using regions for parallel memory allocation [16]. Regions can be allocated in a tree hierarchy dynamically based on the structure of the parallelism. Their motivation, however, is very different. In particular they use typing rules on the regions to statically prevent any race conditions. They have no discussion of garbage collection within or across regions.

Morrisett et al. used an operational semantics for abstracting garbage collection in a way similar to our operational semantics [40]. However, it was in a purely sequential context, and is more abstract than ours, e.g. not even capturing the distinction between copying and mark-sweep collection.

The results presented here build on our recent paper [4], which proposed coupling computation and memory management.

Scheduling. Nearly all modern parallel programming languages rely on a run-time scheduler to map tasks to processors. Determining an optimal schedule is actually NP-hard [50], but a classic result by Brent [17] shows that a 2-factor approximation can be computed by using a greedy-scheduling principle. Brent’s theorem, however, does not take into account the cost of the scheduling algorithm itself. Over the years, many efficient scheduling algorithms implementing Brent’s greedy principle have been developed and bounds have been proven on runtime and space use [3, 15, 29, 30, 41], and locality [1, 14]. These schedulers perform well in practice [3, 29, 48] when the granularity of parallel tasks can be controlled so as to prevent the creation of tiny tasks [2, 12, 49] Much of this work on schedulers, however, does not try to couple the structure of memory with that of the schedule as we do here.

7. Conclusion

This paper presents a memory management technique for parallel functional programs that is based on the general strategy of closely coupling the structure of the memory with that of the computation. As we show, such close coupling enables using the invariants of the semantics of the language, such as the disentanglement property, in managing memory.

There are many interesting directions for future research. Here, we consider fork-join programs, which include a relatively large class of computations. Extending the technique to include other forms of parallelism such as async-finish and futures may be interesting and desirable. Another natural direction for future research would be to extend the techniques to allow some mutable state. We present an approach to realizing the semantics in practice by considering a work-stealing scheduler, by describing the data structures for implementing hierarchical heaps, and by describing an algorithm for performing garbage collection. The initial experimental results for our extension of MLton suggest that the proposed techniques can be implemented efficiently. The implementation, however, is incomplete—it employs local collection only—and is not carefully optimized. A natural next step would be to finish a complete and optimized implementation that can perform both local and nonlocal collections. Finally, it would be interesting to investigate if the techniques described here can be applicable to other parallel languages and systems [19, 28, 29, 33–36, 46].

Acknowledgments

This research is partially supported by the National Science Foundation under grant numbers CCF-1320563 and CCF-1408940, by the European Research Council under grant number ERC-2012-StG-308246, and by Microsoft Research. We also thank the reviewers for their insightful comments and in particular the idea of using flattening to prove correctness.


