

Correctness-Preserving Derivation of Concurrent Garbage Collection Algorithms

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Abstract

Constructing correct concurrent garbage collection algorithms is notoriously hard. Numerous such algorithms have been proposed, implemented, and deployed – and yet the relationship among them in terms of speed and precision is poorly understood, and the validation of one algorithm does not carry over to others.

As programs with low latency requirements written in garbage-collected languages become part of society’s mission-critical infrastructure, it is imperative that we raise the level of confidence in the correctness of the underlying system, and that we understand the trade-offs inherent in our algorithmic choice.

In this paper we present correctness-preserving transformations that can be applied to an initial abstract concurrent garbage collection algorithm which is simpler, more precise, and easier to prove correct than algorithms used in practice — but also more expensive and with less concurrency. We then show how both pre-existing and new algorithms can be synthesized from the abstract algorithm by a series of our transformations. We relate the algorithms formally using a new definition of precision, and informally with respect to overhead and concurrency.

This provides many insights about the nature of concurrent collection, allows the direct synthesis of new and useful algorithms, reduces the burden of proof to a single simple algorithm, and lays the groundwork for the automated synthesis of correct concurrent collectors.

1. Introduction

As garbage-collected languages like Java and C# become more and more widely used, the long pauses introduced by traditional synchronous (“stop the world”) collection are unacceptable in many domains. This is true both at the high end, where the collection of multi-gigabyte heaps causes very long pauses, and at the low end, where systems are used for real-time, embedded, and sensor applications requiring very low latency. As a result, concurrent collectors are now available in most major production virtual machines.

However, concurrent collectors are extremely complex and error-prone. Since such collectors now form part of the trusted computing base of a large portion of the world’s mission-critical software infrastructure, such unreliability is unacceptable.

The study of concurrent collectors began with Steele [40], Dijkstra [21], and Lamport [31].

Concurrent collectors were quickly recognized as paradigmatic examples of the difficulty of constructing correct concurrent algorithms: Steele’s algorithm contained an error which he subsequently corrected [41], and Dijkstra’s algorithm contained an error discovered and corrected by Stenning and Woodger [21]. Doligez and Leroy developed a multiprocessor collector for ML [23] which was subsequently found to contain an error [22]. Furthermore,

some correct algorithms [9] had informal proofs that were found to contain errors [36].

Much later, Yuasa [45] introduced the snapshot-based algorithm, which is conceptually simpler and trades earlier termination and increased concurrency for reduced precision.

Many additional incremental and concurrent algorithms have been introduced over the last 30 years [30, 1, 2, 3, 4, 5, 6, 11, 13, 16, 17, 24, 26, 27, 29, 32, 34, 35, 39, 44], but there has been very little experimental comparison of the algorithms and no formal study of their relative merits. While there is now a well-established “bag of tricks” for concurrent collectors, each algorithm composes them differently based on the intuition and experience of the designer. However, because of the complex interactions of the invariants required by the different “tricks,” many potential combinations of techniques are not used, leading to an underexplored design space. Furthermore, since each algorithm is different, a correctness proof for one algorithm cannot be re-used for others.

All concurrent collectors must decide how to answer the following basic questions:

- Where is the collector in its progress through the heap?
- Which objects must be traced to guarantee that all live objects will be found?
- How does the collector terminate in spite of allocation?
- Which interleavings are allowed between the mutator and the collector?

Our long-term research agenda is to generate a provably correct concurrent garbage collector from a simple, declarative specification of the techniques that should be applied to meet the particular needs of the target system with respect to latency, throughput, and space consumption.

In previous work [43] we hypothesized that the way in which the above questions are answered could be expressed as transformations of a single base algorithm, informally described some such transformations, and evaluated their relative performance experimentally.

In this work we substantiate that hypothesis: we present a single algorithm (which we call the *Apex* algorithm) and a set of composable transformations corresponding to each of the above questions. Each transformation can be applied to arbitrary subsets of the objects, or to restricted subsets for which we present precise formulations.

Transformations can be applied at the granularity of a single object in the heap. This simplifies the proof methodology since we can reason inductively about the transformation, but also allows enormous flexibility since different transformations can reliably be applied to different objects depending on their characteristics.

Furthermore, we formalize for the first time the notion of the *relative precision* of concurrent collectors, and express the transformations as correctness-preserving and precision-reducing. We also

discuss informally how the reduction in precision provides useful tradeoffs in terms of implementation cost, speed of convergence, and level of concurrency.

Our transformational approach yields a wide range of algorithms. We show derivations for some well-known existing algorithms, and also derive some new algorithms which we expect will have desirable properties in practice. In particular, our generalization and formalization of the tradeoff between incremental-update and snapshot-at-the-beginning approaches allows a novel approach to newly allocated objects which yields high precision combined with rapid termination.

The contributions of this paper are:

- A formal framework for describing concurrent garbage collection algorithms;
- The simple “Apex” algorithm from which all others are derived;
- a set of transformations that can be applied to it to yield an enormous number of potential algorithms with different precision, concurrency, and efficiency properties;
- A formal definition of what constitutes relative precision of concurrent collectors;
- a proof of correctness for the transformations, which are shown to be precision-reducing (while improving other aspects of the algorithm);
- The application of the methodology to yield incremental update collectors in the styles of Dijkstra and Steele, snapshot collectors like that of Yuasa, as well as previously unknown algorithms with high precision, rapid termination, and high concurrency.

This work is presented in the context of a mark-and-sweep style of collector. While we show the synchronization with the sweep phase, we do not consider the details of its implementation, which contains a number of its own complexities. We have also simplified the design space by using a write barrier which is always atomic. This corresponds to some but not all implementations used in practice (for instance on a uni-processor safe point based virtual machine, or for write barriers whose synchronization can be implemented with a single compare-and-swap). So in general some manual transformation may still be required to achieve desired performance in the resulting algorithms.

A key aspect of this work is the *modularization* of the proof obligations. Although we have not proved the correctness of the Apex collector, we have proved the correctness of a broad variety of transformations needed for the creation of an efficient algorithm. This breaks the requirements for the creation of a correct algorithm and implementation into small modular proof components which can be re-used across an enormous range of algorithms, rather than requiring a monolithic proof of each new algorithm.

2. A Parametric Concurrent Collector

In this section we present a simple formulation of concurrent collectors using a common algorithm, in which the differences between algorithms are encapsulated in a function (called *expose*) which chooses objects to use as the starting points for additional transitive closure operations which find and mark live data that would otherwise have been hidden by concurrent mutation.

The *expose* function makes these choices based on a log of mutator and collector operations, which is the formal analogue to the information captured by the write barrier in real-world implementations.

2.1 A Trace Model for Concurrent Collectors

We model the heap as an infinite set \mathcal{U} of objects. Each object *obj* contains a set $Fields = \{f_1, \dots, f_F\}$ of fields. A field f_i is a

function $f_i : \mathcal{U} \rightarrow \mathcal{U} \cup \{\mathbf{null}\}$ mapping an object to another object, or to a designated value **null**. Without loss of generality we assume that all objects have the same set of fields F . For convenience, we use $obj.f_i$ to denote the value $f_i(obj)$.

All reachable objects are reachable from a finite set of R root objects, denoted $root_1, \dots, root_R$.

For the time being we assume that stack frames are heap-allocated – some systems, especially for functional languages are in fact implemented in this way.

A global state of the program consists of: (i) the heap; (ii) the state of the mutator; (iii) the state of the collector. We model the mutator as a sequence of allocations and mutations of pointers over a given heap. The state of the mutator is its position in the sequence of allocations and mutations. The state of the collector consists of an assignment of values to all its variables.

A program trace is a potentially infinite sequence of program global states, where a *sequence* is defined in a standard way as a map from natural numbers to an alphabet Σ . Given a sequence $S = S_0, S_1, \dots$, we define its finite prefix P of length $|P| = k$, to be the first k letters in the sequence S_0, S_1, \dots, S_{k-1} , and denote it by $\text{pre}(S, k)$.

Given a finite sequence prefix P and a sequence S , we denote the concatenation of P and S by $P \bullet S$. Similarly, given a finite sequence prefix P and a letter $\tau \in \Sigma$, we denote by $P \bullet \tau$ the concatenation of P and τ .

Our algorithm uses an *interaction log* to record information about the combined behavior of the collector and the mutator. This log is used by the collection algorithm to select the objects to be marked. A log of the interleaving of mutator/collector operations is natural for a concurrent collector because it closely matches the use of write barriers in practical implementations: the function of the write barrier is to synchronize with the collector, which in some collectors is in fact done with a log of heap writes.

The interaction log is a sequence of log entries of the following kinds: (i) a *tracing* entry recording a tracing action of the collector as it traverses the heap during the marking phase; (ii) a *mutation* entry recording a pointer redirection action by the mutator; (iii) an *allocation* entry recording an allocation of a new object by the mutator. This is formally defined as follows:

DEFINITION 2.1. A log entry is a tuple $\langle \mathbf{k}, source, fld, old, new \rangle \in \{\mathbf{T}, \mathbf{M}, \mathbf{A}\} \times \mathcal{U} \times Fields \times (\mathcal{U} \cup \{\mathbf{null}\}) \times (\mathcal{U} \cup \{\mathbf{null}\})$ where:

- k identifies the kind of action as one of tracing, mutation, or allocation, denoted by \mathbf{T}, \mathbf{M} , and \mathbf{A} , respectively.
- *source* is the object affected by the action.
- *fld* is the field of source affected by the action.
- *old* is the value of the field *source.fld* prior to the action.
- *new* is the value of *source.fld* subsequent to the action.

Tracing actions do not change the structure of the heap; therefore $old = new$ for all tracing entries. Allocation actions allocate the object *new*, which must not appear previously in the trace.

For convenience, we define selectors for log entry tuples. Given a tuple $\tau = \langle \mathbf{K}, s, f, o, n \rangle$, we define $\tau.kind = \mathbf{K}$, $\tau.source = s$, $\tau.field = f$, $\tau.old = o$, and $\tau.new = n$.

2.2 The Parametric Algorithm

Fig. 1 presents the pseudo-code for a parametric concurrent mark-and-sweep collector. The operation of this collector is defined over a prefix of the interaction log, recording the collector and mutator interaction. Recording mutator actions in the log is performed by the mutator’s write and allocation barriers, as shown in Fig. 2.

Before describing the algorithm in more detail, we first describe the assumptions we have made for clarity of presentation and the assumptions under which the algorithm operates:

```

collect() {
  atomic
  marked ← {root1, ..., rootR}
  pending ← pending ∪ (∪x∈marked fields(x))

  do {
    mark()
    addOrigins()
  } while (?)

  atomic
  addOrigins()
  mark()

  sweep()
}

mark() {
  while (pending ≠ ∅) {
    (obj, fld) ← removeElement(pending)
    atomic
    dst ← obj.fld
    log ← log • ⟨T, obj, fld, dst, dst⟩
    if (dst ≠ null ∧ dst ∉ marked){
      marked ← marked ∪ {dst}
      pending ← pending ∪ fields(dst)
    }
  }
}

addOrigins() {
  atomic
  origins ← expose(log) \ marked

  marked ← marked ∪ origins
  pending ← pending ∪ (∪x∈origins fields(x))
}

```

Figure 1. Parametric Mark-and-Sweep Collector.

```

mutate(source, fld, new) {
  atomic
  log ← log • ⟨M, source, fld, source.fld, new⟩
  source.fld ← new
}

mutateAlloc(source, fld) {
  atomic
  new ← allocate new object
  log ← log • ⟨A, source, fld, source.fld, new⟩
  source.fld ← new
}

```

Figure 2. Mutator write-barrier and allocation-barrier.

- we do not specify how the `sweep()` operation proceeds, except to ensure that there is the proper synchronization between the mark and sweep phases. We also do not consider compaction, which requires the dynamic relocation of objects. While these are both important issues, they are beyond the scope of this work.
- we assume that there is only a single execution of a collection cycle at any given point in time. That is, the **T** operations in the log all belong to a single collection cycle. Multiple (even

overlapped) collections can be performed by differentiating **T** entries accordingly.

The parametric collection algorithm does not explicitly define how objects are selected to be marked as live. This is defined to be a parameter of the collector. The algorithm, however, does restrict concurrency by making two decisions:

- assuming that write barriers are atomic with respect to collector operations. This assumption is inline with practical systems as mutators and the collector are only allowed to interleave at *safe points*, which do not include the write-barriers. Effectively, this means that a collector cannot preempt a mutator during a write barrier. Under this assumption, we can restrict attention to a system with a single mutator thread without loss of generality.
- the collector is able to trace object fields in an arbitrary order, that is, rather than scanning fields of each object in order.

The collection cycle of the algorithm is described in the `collect()` procedure. The collection cycle consists of two phases: (i) the *marking phase*, in which the collector marks potentially live objects; (ii) the *sweeping phase*, in which unmarked objects are reclaimed.

The collection cycle starts by atomically selecting the set of root objects as origins. This operation is executed atomically, and thus no concurrent mutations could be performed by the mutator.

After selecting the root objects as origins, the collector proceeds by repeatedly tracing heap objects and marking them (`mark()` procedure), and adding origins to be considered by the collector due to concurrent mutations performed by the mutator (`addOrigins()` procedure). These two steps are repeated until a non-deterministic choice (denoted by ‘?’ in the figure) triggers a move to an atomic phase in which the remaining origins and objects to be marked are processed atomically. This atomic phase guarantees the termination of the algorithm, and is in line with some practical collector implementations (e.g., [7]). Nevertheless, in Section 5.5, we show how to derive algorithms in which this atomic marking phase can be eliminated.

At the end of the marking phase, the *marked* set contains the objects that are marked as live, and the collection cycle proceeds to the sweep phase, reclaiming unmarked objects and completing the collection cycle.

2.3 Marking Traversal

The `mark()` procedure implements a collector traversal of the heap. In the algorithm, we use a pair (obj, fld) to denote the field *fld* of an object *obj*. We use $obj.fld$ to denote the object pointed to by the field *fld* of the object *obj*, and $fields(obj) = \{(obj, f_1), \dots, (obj, f_F)\}$ to denote the set of all object fields for a given object *obj*.

The procedure uses a set *pending* of pending fields to be traversed, and performs a transitive traversal of the heap by iteratively removing an object field from the set *pending* and tracing from it. Whenever an object field is traced-from, the procedure inserts a tracing entry into the log. When the traced object field points to an unmarked object, the object is marked, and its fields are added to the pending set.

During this traversal, the mutator might concurrently modify the heap. These concurrent mutations might cause reachable objects to be *hidden* from the traversal, and thus remain unmarked by the traversal.

2.3.1 The Collector Wavefront

All collectors discussed in this paper rely on cooperation between the collector and the mutator to guarantee correctness in the presence of concurrency. A key part of the cooperation is tracking the progress of the collector through the heap, since mutations can be treated differently depending on whether they happened in the por-

tion of the heap already scanned by the collector (behind the wavefront) or not yet scanned (ahead of the wavefront). The wavefront consists of the set of object fields (that is, *not* the values of the pointers in those fields) that have been traced by the collector thus far.

DEFINITION 2.2. *Given a log prefix P , the collector wavefront is the set of object-fields that have been traced by collector operations in P , that is:*

$$\mathcal{W}(P) = \{(P_i.source, P_i.field) \mid P_i.kind = \mathbf{T} \wedge 0 \leq i < |P|\}$$

Given a log prefix P , we say that an object field (o, f) is behind the wavefront when $(o, f) \in \mathcal{W}(P)$, and ahead of the wavefront when $(o, f) \notin \mathcal{W}(P)$.

Most practical collectors use conservative abstractions of the wavefront rather than the precise definition provided here. That is, the wavefront is tracked at an object granularity. However, the precise wavefront is not merely theoretical and has recently been used in the hardware-assisted collector for the Azul Java server, which has a “not marked-through” bit in every pointer [18].

EXAMPLE 2.3. Fig. 3 shows an example of a possible mutator and collector interleaving. In this figure, the progress of collector tracing through the heap is shown by the tracing actions T_1, \dots, T_6 , and by using a darker color for traced fields. The sequence of mutations is shown as M_1, \dots, M_7 . For brevity, we only present part of the states and show the effect of multiple operations in a single step.

The interaction log prefix P^e for this example is:

$$\begin{aligned} &\langle \mathbf{T}, r1, f2, A, A \rangle, \langle \mathbf{T}, A, f1, null, null \rangle, \langle \mathbf{T}, r1, f3, null, null \rangle, \\ &\langle \mathbf{M}, r1, f1, null, B \rangle, \langle \mathbf{M}, A, f1, null, B \rangle, \langle \mathbf{M}, r1, f3, null, E \rangle, \\ &\langle \mathbf{M}, A, f2, C, null \rangle, \langle \mathbf{M}, r1, f1, B, null \rangle, \langle \mathbf{T}, A, f2, null, null \rangle, \\ &\langle \mathbf{T}, r1, f1, null, null \rangle, \langle \mathbf{M}, A, f3, D, null \rangle, \langle \mathbf{M}, A, f1, B, null \rangle, \\ &\langle \mathbf{T}, A, f3, null, null \rangle \end{aligned}$$

The wavefront at the end of the shown prefix P^e is: $\mathcal{W}(P^e) = \{(r1, f2), (r1, f3), (A, f1), (A, f2), (r1, f1), (A, f3)\}$

2.4 Adding Origins

The `addOrigins()` procedure uses the log of mutator and collector interaction to select a set of additional objects to be considered as origins, marked as live and traced from. When this procedure is invoked by the collector, it is possible that a number of reachable pointers were hidden by the mutator behind the \mathcal{W} during the `mark()` procedure.

The `addOrigins()` procedure finds those hidden pointers and produces a set of objects to be used as additional origins. This set is a safe over-approximation of the reachable objects, which were hidden during the `mark()`. The core of `addOrigins()` is the function `expose` which takes a log prefix and returns a set of objects that should be considered as additional origins. Each object returned by `expose` is then marked, and its fields are inserted into the *pending* set.

All algorithms effectively differ in their `expose` function. In the next sections, we will present different choices of `expose` which correspond to various garbage-collection algorithms.

2.4.1 Mutator Barriers

Fig. 2 shows the write-barrier and allocation-barrier used by the mutator. The procedure `mutate(source, fld, new)` is called by the mutator to mutate a pointer in the heap. The procedure `mutateAlloc(source, fld)` is called by the mutator to allocate a new object and store it in the given object field. To collaborate with the collector, the mutator barriers record their actions in the interaction log.

When the mutator performs an assignment $source.fld \leftarrow new$ with $new \neq null$, we say that a pointer is *installed* from $(source, fld)$ to new . When the object field $(source, fld)$ is behind the wavefront, we say that the pointer is *installed behind the wavefront*. Similarly, we may say that the pointer is installed ahead of the wavefront.

Similarly, whenever we assign a value to a field $(source, fld)$ containing an existing pointer, we say that the existing pointer is *deleted*. If the field $(source, fld)$ is ahead (behind) of the wavefront, we say that the pointer is *deleted ahead (behind) of the wavefront*.

EXAMPLE 2.4. In Fig. 3 the mutation (M_2) results in a pointer from $(A, f1)$ to B installed behind the wavefront, and the mutation (M_6) results in the pointer from $(A, f3)$ to D being deleted ahead of the wavefront.

3. The Apex Algorithm

This section introduces the **Apex** algorithm, an instance of the parametric collector presented in Fig. 1. The Apex algorithm is the starting point for the derivation steps which follow in the rest of the paper.

The Apex algorithm makes use of a technique called *rescanning*. Rescanning is a technique which given a set of objects, identifies the object fields that were modified behind the collector wavefront. It then returns the pointers to the objects residing in those modified fields. This approach is necessary to *expose* reachable objects that are hidden by a sequence in which: (i) a pointer to an object is stored in a field behind the wavefront; (ii) all other paths to the object ahead of the wavefront are removed before the collector reaches them.

Rescanning provides a high degree of accuracy, since all hidden pointers are identified precisely.

$$\begin{aligned} expose^{apex}(P) = \{o.f \mid P_i.source = o \wedge P_i.field = f \wedge 0 \leq i < |P| \\ \wedge P_i.kind \in \{\mathbf{M}, \mathbf{A}\} \wedge (o, f) \in \mathcal{W}(\text{pre}(P, i))\} \end{aligned}$$

Given a log prefix P , $expose^{apex}(P)$ returns the current contents of all of the mutated fields behind wavefront.

Note that because the execution of `expose` is performed inside of an **atomic** block in the algorithm in Figure 2, the rescanning of *all of the fields of all of the objects* that were modified are scanned atomically. This means that a pure rescanning algorithm has very low concurrency.

EXAMPLE 3.1. Consider the example interleaving of Fig. 3. The function $expose^{apex}(P^e)$ atomically performs rescanning of the fields $(A, f1)$ and $(r1, f3)$. This results with $expose^{apex}(P^e) = \{E\}$. Unless returned by $expose^{apex}$, object E would have been lost. Assuming there are no further mutations, the objects marked by the Apex algorithm in this collection cycle will be: $\{r1, A, E\}$. In future sections, we see that other (less precise) collectors will consider additional objects as live.

Conceptually, Apex is very similar to the Steele algorithm [40] algorithm, but with an accurate \mathcal{W} definition.

The Apex algorithm is used as the base algorithm in our framework from which all other algorithms are derived using correctness-preserving transformations. Fig. 4 shows part of the derived algorithms, ordered by relative precision.

In this paper, our focus is on unifying the various collection algorithms, and relating them in terms of precision. We show that our transformations are correctness-preserving, but assume that the Apex algorithm is a correct starting point.

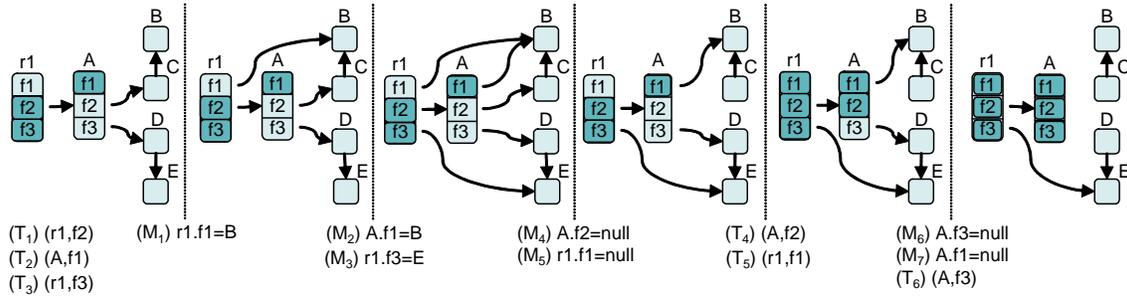


Figure 3. Example interleaving of mutations and tracing operations.

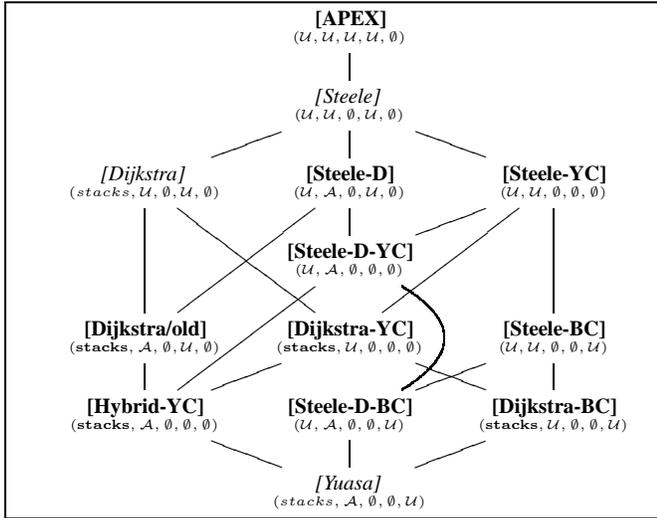


Figure 4. Relative precision of part of the existing and newly derived algorithms. New algorithms are shown in boldface. The most precise algorithm is shown at the top. For readability, only derived algorithms with abstracted wavefront are shown. Tuples are of the form (SR, IS, FL, WC, BC) .

4. Precision of Collection Algorithms

In this section, we introduce the notion of relative precision of concurrent collection algorithms. This allows us to formally relate the various algorithms which are instantiated in our parametric framework.

A correct marking algorithm must satisfy both a *safety* property that it mark at least all live objects, and a *liveness* property that it terminates. These requirements can be satisfied by a variety of correct collectors with varying degrees of precision, efficiency, and atomicity.

The algorithmic differences between the various collectors are manifested in the additional unreachable objects that they mark (and thus retain). It is therefore natural to define relative precision between two collectors by comparing their *marked* sets at the end of the marking phase.

There is a trade-off between the precision of a collector and the degree of concurrency it provides. For example, a stop-the-world collector trades all concurrency for obtaining maximal precision (all unreachable objects are collected). Other algorithms provide a higher degree of concurrency (finer grained atomicity), at the expense of retaining more unreachable objects at the end of the marking phase.

In this paper we focus on the relative precision of algorithms under a given predetermined kind of atomicity. That is, the atomicity constraints of the generic algorithm are fixed and all instantiated algorithms in our parametric framework follow the same atomicity restrictions, namely the **atomic** blocks used in Figures 1 and 2. The precision-reducing transformations presented in the next section create opportunities for concurrency-increasing transformations. Although we do not deal with concurrency transformations which alter atomic sections, the instantiated algorithms do have a shorter duration of these atomic sections. As future work we plan on extending the parametric framework in this paper to include atomicity transformations.

An *algorithm* in our framework consists of the skeletons of Figures 1 and 2, instantiated with an arbitrary *expose* function. Note that this may or may not be a correct algorithm.

We now consider the question of the relative precision of two algorithms. Intuitively, a more precise algorithm should always mark fewer objects. However, this is *not* the case because the actual set of objects marked depends on the specific interleaving of mutator and collector. In fact, there exist algorithms and interleavings such that an algorithm that *always* selects more objects from the mutation log as origins for transitive marking will in fact mark fewer objects during the collection as a whole.

This is because what such a notion of precision compares is not necessarily the effect of the algorithm but may be the effect of arbitrary interleavings. Thus, a meaningful comparison must factor out the non-deterministic effects of particular executions caused by such random variations as temperature-induced changes in CPU speed, rather than the action of the collector itself.

We therefore consider algorithm C_2 to be less precise when for *any given global state* it exposes more objects for marking than algorithm C_1 . We therefore consider the effect of $expose_{C_1}$ on any interaction log obtained by C_2 , and show that when transitive marking is complete (when $pending = \emptyset$), $expose_{C_1}$ returns a subset of the objects returned by $expose_{C_2}$.

DEFINITION 4.1. *Given two collection algorithms C_1 and C_2 , we say that C_1 is more precise than C_2 , denoted $C_1 \sqsubseteq C_2$, when given any global state of C_2 with an interaction log l and where the set pending is empty, $expose_{C_1}(l) \subseteq expose_{C_2}(l)$.*

In the following section we present precision-reducing and correctness-preserving transformations of algorithms in our framework, and show that if the initial algorithm C_1 is correct, then the resulting algorithm C_2 is also correct.

5. Correctness-preserving Transformations

In this section, we present various transformations that can be combined to systematically derive safe collection algorithms from the

Apex collector. For each transformation, we show that its application is *correctness-preserving* and *precision-reducing*.

Each transformation is applied across a *dimension*. Dimensions are the formal analogue of the basic variables in the design of a collector as presented informally in the introduction. Specifically, we parameterize the collector with the following dimensions:

- **Wavefront:** how far has the collector progressed?
- **Policy:** how are modified objects behind the wavefront treated?
- **Threshold:** how large are cross-wavefront counts allowed to grow before they are “stuck”?
- **Protection:** which objects are traced to guarantee that all live objects are found?
- **Allocation:** how does the collector handle newly allocated objects to ensure timely termination?

A dimension is described by an ordered partition $\langle P_1, \dots, P_n \rangle$ of the objects in \mathcal{U} , where each subset of the partition corresponds to a different manner of handling objects in the dimension.

The subsets of a partition have the property that moving an object to a subset “to the right” yields an algorithm of lower precision. Formally, the relation between subsets is such that if $i < j$, $x \in P_i$, algorithm C uses partition $\langle \dots, P_i, \dots, P_j, \dots \rangle$ and algorithm C' uses partition $\langle \dots, P_i \setminus \{x\}, \dots, P_j \cup \{x\}, \dots \rangle$, then $C \sqsubseteq C'$.

For each dimension, we have generalized the algorithm so that all of the mechanisms represented by the subsets can be used simultaneously within a collector. In some dimensions, there is no restriction on the partitioning of \mathcal{U} ; for others, we specify a restriction formally.

All theorems stated in the paper are proved, but due to space restrictions are provided in an online supplement [42].

5.1 The Wavefront Dimension

The wavefront denotes the progress of the collector through the heap. We defined the precise wavefront in Section 2.3.1: it consists of the set of object fields that have thus far been traced by the collector. The *expose* functions determine how to handle mutations to the heap depending on whether they occur behind or in front of the wavefront.

The wavefront dimension represents different choices for the granularity at which the collector progress is tracked. Tracking the wavefront precisely may be inefficient because it requires per-field information, so it is sometimes desirable to sacrifice some precision in exchange for a more efficient implementation.

The wavefront dimension is an ordered partition of the objects in \mathcal{U} into: $D_W = \langle FL, OL \rangle$, where objects in the *FL* subset have their wavefront tracked precisely (at “Field Level”) as in Definition 2.2, while objects in the *OL* subset do not distinguish between fields within an object (and are tracked at “Object Level”).

One could further generalize this dimension to include all possible subsets of fields for all objects, but we do not consider this here for simplicity of presentation.

There are no restrictions on how the objects may be partitioned in the wavefront dimension.

5.1.1 Wavefront Abstraction

The wavefront abstraction transformation abstracts the exact collector wavefront, and tracks the collector’s progress at the granularity of an object rather than at the granularity of individual fields of objects. Given the set *FL* of objects for which the wavefront is to be maintained at the field level, we define the abstracted wavefront

as follows:

$$\begin{aligned} \mathcal{W}^>(P) &= \{(o, f) \mid (o, f) \in \mathcal{W}(P) \wedge o \in FL\} \cup \\ &\quad \{(o, f) \mid \exists f' \in Fields : (o, f') \in \mathcal{W}(P) \wedge \\ &\quad f \in Fields \wedge o \in OL\} \\ \mathcal{W}^<(P) &= \{(o, f) \mid (o, f) \in \mathcal{W}(P) \wedge o \in FL\} \cup \\ &\quad \{(o, f) \mid \forall f' \in Fields : (o, f') \in \mathcal{W}(P) \wedge \\ &\quad f \in Fields \wedge o \in OL\} \end{aligned}$$

The abstracted wavefront consists of two functions, $\mathcal{W}_{FL}^>(P)$ and $\mathcal{W}_{FL}^<(P)$. The function $\mathcal{W}_{FL}^>(P)$ over-approximates the set of object fields behind the wavefront. The function $\mathcal{W}_{FL}^<(P)$ under-approximates the set of objects fields behind the wavefront (and thus over-approximates the set of object fields ahead of the wavefront). Both functions are needed because *expose* functions that depend on an object’s field being behind or ahead of the wavefront must each use a conservative approximation.

EXAMPLE 5.1. Consider a prefix of the example interleaving of Fig. 3 just before tracing action (*T6*) is performed and the field $(A, f3)$ is traced. For this prefix P assuming $FL = \emptyset$ we get the following:

$$\begin{aligned} \mathcal{W}(P) &= \{(r1, f2), (r1, f3), (A, f1), (A, f2), (r1, f1)\} \\ \mathcal{W}^>(P) &= \{(r1, f2), (r1, f3), (A, f1), (A, f2), (r1, f1), (A, f3)\} \\ \mathcal{W}^<(P) &= \{(r1, f2), (r1, f3), (r1, f1)\} \end{aligned}$$

The wavefront abstraction transformation moves an object o from *FL* to *OL*.

THEOREM 5.2. *The wavefront abstraction transformation is a correctness-preserving and precision-reducing transformation.*

5.2 The Policy Dimension

Traditionally, when deciding how to protect against lost objects, implementers have thought in terms of the three classical types of write barriers: those of Dijkstra [21], which records the new pointer stored into an object; of Steele [40], which records the pointer to the object being modified; and of Yuasa [45], which records the old pointer that was overwritten.

However, while this decomposition may seem intuitive it does not in fact capture the essential properties of the design space in an orthogonal manner. Therefore, we introduce two separate and orthogonal dimensions which determine how objects are protected and do so in a manner that allows the different mechanisms to be composed.

In the Apex collector of Section 3, we used *rescanning* as a uniform policy for protecting all objects. The key to the simplicity of rescanning is that it finds all pointers to traverse in an atomic step of the *expose* operator.

However, this atomicity is costly and therefore rescanning is generally applied to some minimal portions of the memory, such as the stacks, for which it is practical to do so. The rest of the objects are processed incrementally from the log.

The policy dimension determines whether the modifications to a field are found by atomically scanning the heap (called “Scan-based Reachability”) or by examining the log (called “Log-based Reachability”).

This dimension is an ordered partition of objects in \mathcal{U} into

$$D_P = \langle SR, LR \rangle$$

The objects in *SR* are rescanned as described previously; the objects in the *LR* are discovered solely by processing the log, without accessing the contents of the heap.

There are no restrictions on how the objects may be partitioned in the policy dimension.

5.2.1 Rescanning

In order to define an *expose* function that works along this dimension, we first have to refine the simplified definition of a rescanning collector we presented in Section 3 and parameterize it according to the potentially imprecise wavefront and the partitioned policy dimension:

$$\begin{aligned} \text{expose}^r(P) &= \{o.f \mid P_i.\text{kind} \in \{\mathbf{M}, \mathbf{A}\} \\ &\quad \wedge P_i.\text{source} = o \wedge P_i.\text{field} = f \\ &\quad \wedge (o, f) \in \mathcal{W}^>(\text{pre}(P, i)) \wedge o \in SR \wedge P_i.\text{new} \in IS \\ &\quad \wedge 0 \leq i < |P|\} \end{aligned}$$

For the time being the set $IS = \mathcal{U}$; non-trivial use will be made of the set IS below for the protection dimension.

5.2.2 Maintaining Cross-Wavefront Counts

With rescanning, if a field of an object is modified repeatedly, rescanning will only see the final value when it processes the log. That is, there may have been intermediate values stored by the mutator and subsequently overwritten. We now describe a different way of discovering the resulting pointers, which is based on reference counting of pointers from behind the wavefront.

In particular, we observe that if a field initially contains a pointer p_0 and then has a sequence of pointers p_1, \dots, p_n written to it, then rescanning will find only p_n . If we were to apply a specialized form of reference counting, then the reference counts of pointers p_1, \dots, p_{n-1} would remain unchanged: they would be first incremented and then decremented. In the end, only the reference counts of p_0 and p_n would change, being decremented and incremented, respectively. This means that the intermediate reference counting operations can be ignored.

This observation is originally due to Barth [8], and is central to the multiprocessor reference counting algorithm of Levanoni and Petrank [33]. In our formulation, we use this approach to show how rescanning can be replaced by log-based reachability which keeps a count of references from behind the wavefront in order to identify exactly the same objects as are found by rescanning.

Note that this is *not a general form of reference counting*. Our framework only covers tracing algorithms. In particular, since counts are only maintained from behind to in front of the wavefront, there can be no cycles of objects with non-zero counts.

Existing approaches will subsequently be shown to be degenerate cases of reference counting in which the cross-wavefront count is a single sticky bit (expressed by the threshold dimension).

The counting-based approach has the enormous advantage of not requiring the synchronization of the *expose* function; on the other hand it will perform n steps for the mutated field described above, whereas rescanning would perform exactly one. One area in which this tradeoff manifests itself is in the treatment of stacks, whose high mutation rate makes them unsuitable for write barriers – instead, they are rescanned atomically.

Mutator Count: The mutator count is the number of pointers to an object from object fields behind the wavefront. This quantity is computed with respect to a given wavefront. We assume that some objects in the heap are *rescanned objects* that do not affect the mutator count. The mutator count computation is therefore parameterized by a set of objects LR from which the count should be computed. To compute the mutator count from a given log prefix P , we define the mutator-count increment and decrement as follows:

$$\begin{aligned} M^+(o, P) &= |\{P_i \mid P_i.\text{kind} \in \{\mathbf{M}, \mathbf{A}\} \wedge P_i.\text{new} = o \\ &\quad \wedge (P_i.\text{source}, P_i.\text{field}) \in \mathcal{W}^>(\text{pre}(P, i)) \\ &\quad \wedge P_i.\text{source} \in LR \wedge 0 \leq i < |P|\}| \end{aligned}$$

$$\begin{aligned} M^-(o, P) &= |\{P_i \mid P_i.\text{kind} \in \{\mathbf{M}, \mathbf{A}\} \wedge P_i.\text{old} = o \\ &\quad \wedge (P_i.\text{source}, P_i.\text{field}) \in \mathcal{W}^<(\text{pre}(P, i)) \\ &\quad \wedge P_i.\text{source} \in LR \wedge 0 \leq i < |P|\}| \end{aligned}$$

The value $M^+(o, P)$ is the number of new references introduced by the mutator from object fields that are behind the wavefront. Similarly, the value $M^-(o, P)$ is the number of references removed by the mutator from object fields behind of the wavefront. The mutator count is computed by combining the mutator-count increments and decrements as follows:

$$M(o, P) = M^+(o, P) - M^-(o, P)$$

EXAMPLE 5.3. Consider the example of Fig. 3 and its corresponding interaction $\log P^e$ as shown in Example 2.3. Assuming that $LR = \{A\}$ and $FL = \mathcal{U}$, the mutator count for B increases to 1 when the pointer from $(A, f1)$ is installed, and is decreased back to 0 as the pointer is deleted. Therefore, at the end of the prefix P^e , $M(B, P^e) = 0$. Note that the installation from $(r1, f1)$ does not increment the count, as the installation takes place ahead of the wavefront.

Collection by Counting: Using the formulation of Section 2.2, a counting-based collector can be instantiated using the following *expose* function. We use the superscript c to denote the fact that this function is based on counting, and name the function expose^c .

$$\begin{aligned} \text{expose}^c(P) &= \{n \mid n = P_i.\text{new} \wedge M(n, P) > 0 \\ &\quad \wedge n \in IS \wedge 0 \leq i < |P|\} \end{aligned}$$

EXAMPLE 5.4. Consider the example of Fig. 3. Assuming $LR = \mathcal{U}$, $FL = \mathcal{U}$, the function $\text{expose}^c(P^e) = \{E\}$.

However, since counting depends on the wavefront, taking a less precise wavefront can result with more objects being exposed. For example, taking $FL = \mathcal{U} \setminus \{r1\}$ results with $\text{expose}^c(P^e) = \{E, B\}$ as the count for B is incremented on the installation from $(r1, f1)$ that is behind the (overapproximated) wavefront, but cannot be decremented when the mutation $r1.f1 = \text{null}$ takes place, as $(r1, f1)$ is not behind the (underapproximated) wavefront.

Note that using a less precise wavefront resulted with additional objects exposed by the algorithm. In particular, in this case $\text{expose}^c(P^e) = \{E, B\}$ is a superset of the origins exposed by Apex algorithm on the same prefix as (see in Example 3.1).

We now formally define an *expose* function that works along the D_P dimension:

$$\text{expose}^{rc}(P) = \text{expose}^r(P) \cup \text{expose}^c(P)$$

The following theorem shows that moving along the D_P dimension is a precision reducing transformation.

THEOREM 5.5. *The rescanning to counting transformation, moving an object from SR to LR, is a correctness-preserving and precision-reducing transformation.*

It is interesting to note that in the special case in which a precise wavefront is maintained for all objects, and under an infinite mutator count, the precision of any partition along the D_P dimension is identical.

5.3 The Threshold Dimension

The threshold dimension represents different choices for the precision of maintaining the mutator count introduced in the previous section. In real systems, reference counts are usually very low, and the mutator count, which only counts references installed from behind the wavefront, will be even lower. Therefore, it would be very wasteful to have a reference count per object each capable of indexing the number of live objects in the heap.

The threshold limits the mutator count to a maximum value, at which it “sticks” and is not subsequently decremented. This allows the count to be represented with a fixed (small) number of bits while maintaining the correctness properties provided by reference counting.

The threshold dimension is an ordered partition of the objects in \mathcal{U} into:

$$D_T = \langle C_\infty, \dots, C_k, \dots, C_1 \rangle$$

where the subsets represent the count with successively less and less precision, which leads to collectors which are successively less precise, as we will show below. There are no restrictions on how the objects may be partitioned in the threshold dimension.

5.3.1 Abstracting Mutator Count

In practice, it is not possible to maintain an infinite mutator count. That is, it is beneficial to limit the mutator count $M(o, T)$ to a small bounded range.

Generally, the mutator count increment $M(o, T)$ can be abstracted to range over an interval $[0, k)$ and ∞ , defined as follows:

$$M_k(o, P) = \begin{cases} \infty, & \exists i. M(o, \text{pre}(P, i)) \geq k; \\ M(o, P), & \text{otherwise.} \end{cases}$$

To the best of our knowledge, all existing algorithms use the degenerate case where $k = 1$ and the mutator count is either 0 or ∞ , in which case the count is simply a flag that indicates whether a pointer to the object has been stored behind the wavefront. That is, immediately after a pointer to an object is stored behind the wavefront, the mutator count is set to a value that cannot be decremented.

Thus we will use this special case when presenting transformations to pre-existing algorithms such as those of Dijkstra, but it should be noted that 2- or 3-bit counts ($k = 3$ or $k = 7$) could be implemented efficiently and would likely provide most of the potential increase in precision available in practice.

EXAMPLE 5.6. Consider the example of Fig. 3 and its corresponding interaction log P^e . Assuming that $LR = \mathcal{U} \setminus \{r1\}$, $IS = \mathcal{U}$, and $FL = \mathcal{U}$, the function $\text{expose}^e(P^e)$ using $M_1(o, P)$ exposes B as $M(B, \text{pre}(P^e, 3)) = 1$.

We now show that the mutator count abstraction transformation preserves correctness and reduces precision.

THEOREM 5.7. *The mutator count abstraction, moving an object from C_k to C_{k-1} is a correctness-preserving and precision-reducing transformation.*

5.4 The Protection Dimension

Fundamentally, the synchronization problem of a concurrent collector is to prevent the mutator from “hiding” objects by moving pointers that are ahead of the wavefront to locations behind the wavefront, and then deleting the original pointer. Therefore, safety can fundamentally be guaranteed either by considering pointers installed behind the wavefront (installation-based protection), or by considering pointers deleted ahead of the wavefront (deletion-based protection).

Previously known collectors treat all pointers uniformly: so-called incremental update collectors (such as that of Dijkstra) use installation-based protection; snapshot collectors (such as that of Yuasa) use deletion-based protection. However, our framework allows the two approaches to be mixed, subject to some restrictions necessary for correctness.

The protection dimension is an ordered partition of \mathcal{U} into an Installation Set IS and a Deletion Set DS :

$$D_\pi = \langle IS, DS \rangle$$

The objects in IS are said to be *I-protected*, while the objects in DS are said to be *D-protected*.

The partition is restricted such that every live D-protected object is reachable from a sequence of D-protected objects (this is formalized below).

5.4.1 Snapshot-Based Collector

A snapshot-based collector marks as live all objects that were reachable at the start of the collection cycle; objects that become unreachable during the collection cycle are still treated as live [45].

Using the formulation of Section 2.2, a snapshot-based collector can be defined using the following *expose* function. We use the superscript d to denote the fact that this function is based on deletion, and name the function expose^d .

$$\text{expose}^d(P) = \{o \mid P_i.\text{kind} = \mathbf{M} \wedge P_i.\text{old} = o \\ (P_i.\text{source}, P_i.\text{field}) \notin \mathcal{W}^<(\text{pre}(P, i)) \wedge o \in DS \wedge 0 \leq i < |P|\}$$

Given a log prefix P , $\text{expose}^d(P)$ returns all objects in DS that were pointed-to by a field that was assigned a new value (possibly **null**) before it was scanned by the collector.

EXAMPLE 5.8. Consider the example of Fig. 3. Assuming that $DS = \mathcal{U}$, and $FL = \mathcal{U}$, $\text{expose}^d(P^e) = \{B, C, D\}$. Note that this exposes a superset of the origins exposed by the collector in Example 5.4 and by the Apex collector in Example 3.1.

5.4.2 Combinations of I-protected and D-protected Objects

Using the formulation of Section 2.2, a collector combining protection policies at the granularity of objects can be defined using the following *expose*:

$$\text{expose}^{rcd}(P) = \text{expose}^{rc}(P) \cup \text{expose}^d(P)$$

More importantly, we introduce a transformation which changes an object from I-protected to D-protected.

However, we must place an additional constraint on which objects in a given graph can be transformed from I-protected to D-protected. To guarantee that an object can be safely transformed from I-protected to D-protected, the object has to be transitively protected by a path of D-protected objects.

DEFINITION 5.9 (Valid Protection Sequence). *A valid protection sequence to an object x is a sequence of objects $o_1, \dots, o_k = x$ such that o_1 is a root object, and for every $1 \leq i < k$, there is a field f of o_i such that $o_i.f = o_{i+1}$ and $o_i \in DS$.*

DEFINITION 5.10 (Eligibility). *Given an object $x \in IS$, we say that x is eligible for membership in DS if there exists a valid protection sequence to x .*

Transformation along the protection dimension is significantly more complex than previous transformations because the transformed algorithm makes decisions in its *expose* function which may be locally more precise and yet are globally less precise. In particular, if a pointer to object X is stored behind the wavefront, and X is I-protected, then X will be exposed. But if X is D-protected, it will not be exposed. But since the object is D-protected, it will either be discovered through tracing either directly or through an overwritten pointer in its protection sequence.

In order to consider relative precision of algorithms with non-local effects, we need to refine Definition 4.1:

DEFINITION 5.11 (Weak Precision). *Given two collection algorithms C_1 and C_2 , we say that C_1 is weakly more precise than C_2 , denoted $C_1 \trianglelefteq C_2$, when given any global state of C_2 with an interaction log l and where the set pending is empty, $\text{expose}_{C_1}(l)^* \subseteq \text{expose}_{C_2}(l)^*$.*

That is, if the transitive closure of the objects exposed by C_1 is a subset of the transitive closure of the objects exposed by C_2 .

Weak precision is implied by the strong precision of Definition 4.1, which only consider the exposed objects and not their transitive closure. Since the previous transformations have been shown to be strongly precision reducing, they are also weakly precision reducing.

Note that there is a direct analogue between strong and weak precision, and the strong and weak white-black invariants of incremental update and snapshot collectors. In incremental update collectors all objects are I-protected; in snapshot collectors all objects are D-protected.

Under this refined definition, we show that the protection transformation is weakly precision reducing.

THEOREM 5.12. *Given an eligible object $x \in IS$, changing D_π from $\langle IS, DS \rangle$ to $\langle IS \setminus \{x\}, DS \cup \{x\} \rangle$ is a weakly precision-reducing and correctness-preserving transformation.*

5.5 The Allocation Dimension

To guarantee termination, an algorithm must provide a certain level of progress on each collector marking step. In the parametric algorithm of Fig. 1, the `mark()` procedure marks at least one object each time it is called. Therefore, if no new objects are allocated, the termination of the algorithm is guaranteed.

In the presence of allocation, however, additional mechanisms are required to guarantee termination. The impact of allocation on termination is that if the mutator allocation rate is faster than the collector tracing rate, the collector might need to trace through the newly allocated objects and thus does not terminate in a predictable manner.

The parametric collector presented in figure Fig. 1 solves this problem by introducing a synchronous termination phase after the main `while` loop in `collect`. Such approach is also used in [11, 7]. As noted in [11] the worst case of this approach can degenerate to an atomic processing of the whole heap in the case of Steele-like barriers.

However, although in practice this rarely happens it is important to explore options which guarantee timely termination for the worst case as well. This is particular important for real-time collectors where it is vital to guarantee a worst-case pause time.

Objects have typically been allocated white or black; we consider an additional kind of allocation (denoted by the color yellow) that provides an intermediate point in the trade-off space between precision and termination.

The dimension is an ordered partition of objects in \mathcal{U} into:

$$D_A = \langle WC, YC, BC \rangle$$

In our framework, all newly allocated objects are considered to be members of IS , that is I-protected objects.

5.5.1 Allocating Unmarked Ahead of the Wavefront (White)

The first approach is the least conservative towards towards marking allocated objects. Effectively, they are allocated unmarked, that is, *white*. In this case, the `mutateAlloc` need not make any special provisions for the object, that is, `mutateAlloc` simply invokes the `mutate` procedure.

The Apex algorithm presented earlier allocates *all* objects white. However, the negative impact on termination when allocating white is that the collector may need to trace through the objects. Allocating white is the primary reason for allowing the collector to non-deterministically enter the synchronous termination phase at the end of the `while` loop in Fig. 1. It could enter that phase after a fixed number of iterations of the while loop. In the Apex algorithm, as mentioned already, allocating white results

in the worst-case pause time being proportional to the size of the heap.

5.5.2 Allocating Unmarked Behind of the Wavefront (Yellow)

The termination problem introduced by objects which are allocated white is that the collector needs to trace through these objects to find other objects which are allocated white.

In order to avoid tracing through these white allocated objects, we introduce *yellow* objects. Yellow objects are allocated unmarked, but any references to objects in IS stored into a yellow object will be treated as if the yellow object is behind the wavefront. That means that pointers can not escape into a yellow object. In particular, it means that it is not possible for the mutator to create chains of unmarked objects that the collector must “chase”. Nevertheless, in contrast to grey objects, yellow objects can die in the collection cycle in which they were allocated.

The WC to YC transformation can significantly reduce the termination problems associated with allocating white: it eliminates the requirement on the `mark` to trace through a yellow object. To that end, when the collector encounters a yellow object, it does not place its fields in *pending*.

In the case where all objects are allocated yellow, the synchronous termination phase will not need to perform any transitive marking, since all objects reachable from live yellow objects will have already been marked. It need only consider the directly reachable yellow objects, greatly reducing the cost of the termination phase.

To understand why the WC to YC transformation leads to a less precise algorithm, consider the following simple example (object A is the object onto which the WC to YC transformations is applied): (i) allocate object A; (ii) store a pointer to an unmarked object $B \in IS$ into object A; (iii) delete all other pointers to B except the pointer from A. (iv) delete all pointers to object A; making A unreachable.

After object A becomes unreachable in the last step, if A was treated as a yellow object, object B will be retained. However, object B will not be retained if A is allocated white.

THEOREM 5.13. *The WC to YC transformation is a precision-reducing and correctness-preserving transformation.*

5.5.3 Allocating Marked (Black)

Termination can be further improved by taking allocated objects completely out of consideration. This approach is to allocate objects *black*. Black objects can be informally thought of as yellow objects which are allocated marked. That is, the `mutateAlloc` procedure is the same as the procedure for a yellow object, with the addition that the black object is marked upon allocation.

Unlike yellow objects, black allocated objects are always considered live for this collection cycle.

The main insight for the introduction of yellow and black objects is that they are not traced through. We therefore consider these objects outside SR .

The YC to BC will lead to even less work for the collector as it does not need to do additional work for these objects. Real-time collectors such as Metronome choose to allocate all objects black.

THEOREM 5.14. *The YC to BC transformation is a precision-reducing and correctness-preserving transformation.*

6. Collector Derivations

In this section, we explore a small subset of the space of concurrent collection algorithms along the dimensions of Section 5. The space we consider is depicted in Fig. 4. We will typically explore algorithms at the end points of a dimension. That is, we consider

the sets to be either \mathcal{U} or \emptyset . For presentational simplicity, we also assume that all derived algorithms use an abstracted wavefront (see Section 5.1.1) and mutator count abstracted to $k = 1$.

We first describe several well-known collector algorithms and where they fit into the lattice. We then discuss a few of the new practical collectors. The names of the new algorithms are depicted in boldface in Fig. 4.

In our parametric framework constructing new algorithms is a matter of choosing values over the various dimensions. For example, only recently a collector which uses a precise wavefront definition has been introduced in [18]. We can instantiate similar collectors which use a precise wavefront definition by setting \mathcal{W} to \mathcal{U} .

In the figure, we use a tuple of the form (SR, IS, FL, WC, BC) to define the point of the algorithm along the dimensions of Section 5. The values for other sets along each dimension are defined as complements using the values in the tuple, e.g., if $WC = \emptyset$ and $BC = \emptyset$, then $YC = \mathcal{U}$. In the tuple, we use the set *stacks* to denote the set of stack objects, and \mathcal{A} to denote the set of newly allocated objects. Additionally, every edge represents a precision order relation.

6.1 Existing Algorithms

A Steele-style algorithm can be derived from the Apex collector by applying the *wavefront abstraction* transformation to all objects.

A Dijkstra-style algorithm is derived from the Steele-style collector by moving all objects except stacks along the D_P dimension, that is, from SR to LR . It is less precise than the Steele-style collector.

The Yuasa algorithm is the least precise of the three existing algorithms, it allocates all objects black (i.e. $BC = \mathcal{U}$) and in addition all existing objects are D -protected.

6.2 New Algorithms

Fig. 4 contains several new algorithms of practical importance. In this section we informally describe some of those new collectors.

The Steele-YC collector is derived from the Steele-like algorithm by applying the WC to YC transformation to all allocated objects. This algorithm bounds the duration of the synchronous termination phase, thus addressing the main issue of algorithms that use a Steele-like write barrier (regardless of the granularity of rescanning) such as [11, 7]. The disadvantage of this collector is that it might retain more unreachable objects than Steele.

The Steele-BC algorithm makes an even more conservative assumption in regards to allocated objects. This leads to an opportunity to reduce the work for termination even further while still retaining relatively high precision for existing heap objects. This algorithm could be beneficial for applications where most of the allocated objects are long lived (i.e. do not die during the collection cycle), such as the mature space of generation collectors.

The precision of the Steele-BC algorithm can be reduced further (and hence the potential for concurrency is increased) by moving along the D_P dimension, moving all objects but stacks from SR to LR resulting with the Dijkstra-BC algorithm.

The Steele-D algorithm, derived from Steele by moving “to the right” on the D_π dimension, uses synchronous termination and at the same time considers all existing objects as live. However, the algorithm Steele-D-YC which is derived from Steele-D bounds that synchronous phase.

7. Related Work

In previous work [43] we observed a common structure between concurrent collectors and suggested that they can possibly be viewed as instances of a more abstract collector. However, the

paper effectively contained two very complex abstract algorithms, and a few discontinuous “transformations” where their application was only described informally. Moreover, the resulting collectors could not be related.

In [10], separation logic is used to prove the correctness of a stop-the-world copying garbage collector. However, with the extension of separation logic to concurrency [14], it may be possible to formally prove the Apex algorithm presented in this work.

Another work modelling collectors is [12]. In this paper, the authors use CCS to specify a stop-the-world collector, and temporal logic to specify its liveness and safety properties. However, the presented algorithm is not concurrent and although the collector is specified in CCS, there is no attempt at verifying the presented algorithm. The authors do note however that proving the correctness of a concurrent collector would be even more challenging.

Several works formally verify the correctness of Ben-Ari’s and Dijkstra’s algorithms [9, 21]. The focus of Ben-Ari’s algorithm is correctness rather than efficiency. However, both of these algorithms are not practical because their worst-case time complexity is quadratic in the size of the heap.

In [37], Ben-Ari’s algorithm is verified for both single and multi-mutator systems using Owicki-Gries’s logic in the HOL theorem proving system. In the work of [25], again Ben-Ari’s algorithm is verified using the PVS theorem proving system. Similar work has been done by [38], where he proves Ben-Ari’s algorithm but this time in Boyer-Moore’s theorem prover. In [28], Dijkstra’s algorithm has been verified again in the PVS theorem prover. The paper of [15], proves Ben-Ari’s algorithm using the B and Coq systems. These works are complementary to ours in the sense that they concentrate on formally proving a particular collector algorithm. The formal systems used for that purpose may be utilized in proving Apex correct. The works of [19, 20] define a framework to describe generational and conservative collectors. However, it only deals with stop-the-world algorithms. Another transformational approach to collectors can be found in [20]. The authors use the SETL wide spectrum language to specify an initially correct and inefficient implementation of a stop-the-world collector. Through loop fusion and formal differentiation transformations, they obtain a more precise implementation of a well-known stop-the-world algorithm. The transformations in our work are specific to the world of concurrent marking collectors.

8. Conclusions and Future Work

In this paper we presented a mechanism to automatically generate various correct concurrent garbage collector algorithms.

The mechanism consists of starting with a correct initial algorithm and applying a set of correct object-level transformations, with each application yielding a new concurrent algorithm.

We also introduced a definition of precision which allows us to formally relate the various algorithms. For understanding purposes, it is important to follow a structured approach in designing such complex algorithms.

In the future, we plan on working on how to automatically arrive at practical synchronization skeletons from our trace-based collectors, as well as relax the atomicity constraints in the current skeleton algorithm.

A. Proofs

In the following, we abuse notation and use the Kleene star to denote the transitive closure of objects reachable from a given set of objects. For example, we write S^* to denote all objects that are transitively reachable from the set of objects S .

A.1 Collector Invariants and Proof Methodology

The correctness of algorithms in our framework hinges on $expose(l)$ exposing all hidden origins. An algorithm in our framework is correct if and only if the following invariants hold immediately after computing the set of origins by performing $expose(l)$:

I-Invariant $expose(l)$ contains all unmarked objects in IS pointed-to by a marked object.

D-Invariant any unmarked object in DS is either returned by $expose(l)$ or is reachable from a DS object in $expose(l)$ by a path of objects in DS .

These invariants imply the intuitive notion of the algorithm marking a superset of all objects which are required to determine correct transitive reachability.

Our proofs work by showing that the transformations preserve the above invariants. We will denote as C_1 the algorithm before the transformation is applied to a single object o , and C_2 , the algorithm after the transformation is applied. The proofs use the correctness of C_1 to show the correctness of C_2 . Moreover, the correctness proofs also show that C_1 is more precise than C_2 .

We compare the executions of C_1 and C_2 starting from the same initial heap H , sequence of mutations M , and set of roots R .

In our framework, algorithms only differ in their $expose(l)$ function and therefore we compare their executions by comparing the corresponding $expose(l)$ at the point of divergence of the two traces (that is, when $pending$ is empty). For all transformations except white-to-yellow, the divergence point occurs when the $mark()$ procedure has finished and the algorithms proceed into the $addOrigins()$ procedure and compute $expose(l)$.

Let $expose_{C_1}$ denote the origins computed by C_1 and $expose_{C_2}$ the origins computed by C_2 with the same $log\ l$.

Certainly, if $expose_{C_1} = expose_{C_2}$, then the algorithms do not diverge and they can continue to execute in a lock-step. If throughout the entire execution, the algorithms do not diverge, then the execution of C_2 is identical to the execution of C_1 and hence is shown to be correct. Intuitively, in this case, the algorithms are of the same precision.

However, if $expose_{C_1} \neq expose_{C_2}$, the following invariants hold (we use the subscripts C_1 and C_2 to denote the various sets in C_1 and C_2 respectively):

- $marked_{C_1} = marked_{C_2}$
- the I-invariant and the D-invariant hold for C_1

The first step to showing that C_2 is correct is proving that I- and D- invariants hold for C_2 at the point of $expose$. For all transformations but the protection-transformation, we prove this by showing that $expose_{C_1} \subset expose_{C_2}$. This also shows that $C_1 \sqsubseteq C_2$.

The second step of the proof involves reasoning about the continuation of C_2 . That is, we need to find a corresponding witness trace which is also correct so that at the next point of $expose$, we can repeat this process. In all proofs except the protection and white-to-yellow transformations, the correct witness trace is basically the restart of C_1 with the *origins* resulting from $expose_{C_2}$ of C_2 . The C_1 algorithm can be restarted with the new state, because the I- and D- invariants are preserved by the transformations.

A.2 Wavefront Abstraction

The following proof shows the correctness of the wavefront abstraction transformation which takes a single object $o \in FL$ in C_1 and moves that object so that $o \in OL$ in C_2 . It also shows that C_2 is less precise than C_1 according to 4.1.

Proof:(Theorem 5.2) (sketch)

The application of this transformation on object o potentially affects the marking decision for heap objects other than o , but does not affect the marking decision for o itself. This is because the transformation takes effect once object $o \in marked$ and at least one $(o, field) \in \mathcal{W}^>$.

For this transformation, the divergence point always occurs after a call to $expose$ in $addOrigins()$.

Let P be the common prefix of C_1 and C_2 just before $expose$ is called with an interaction $log\ l$. At the point after the call to $expose$ where $expose_{C_2} \neq expose_{C_1}$, the computation $M^-(x, P)$ for any object $x \in IS$ indicates that any pointer installation to x into o will be returned by $expose$, provided that at least one $(o, field) \in \mathcal{W}$ and not all fields of $o \in \mathcal{W}$ at the time of the mutator operation. When object o is in that state, no destruction of a pointer to x in o can affect $M^-(x, P)$. This is indicated by the requirement for $(P_i.o, P_i.field) \in \mathcal{W}^<(pre(P, i))$ in $M^-(x, P)$.

In addition, operations on objects in DS also affect the return of $expose$. If a mutator stores a pointer to an object $d \in DS$ into $(o, field) \in \mathcal{W}$, and the mutator subsequently destroys that pointer to d , then this pointer would not be returned by $expose_{C_1}$. However, such a mutation sequence would cause the object d to be returned by $expose_{C_2}$.

The above discussion applies when $o \in LR$. In the case where $o \in SR$, and we are manipulating a pointer to an object in IS , $expose_{C_1} = expose_{C_2}$. It is only possible that more rescanning work will be done in $expose_{C_2}$ for o , but this will not affect the result of $expose_{C_2}$. However, this is not the case when target objects in DS are manipulated. Even if $o \in SR$, all $d \in DS$ removed from o will be returned by $expose_{C_2}$ as indicated by $expose^d(P)$.

Either way, at the point of divergence (where $pending(C_2)$ is empty), $expose_{C_1} \subseteq expose_{C_2}$. This is also the necessary condition to establish that $C_1 \sqsubseteq C_2$. However, the converse statement clearly does not hold, that is, $C_2 \not\sqsubseteq C_1$.

It is worth noting that although $expose_{C_1} \subset expose_{C_2}$, we cannot detect whether the objects in $expose_{C_2} - expose_{C_1}$ are unreachable. We have constructed examples which show that it is possible for all such objects to be unreachable or for all objects to be reachable or for the mix of the two to occur. However, because $expose_{C_1} \subset expose_{C_2}$, we can deduce that at the point in the trace right after the call to $expose$ in C_2 , the I- and D- invariants are satisfied. This is clear because the invariants are satisfied at the same point in $expose$ of C_1 and by the subset relation we can now trivially conclude that they also hold at the corresponding point in C_2 since adding pointers cannot cause an invariant violation.

Because the invariants are satisfied at this point, we can restart the execution of C_1 with the origins as returned by $expose_{C_2}$. Additional origins cannot violate the invariants and we therefore consider this to be a safe witness trace. We compare the possible continuations of C_2 to the restarted trace of C_1 with additional origins. However, from this point on, all fields of $o \in \mathcal{W}^<$ and $o \in marked$. That is, o will behave identically in both, the set of restarted C_1 traces versus the set of traces representing the continuations of C_2 . Subsequently, the traces of the safe restarted algorithm of C_1 are exactly the same as the continuations of C_2 and hence we can deduce that C_2 is correct. Additionally, because the continued traces are the same, $expose_{C_1}$ on any subsequent state of C_2 will be the same as $expose_{C_2}$ when $pending(C_2)$ is empty and thus satisfy $C_1 \sqsubseteq C_2$.

A.3 Collection by Counting

The following proof shows the correctness of the rescanning-to-infinite counting transformation which takes a single object $o \in SR$ in C_1 and moves that object so that $o \in LR$ in C_2 assuming

$IS = C_\infty$. It also shows that C_2 is less precise than C_1 according to 4.1.

Proof:(Theorem 5.5) (sketch)

Similarly to the wavefront abstraction, this transformation affects the marking decision for heap objects other than o , but does not affect the marking decision for o . The transformation takes effect when pointers to IS objects are stored in o .

Let P be the common prefix between C_1 and C_2 right before the point of divergence (right before $expose$) with an interaction $\log l$. If $o \in SR$, the M^+ and M^- computations are not activated. That is, the proposition $P_i.o \in LR$ will be evaluated to false. Therefore, the intermediate installations to objects $\in IS$ into a rescanned object o or intermediate deletions from o of objects $\in IS$ do not affect the mutator count (M) computation for these objects. That is, $expose^r(l)$ returns only the objects $\in IS$ pointed to from o at the time of the call.

In the case where $o \in LR$ and $o \in FL$, $expose_{C_1} = expose_{C_2}$. This is the case where o maintains a precise wavefront. Although, M^- and M^+ computations are activated because $o \in LR$, the intermediate installations cannot affect the result of $expose_{C_2}$. That is, $expose_{C_2}$ will always return precisely the pointers to IS objects residing in o at the time of the call to $expose$.

In the case where $o \in OL$, that is, when the wavefront for o is imprecise, a superset of the pointers to IS objects stored in o will be returned by $expose_{C_2}$. That is because, even if infinite count is available for an IS object stored in o , we can no longer decrement its count if the removal occurs when object o is being processed (i.e. in a state where some of its fields are processed but not all fields of o are processed yet). With C_1 , where $o \in OL$ and $o \in SR$, an imprecise wavefront can only cause additional work, but cannot affect the result of $expose_{C_1}$.

Therefore, we obtain $expose_{C_1} \subseteq expose_{C_2}$ at the divergence point and I- and D- invariants are satisfied. We have now shown the required condition for precision and therefore we can establish that $C_1 \sqsubseteq C_2$. The converse statement does not and therefore $C_2 \not\sqsubseteq C_1$.

Similarly to the wavefront abstraction, we cannot detect whether the objects in $expose_{C_2} - expose_{C_1}$ are reachable or not. However, one difference is that in this transformation, objects in DS are not affected. That is, $expose_{C_1} \wedge DS = expose_{C_2} \wedge DS$. Effectively, $expose_{C_2} - expose_{C_1} \in IS$. This is because both C_1 and C_2 are working with the same wavefront and hence $expose^d(l)$ in C_1 and C_2 are identical.

After the first divergence point where the invariants still hold, the reasoning is the same as in the wavefront abstraction since object $o \in marked$ and $o \in \mathcal{W}^<$ and hence even if $o \in OL$, o cannot create further differentiation in the continuations of C_2 versus the restarted traces of C_1 (recall that the 'difference' for this transformation occurs only because of an imprecise wavefront for object o). Similarly, because the continued traces are the same, $C_1 \sqsubseteq C_2$ cannot be violated.

A.4 Abstraction of Mutator Count

The following proof shows the correctness of the mutator count abstraction on a single object $o \in IS$. The transformation takes a single object $o \in C_k$ and moves that object to the set C_{k-1} in C_2 . It also shows that C_2 is less precise than C_1 according to 4.1.

Proof:(Theorem 5.7) (sketch)

The proof follows the same lines as the proofs for the previous two transformations. The difference however is that at the point of divergence, $expose_{C_2} - expose_{C_1} = o$. That is, the difference is a single object rather than a set of objects.

The traces which cause a divergence point between C_1 and C_2 are due to installations of a pointer to o into $(field, source) \in \mathcal{W}^>$. At the point before calling $expose$, there exists a common prefix P with a $\log l$, which results in $M_k(o, l) = 0$ and

$M_{k-1}(o, l) = \infty$ in which case $o \in expose_{C_2}$ and $o \notin expose_{C_1}$ and hence $expose_{C_1} \subset expose_{C_2}$. Hence, at the point where $pending$ is empty, $expose_{C_1} \subseteq expose_{C_2}$ and therefore we can establish that $C_1 \sqsubseteq C_2$. The converse statement does not and therefore $C_2 \not\sqsubseteq C_1$.

The restarting portion is similar to the wavefront in the sense that because object o is returned by $expose_{C_2}$, it will be marked in C_2 and hence it will have no further effect when comparing continuations of C_2 versus safe restarted traces of C_1 . Similarly, because the continued traces are the same, $C_1 \sqsubseteq C_2$ cannot be violated.

A.5 Counting to Snapshot

The following proof shows the correctness of the protection transformation on a single object o . In C_1 , $o \in IS$, while in C_2 , $o \in DS$. This transformation is the only one which changes the invariant mapping between objects. That is, in C_1 , object o satisfies the I- invariant while in C_2 , o satisfies the D- invariant. It also shows that C_2 is weakly less precise than C_1 according to 5.11.

Proof:(Theorem 5.12) (sketch)

Let P denote the common prefix trace with $\log l$ before the divergence point. With this transformation, it is possible to obtain two cases: one where $expose_{C_1} \subseteq expose_{C_2}$ and another where $expose_{C_2} \subseteq expose_{C_1}$.

Case I - $expose_{C_1} \subseteq expose_{C_2}$: This case occurs when a pointer to the transformed object $o \in DS$ has been removed from a $(source, field) \notin \mathcal{W}^<$. Therefore, $o \in expose_{C_2}$. However, in C_1 , $o \in IS$ and o will be returned by $expose_{C_1}$ only if $M_k(o, P) > 0$. That is, in C_2 , o is returned by $expose$ unconditionally, while in C_1 , there is a condition that o 's count is positive.

Therefore, in this scheme, if $expose_{C_1} \subset expose_{C_2}$, then $o = expose_{C_2} \setminus expose_{C_1}$ and since the invariants are satisfied in C_1 at this point, it is clear that they are satisfied in C_2 as well. In addition, o satisfies the D-invariant in C_2 . Since $expose_{C_1} \subseteq expose_{C_2}$ when $pending$ is empty, in this case, we can say that $C_1 \sqsubseteq C_2$ and $C_2 \not\sqsubseteq C_1$.

Similarly to the previous transformations we cannot actually detect whether o is actually reachable. We have constructed complex examples showing that o can be both floating garbage as well as live at the divergence point.

The restarting argument is the same as the previous transformations : object o will be marked by C_2 and therefore cannot affect the invariants and restarting C_1 with the roots returned by $expose_{C_2}$ is safe. In effect, the original algorithm C_1 continues with the additional roots returned by $expose_{C_2}$. Since o is marked in both the continuations of C_2 and the restarted traces of C_1 , there can be no further divergence points and therefore C_2 is safe. Similarly, because the continued traces are the same, $C_1 \sqsubseteq C_2$ cannot be violated.

Case II - $expose_{C_2} \subseteq expose_{C_1}$: This case occurs when $o = expose_{C_1} - expose_{C_2}$. This situation can arise when the mutators have *not* removed a pointer to the transformed object o ahead of the wavefront, that is, from $(source, field) \in \mathcal{W}^<$. However, the mutators have installed pointers to the transformed object o behind the wavefront, that is, in $(source, field) \in \mathcal{W}^>$, which results in $M_k(o, P) > 0$.

Because of the D* invariant in C_2 , we know that there must exist at least one object $p \in expose_{C_2}$ such that $p \in DS$ and o is reachable from p via a sequence of objects $\in DS$. Therefore, because $expose_{C_2} = expose_{C_1} \setminus o$ and the D- invariant holds for o , it is clear that C_2 satisfies both invariants at this point. However, we cannot say that at this point $C_1 \sqsubseteq C_2$ and therefore we cannot apply the precision definition 4.1 to this case. Instead, because of

the D^* invariant, we know that $expose_{C_1}(l)^* = expose_{C_2}(l)^*$ and therefore $C_1 \sqsubseteq C_2$.

However, unlike case I, it is not possible to restart C_1 with the origins as returned by $expose_{C_2}$. Restarting C_1 requires placing o back into IS and this would be incorrect since it would lead to a violation of the I-invariant in C_1 . This is because $M_k(o, P)$ may be > 0 at this point, i.e. o could be reachable from $(source, field) \in \mathcal{W}^<$ yet $o \notin expose_{C_2}$. Therefore, moving o from DS into IS in order to restart C_1 is unsafe.

Instead of restarting at the point of divergence, C_2 can continue its operation in a lock step with C_1 despite the fact that C_1 will mark the transformed object o right after the call to $expose_{C_1}$, while C_2 will not mark that object in $addOrigins()$. However, besides the marking of o , the traces have no other difference. At some point in the execution of C_2 before the next $expose$ point is reached, either C_2 will encounter o and mark it, or a path from an object in DS will be destroyed and C_2 will not encounter o . In the case where C_2 marks o , the traces of C_1 and C_2 will become identical after this point. In the case where C_2 reaches $expose$ before marking o , C_1 would not have yet reached $expose$ as it has to process o before doing so. In this execution however, we know that the invariants hold just before the call to $expose_{C_2}$. This is because C_1 and C_2 have been running nearly identical traces except with the marking of o in C_1 . However, C_1 has not yet started processing the object, i.e. $fields(o) \notin \mathcal{W}^>$ and therefore it could not have had transitive affect on other objects stored in o or deleted from o . Also, even if object o became unreachable, it is already marked and will be processed in C_1 . At any rate, the result of $expose_{C_2}$ at the point where $pending$ is empty will be a superset to the result of $expose_{C_1}$ if we had executed it on the trace of C_2 up to that point. That is, the invariants after the call to $expose_{C_2}$ are satisfied.

Clearly, because the invariants are satisfied in C_1 as well, we can continue the C_1 execution with the additional roots returned by $expose_{C_2}$ as additional pointers cannot violate the invariants. This process continues until either o is reached in the tracing phase and marked by C_2 or o is returned by $expose_{C_2}$ in which case it will be marked. Once o is marked, the next $expose$ of C_1 and C_2 will return identical results the algorithms will behave identically.

A.6 White to Yellow

The following proof shows the correctness of the white to yellow transformation which takes a single object $o \in WC$ in C_1 and moves that object so that $o \in YC$ in C_2 . Also, $o \in IS$ for both algorithms and $o \in FL$. It also shows that C_2 is less precise than C_1 according to 4.1.

Proof:(Theorem 5.13) (sketch) Similarly to the wavefront abstraction, the application of this transformation on object o can potentially affect other heap objects. However, unlike wavefront, this transformation is activated before $o \in marked$. That is, the effect of the transformation can take place immediately after the object is allocated. This implies that unlike wavefront, object o can be affected as well (i.e. a cyclic pointer can be installed in o).

For this transformation, the divergence points can occur in two places: (i) after the call to $expose$ in $addOrigins()$, and (ii) when the collector marks o in its tracing phase.

In the first case, because $o \notin marked$, no $(o, field) \in \mathcal{W}^>$. However, it is possible to obtain a divergence point in C_2 if \exists an object q such that $q \in IS$ and $q \in \mathcal{C}_k$. It is then possible to find a common prefix P such that $M_k(q, P) = \infty$ in C_2 while $M_k(q, P) = 0$ in C_1 . Such traces lead to $expose_{C_1} \subset expose_{C_2}$ because in C_1 no fields of o have been processed yet and therefore o cannot affect $M_k(q, P)$.

In the second case, o is encountered and marked in the tracing phase or is returned by $expose$ in $addOrigins()$. In either case,

both, C_1 and C_2 will mark o , but C_2 will not place o 's field in $pending$ as yellow allocated objects are not traced through in order to guarantee termination without requiring a stop-the-world phase to retrace the whole heap.

At this stage, similarly to case (ii) of the protection transformation, although a divergence point has occurred, the algorithms can keep running in a lock-step. In effect, for every trace of $t_2 \in C_2$, before $expose_{C_2}$, there exists a trace $t_1 \in C_1$ such that t_2 is a subtrace of t_1 , where the only difference between t_1 and t_2 is the value of the $pending$ set, that is $pending(C_1) = pending(C_2) \cup fields(o)$. When C_2 reaches $expose$, the invariants certainly hold for both C_1 and C_2 at their corresponding points in the trace. For C_1 , that point will be just before fetching the next field of o from $pending$. The reason is that $marked(C_1) = marked(C_2)$ and $expose_{C_2}$ contains all objects $\in IS$ pointed to from o . An important observation here is that we are not interested in objects in DS pointed to from o . Those objects are installed in o during this cycle and hence not rescanning o in order to return them in $expose_{C_2}$ cannot violate the D-invariant. That is, their destruction cannot cause the violation of the D^* invariant for objects $\in DS$.

Because the invariants are satisfied for C_2 , we can restart C_1 with the new state as returned by $expose_{C_2}$. After the restart, the algorithms behave identically because o is marked and processed.

The precision reasoning here is very similar to other transformation which affect only objects in IS such as wavefront. When we reach $pending$ being empty in C_2 , $expose_{C_1} \subseteq expose_{C_2}$ and hence $C_1 \sqsubseteq C_2$.

A.7 Yellow to Black

The following proof shows the correctness of the yellow to black transformation which takes a single object $o \in YC$ in C_1 and moves that object to BC in C_2 . Also, $o \in IS$ for both algorithms as o is a newly allocated object for this collector cycle. It also shows that C_2 is less precise than C_1 according to 4.1.

Proof:(Theorem 5.14) (sketch) The proof for this transformation is simpler than the previous transformations. It shows how given any trace $t_2 \in C_2$, we can find a corresponding trace t_1 in C_1 such that $marked(t_1) \subset marked(t_2)$.

After o is allocated, the effect of o on other objects in the heap is identical for both C_1 and C_2 . However, the traces for C_2 and C_1 diverge since $marked(t_1) = marked(t_2) \setminus o$ at that point. From this point on however, we can match every step of t_2 to a step in t_1 .

If C_1 encounters o and marks it, then, from that point, the suffixes of t_1 and t_2 are identical and $marked(t_1) = marked(t_2)$.

However, after allocating o , it is possible that the object is made unreachable and will not be encountered further by C_1 . For example, o can be stored in $(source, field) \notin \mathcal{W}^>$ and subsequently removed from that same field $(source, field) \notin \mathcal{W}^>$. In that case $expose_{C_1} = expose_{C_2}$ for all subsequent calls to $expose$ and $marked(t_2) = marked(t_1) \setminus o$ after the algorithm terminates.

However, it is possible to obtain a case where $o \in expose_{C_1}$ and $expose_{C_1} = expose_{C_2}$, but o has not yet been marked in C_1 . In that case, o will be marked subsequently after the call to $expose$ in $addOrigins()$ in C_1 and right after marking it $marked(t_1) = marked(t_2)$. The suffixes of t_1 and t_2 after this point will be identical.

Therefore, for any trace $t_2 \in C_2$ we can find the corresponding trace $t_1 \in C_1$ which marks less objects. More precisely if t_1 marks less objects than t_2 , then $marked(t_2) = marked(t_1) \cup o$.

Because object o is marked as soon as it is allocated in C_2 , $expose_{C_2}$ can return that object at any point (it will be eliminated immediately from consideration when $origins$ are computed. Alternatively, $expose$ can be extended to return a marked o as part of its result. Either way, when $expose_{C_1} \subset expose_{C_2}$, then

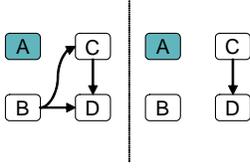
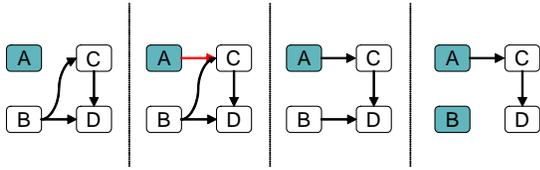


Figure 5. CASE I

$expose_{C_1} = expose_{C_2} \cup o$ and otherwise $expose_{C_1} = expose_{C_2}$. Therefore, when $pending(C_2)$ is empty, $expose_{C_1} \subseteq expose_{C_2}$ and hence $C_1 \sqsubseteq C_2$.

B. Examples: Counting to Snapshot

The following two examples illustrate the two cases in the counting-to-snapshot transformation. The first example shows how one can obtain $expose_{C_1} \subset expose_{C_2}$ (case I), and the second example shows how we can arrive at $expose_{C_2} \subset expose_{C_1}$ (case II).

B.1 CASE I

Consider the top sequence depicted in Fig.5 where $A \in DS$, $B \in DS$, $C \in IS$ and D is the transformed objects, that is, $D \in IS$ in C_1 and $D \in DS$ in C_2 . Object A has been marked and scanned.

In the first interleaving (top sequence), at the divergence point (after the fourth step), we get $expose_{C_1} = \{C\}$ and $expose_{C_2} = \{C, D\}$. Note that $expose_{C_2} \setminus expose_{C_1} = \{D\}$. In this example, object D is live.

In the second interleaving (bottom sequence), at the divergence point (after the second step), we get $expose_{C_1} = \emptyset$ and $expose_{C_2} = \{D\}$. Note again that $expose_{C_2} \setminus expose_{C_1} = \{D\}$. However, in this case object D is dead.

These two examples demonstrate how we can arrive at $expose_{C_1} \subset expose_{C_2}$ and moreover that we cannot deduce whether the converted object is live or not at the point where $expose$ is performed.

B.2 CASE II

Consider the sequence depicted in Fig.6 where $A \in DS$, $B \in DS$, $C \in DS$ and D is the transformed object, that is, $D \in IS$ in C_1 and $D \in DS$ in C_2 . Object A has been marked and scanned.

At the divergence point (the pointer after the 3rd step in the figure), we get $expose_{C_1} = \{C, D\}$ because a pointer to C is removed ahead of the wavefront and a pointer to $D \in IS$ is installed behind the wavefront. We also get $expose_{C_2} = \{C\}$. We do not get D because D in C_2 is $\in DS$ and no pointer to it has been destroyed. Note that $expose_{C_1} \setminus expose_{C_2} = \{D\}$.

In this example, object D is live because it is pointed to from object A . However, if the pointer from A to D was destroyed, $expose_{C_2}$ would not have returned D since that pointer is behind the wavefront and it is possible that D has become dead.

This example shows how we can arrive at $expose_{C_2} \subset expose_{C_1}$. It also demonstrates that we cannot reason whether

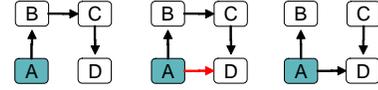


Figure 6. CASE II

the converted object is live or not at the point where $expose$ is performed.

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