

Inlining-Benefit Prediction with Interprocedural Partial Escape Analysis

Matthew Edwin Weingarten* Oracle Labs Zurich, Switzerland matthew.weingarten@inf.ethz.ch Theodoros Theodoridis

Swiss Federal Institute of Technology Zurich, Switzerland theodoros.theodoridis@inf.ethz.ch Aleksandar Prokopec Oracle Labs Zurich, Switzerland aleksandar.prokopec@oracle.com

Abstract

Inlining is the primary facilitating mechanism for intraprocedural Partial Escape Analysis (PEA), which allows for the removal of object allocations on a branch-by-branch basis and is critical for performance in object-oriented languages. Prior work used interprocedural Escape Analysis to make inlining decisions, but it discarded control-flow-sensitivity when crossing procedure boundaries, and did not weigh other metrics to model the cost-benefit of inlining, resulting in unpredictable inlining decisions and suboptimal performance. Our work addresses these issues and introduces a novel Interprocedural Partial Escape Analysis algorithm (IPEA) to predict the inlining benefits, and improve the cost-benefit model of an existing optimization-driven inliner. We evaluate the implementation of IPEA in GraalVM Native Image, on industry-standard benchmark suites Dacapo, ScalaBench, and Renaissance. Out of 36 benchmarks with a geometric mean runtime improvement of 1.79%, 6 benchmarks achieve an improvement of over 5% with a geomean of 9.10% and up to 24.62%, while also reducing code size and compilation times compared to existing approaches.

CCS Concepts: • Software and its engineering \rightarrow Source code generation; Runtime environments; Just-in-time compilers.

Keywords: escape analysis, inlining, compilers, graalvm

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*Also with Swiss Federal Institute of Technology.

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1 Introduction

Partial Escape Analysis (PEA) [50] is a control flow sensitive variant of Escape Analysis [7–9, 11, 12, 18, 25, 44, 55]. The analysis computes whether an object escapes outside the scope of the procedure and performs optimizations such as stack allocation for objects, replacing object fields with scalars, and eliding locks. Flow-sensitivity significantly increases the escape-analysis effectiveness, and decreases the dynamic allocation rate compared to flow-insensitive analysis [50]. This is achieved by allocating, or *materializing*, objects on some control-flow paths and keeping them *virtual* on others. PEA is an intraprocedural phase, so its potency is limited by calls to other procedures since almost all objects escape into at least one other procedure [25, 50].

In existing work, this limitation is mitigated with interprocedural analysis that foregoes control-flow sensitivity when crossing procedure boundaries [7, 11, 25]. The results of the analysis are then used to either guide inlining decisions [25] and then applying PEA within the scope of inlined procedures, or to employ alternative schemes such as stack allocation or region-based memory management, but without doing the inlining [11].

Although not mutually exclusive with interprocedural analyses, intraprocedural analysis is the more pragmatic choice for most compiler optimizations. In most optimizing compilers, optimization passes are intraprocedural [1, 15, 27, 32], and inlining kickstarts their symbiotic interplay. This is particularly prevalent for the optimizations performed by PEA, as one can easily see how scalar replacement consequently enables constant folding, conditional branch elimination, value numbering, and many others [36]. Most compilers therefore perform inlining as one of the earliest optimization passes, before optimizations such as PEA.

While delegating the responsibility of interprocedural escape analysis to inlining is powerful, it introduces its own set of challenges. Inlining can be modeled as the NP-hard Knapsack problem if a *cost-benefit* assignment is available at each callsite [45]. However, the cost-benefit assignment is a best-effort approximation of a ground-truth cost-benefit – to correctly assess cost-benefit, the compiler must predict the optimizations that follow inlining, either by inferring them from the procedure body [4, 46], or by using inlining trials [13]. Profiling data, such as the callsite frequencies and

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call-target distributions, is then used to further assess the performance impact of each inlining decision [20, 21, 35].

Most prior work in the domain of inlining and interprocedural escape analysis indiscriminately applies ad-hoc heuristics to make inlining decisions whenever an allocation can be removed [5, 8, 25, 26, 33], and without ostensibly weighing these decisions against other considerations such as code size, compilation unit clustering, and instruction cost, which is what inliners traditionally do [2-4, 13, 35, 45, 46, 51]. This leads to inlining decisions that are based on a fragmented view of available information, leading to inconsistent performance metrics at best and adverse effects at worst.

This paper aims to address the following question -howcan interprocedural partial escape analysis improve inlining decisions, and consequently peak performance of user programs? The contributions in the paper are as follows:

- A novel approach that interleaves interprocedural partial escape analysis with inlining and computes the allocation frequency of escaping objects, with the aim of improving inlining decisions. The approach maintains flow-sensitivity across procedure boundaries and utilizes the collected profiling information (Section 3.1).
- Integration of materialization frequency metrics (Section 3.2) into an existing optimization-driven incremental inlining algorithm [35] used in GraalVM Native Image, yielding a single point of accountability for inlining (Section 3.3).
- A production-ready implementation in state-of-the-art optimizing compiler Graal [15, 16], in ahead-of-time compilation mode on GraalVM Native Image [57].
- A comprehensive performance evaluation that consists of peak-performance, allocation-rate, code-size and compile-time analysis on DaCapo [6], Scalabench [47], and Renaissance benchmark suites [39-41], as well as the extensive tuning of the proposed heuristics. We report a geometric mean performance improvement of 1.79% up to 24.62% across all 36 benchmarks, and geomean code size improvement of 1.81%. In total there are 6 benchmarks with >5% speedup and geometric mean of 9.10% (Section 4).

2 Motivation & Background

Consider the method put of the IntListMap class in Listing 1, which closely corresponds to the class ListHashMap from the Groovy core library [19]. This collection stores key-value pairs as two consecutive entries in an array-

```
class IntListMap<V> {
 1
   Object[] list:
2
    int last:
   V put(int key, V val) {
 5
     var boxed = new Integer(key);
     for (int i=0;i<last;i+=2)</pre>
      if (list[i].equals(boxed))
8
9
       return list[i + 1] = val;
10
     list[last++] = boxed:
     return list[last++] = val;
11
12
   }
13 }
```

Listing 1. IntListMap

based list. Keys are always primitive integers, but values are

of any generic object type V. The list array is of Object[] type because JVM cannot store primitives and objects in the same array. Thus, a key must be boxed in line 6 into an object, before being stored. In the loop in line 7, put then scans the array up to position last, to see if the specified key is already in list. If an equal key is present at i, the existing value at i + 1 is overwritten, but the key itself is not stored. If the key is not present, the boxed key is stored in line 10.

It seems as if the allocation can be delayed until line 10 if a map kept reusing the same keys, this would reduce the allocation rate greatly. But alas - boxed escapes earlier, in the equals call in line 8! Luckily, any decent inliner will exploit the receiver-type profile to decide that equals always calls Integer#equals [14] – it inserts a speculative typecheck, and trivially inlines Integer#equals, which compares the primitive integer fields of the two boxed Integer objects: if (!(list[i] instanceof Integer)) deopt();

if (list[i].integer==boxed.integer) return list[i + 1] = val; If the dynamic typecheck fails, the code is either deoptimized, or the execution gets continued in baseline compiled code, which does not inline equals, but is less efficient [56].

1	for (i = 0; i < last; i += 2)					
2	<pre>if (list[i].integer == key)</pre>					
3	return list[i + 1] = val;					
4	<pre>var boxed1 = new Integer(key);</pre>					
5	list[last++] = boxed1;					
6	<pre>return list[last++] = val;</pre>					
	Listing 2. PEA on put					

PEA can now postpone the boxed allocation – it keeps it virtual in line 2 in Listing 2, where the read boxed.integer can be replaced with just key; and materializes it in line 4. Yet,

a more realistic list-map, one that better resembles Groovy's ListHashMap [19], would keep its key type generic.

class ListMap<K, V> 1

2

4

5

6

In this case, the key parameter has a type K, which is erased to an Object type [10] - a primitive integer key must be boxed at put's callsite, be-

```
return list[i + 1] = val;
    list[last++] = key;
    return list[last++] = val;
7 }
    Listing 3. Generic put
```

for (int i=0;i<last;i+=2)
if (list[i].equals(key))</pre>

V put(K key, V val) {

fore put is even called, so there is nothing PEA can do! The only hope is that the inliner decides to inline put wherever put is called - then PEA could 'see' the allocation of key. The following procedure inserts the primitive key 1 into a ListMap that maps Integer keys to String values:

```
void client(ListMap<Integer, String> m)
  { return m.put(1, "one"); }
```

Since the key-type is Integer, the javac compiler 'autoboxes' the primitive int value that is passed in the put call. The actual compiled code resembles the following:

return m.put(new Integer(1), "one");

If something were to 'convince' the inliner that there is benefit in inlining the put call, the resulting compilation unit would contain both the allocation and its usages. This extended horizon over the program would allow PEA to transform the code to the following:

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```
for (i = 0; i < m.last; i += 2)
    if (m.list[i].integer == 1) return m.list[i + 1] = "one";
var boxed1 = new Integer(1);
m.list[m.last++] = boxed1;
return m.list[m.last++] = "one";</pre>
```

But, despite having delayed the allocation of the Integer, was inlining put really worth it? It depends. For one thing, the put call inside the client procedure should be overall executed often, otherwise, the gains are anyway miniscule. Then, we should be reasonably assured that path to the final return is cold, that is, that most of the time we exit early, in the for loop. Finally, we should make sure that there are no other 'better' inlining candidates – perhaps the inlining budget would be better spent elsewhere?

To deal with such considerations, inliners build cost-benefit models that guide their decisions. These models are usually based on execution profiles, knowledge about code-size and opportunities for other optimizations. Rather than disregarding all these other factors, perhaps we should try to influence an existing inliner by biasing its cost-benefit model?

This deliberation motivates our main research questions:

- **RQ1:** How do we compute the frequency of materializations for objects that escape across procedures?
- **RQ2:** How can we quantify the benefit of reducing materialization-frequency through inlining, such that we improve an existing cost-benefit estimation?

2.1 Intraprocedural Partial Escape Analysis

Let's consider how standard intraprocedural PEA works when viewed at the source code level of the previous example. At each instruction *n*, PEA maintains a mapping O(n) called *virtualization*, which maps each value to the virtualization state of that value. Each allocated object is mapped to either a materialized value $m \in \mathbb{V}$, indicating that the respective object escaped and was allocated as another value *m*; or to a sequence of field-value pairs $f_j \rightarrow x_j$, indicating the state of an object that did not escape (i.e. object that is virtual):

 $O(n) : \mathbb{V} \mapsto \mathbb{V} \cup \{(f_0 \to x_0, \dots, f_i \to x_i) : f_j \in \mathbb{S} \land x_j \in \mathbb{V}\}$ Above, \mathbb{V} represents the set of referable values in the program, f_j represents a field name from the set of all strings \mathbb{S} , and each field is mapped to another value $x_j \in \mathbb{V}$.

In our example with the put(int, V) procedure, after executing the statement $n_0 \equiv var boxed=new Integer(key)$ in line 6, O maps the boxed value to a tuple whose integer field is set to the value of the parameter key:

 $O(n_0) = \{ boxed \mapsto (integer \rightarrow key) \}$

Statements in line 7 and line 8 do not modify the boxed object, in other words $O(n_1) = O(n_0) = O(n_2)$. However, n_2 can exploit the mapping O, namely, n_2 can be replaced with list[i].integer==key, because $O(n_2)$ tells us that boxed is virtual (i.e. did not yet escape), and its value field is equal to key. The next statement $n_3 \equiv list[last++]=boxed$ materializes the boxed object, since the reference to boxed escapes

as part of the array-write. Mapping $O(n_3)$ becomes:

$$O(n_3) = \{ boxed \mapsto boxed1 \}$$

which means that original allocation boxed is on this controlflow path materialized under the name boxed1. PEA thus inserts the instruction var boxed1=new Integer(key) just before n_3 , and makes sure that the boxed1 value is used instead of boxed on the rest of the control-flow path.

PEA on control-flow graphs. Having grasped how PEA uses the mapping O, let's express PEA as a dataflow analysis on control-flow graphs (CFGs). Control-flow graphs will allow us to more concisely define how individual instructions affect the mapping O. Also, our actual implementation was done on Graal IR, which is conceptually a reducible control-flow graph. The CFG that corresponds to the put procedure from Listing 1 is shown in Figure 1.

In the previous discussion of Listing 1, we were somewhat lax about what an *instruction* is – in the CFG, we define an individual instruction as a node. Nodes may be connected with directed edges, which represent the control-flow of the program. For each type of node, we can define two mappings O_{in} and O_{out} , as the virtualization before and after the respective node is executed, and use them to describe how each node changes the virtualization.

Let the set \mathbb{A} be the set of all nodes that represent object allocations. Then, node $a \in \mathbb{A}$, which allocates an object with *i* fields, modifies *O* as follows (\perp is the default value):

$$O_{out}(a) = O_{in}(a)[a \mapsto (f_0 \to \bot, \dots, f_i \to \bot)] \quad a \in \mathbb{A}$$

Next, let \mathbb{W} be the set of nodes that store a value x into a field f of an object allocated at a. For a node $w_{f,x}^a \in \mathbb{W}$ we have two cases – the field f is updated to x only if a is virtual, and otherwise, the mapping is not modified:

$$O_{out}(w_{f,x}^{a}) = \begin{cases} O_{in}(w_{f,x}^{a})[a \mapsto (\dots, f \to x, \dots)] \\ if a \mapsto (\dots, f \to y, \dots) \in O_{in}(w_{f,x}^{a}) \\ O_{in}(w_{f,x}^{a}) & otherwise when w_{f,x}^{a} \in \mathbb{W} \end{cases}$$

Set \mathbb{E} consists of nodes that cause an object to escape to the heap: in intraprocedural PEA, these are calls to other subroutines, writes to global fields, and writes to already materialized objects. We use the relation $n \xrightarrow{\text{esc}} a$ to denote that a node *n* causes an object allocated at node *a* to escape.

Let $e^a \xrightarrow{\text{esc}} a$. Again, *O* changes only if object *a* is virtual, in which case *a* is materialized. For all fields *f*, if *f* points to another virtual object, that object is transitively materialized:

$$\mu(O, x) = \begin{cases} \mu(\dots \mu(O[x \mapsto m], x_0), \dots, x_i) \\ if x \mapsto (f_0 \to x_0, \dots, f_i \to x_i) \in O \\ O \quad otherwise \end{cases}$$

$$O_{out}(e^a) = \mu(O_{in}(e^a), a) \quad \text{when } e^a \xrightarrow{\text{esc}} a$$

Above, the helper function μ recursively traverses the object graph. Each virtual object is replaced with *m*, which represents a newly inserted node that *materializes* the object that had been allocated at node *a* in the original program.



Figure 1. Simplified control flow(\rightarrow) and data flow (\rightarrow) of put. The boxed allocation is moved to latest point before boxed becomes materialized and loads are replaced with scalars from the virtual state.

One may now ask: could the escaping node e^a be used to represent the materialization too, i.e. is $m = e^a$? This is not generally the case – to see why, consider the *merge* node, which has multiple predecessors. A merge node *j* can implicitly trigger an escape in order to enforce a consistent state in its successor node. First, a value only survives the merge if it exists in all predecessor states. Then, if an object is virtual in all predecessors, it can remain virtual after the merge. But, if the object is materialized in at least one predecessor, then it must be materialized in the others too (for simplicity, this is what most PEA implementations do). Formally:

$$\iota(x,y) = \begin{cases} (f_1 \to \iota(x_1,y_1), \dots, f_n \to \iota(x_n,y_n)) \\ if x = (f_1 \to x_1, \dots, f_n \to x_n) \\ \land y = (f_1 \to y_1, \dots, f_n \to y_n) \\ \varphi(m_1,m_2) \quad otherwise \end{cases}$$
$$O_1 \sqcap O_2 = [a \mapsto \iota(x,y) | \forall a.a \mapsto x \in O_1 \land a \mapsto y \in O_2] \\ O_{out}(j) = O_{in,1}(j) \sqcap \dots \sqcap O_{in,p}(j) \end{cases}$$

Above, for brevity, φ is a *Phi function* from static singleassignment (SSA) form, which takes a different value depending on the predecessor by which the merge was reached – it can be trivially translated to a CFG with variables [43].

In the definition of ι , if the allocation was materialized as m_1 on one predecessor, and virtual on the second predecessor, then it must be materialized to a node m_2 on the second predecessor too. Observe the CFG with a merge node. The allocation *a* escapes in the left branch at the store node, but not on the right branch. Here, *only* store $\stackrel{\text{esc.}}{\xrightarrow{}}$ *a*, but *both*

store $\xrightarrow{\text{mat}} a \text{ and } \text{end}_2 \xrightarrow{\text{mat}} a$, so allocations are inserted in both branches.

We can now inspect how the virtualization O changes at the edges between nodes in the CFG of the put procedure in Figure 1, The object behind the boxed allocation is virtual throughout the loop, and allows simplifying the check inside the loop. The store node causes the object to escape after the loop exit ($\frac{esc}{c}$), and the newly inserted



red/bold allocation node is that allocation's materialization node (\underline{mat}) .

3 Algorithm Description

The main idea behind our Interprocedural Partial Escape Analysis (IPEA) algorithm is to predict the materialization frequency after inlining, and use it to adjust the inliningbenefit approximation of a callee. The frequency of materializations determines the number of dynamic allocations at runtime. We first discuss how IPEA locates materializations across the entire call tree instead of inside a single procedure.



Figure 2. Call tree representation of put and client code and key pointing to its interprocedural materialization.

3.1 Interprocedural Partial Escape Analysis

Interprocedural analysis accounts for the relationships between procedures, so we shift our perspective from an individual procedure body to a set of procedure bodies logically assembled in *call tree* representation. Revisiting our generic ListMap#put procedure from Section 2, Figure 2 now shows the call tree of client, put and equals and their respective simplified Intermediate Representations (IR). Recall that PEA could postpone the materialization of the key parameter of put *after* the loop and before it is written to the array. In order to find this materialization location, the virtualization states must be shared across procedure boundaries. We call this *virtual state injection*, or *virtual injection*.

In the example in Figure 2, the parameter key of the put procedure originates from the allocation in the client code. To simulate the effects of inlining, we can intuitively treat the parameter node as an extension of the allocation in the caller. If key is still virtual at the callsite, we can replace the parameter with a virtual object originating from the allocation in the client. From this point on, when a materialization is triggered for the replaced parameter node, we know that this is exactly where PEA will insert the allocation after inlining.

Similarly, when the entire procedure body is completely processed, the callsite must reacquire the virtual state from its callee, i.e. the virtual state must be re-injected back into the caller. Let's assume the equals procedure has yet to be inlined – the key object is also passed to the equals procedure inside put. If the equals procedure materializes its parameter, there is no benefit in inlining put. Our analysis must therefore first process the entire equals procedure, and re-inject the virtual state into put to determine that key would stay virtual until key is materialized by the store node.

Having provided an intuition of IPEA on the client code, we now detail how the intraprocedural analysis is extended to allow the prediction of materializations after inlining. To properly model O across procedure boundaries, invoke nodes no longer cause escaping – instead, invokes trigger a recursive analysis in the subtree of the call tree. All other nodes that trigger escaping are treated as in ordinary PEA.

We also add an additional state component on top of the local analysis state O_n , namely the shared global state *materializations* \mathcal{M} , which maps each allocation in the call tree to the set of its materializations in the call tree. Our new compositional state \mathcal{S} is now defined as follows:

$$S(n) = \langle O(n), \mathcal{M} \rangle$$
$$\mathcal{M} : \mathbb{A} \to \{ n \mid n \xrightarrow{\text{mat}} a \land a \in \mathbb{A} \}$$

where O(n) is a per-node state, and \mathcal{M} is a shared global state across the entire call tree. \mathcal{M} holds the desired analysis results, and is later used to drive inlining decisions. When materializations are determined, i.e. after IPEA finishes processing a procedure (a call-tree node), for any node *n* such that $n \xrightarrow{\text{mat}} a$, we update \mathcal{M} accordingly:

$$\mathcal{M}[a] = \mathcal{M}[a] \cup \{n\}$$

To formalize the notion of the call tree, we rely on the relation $c_{a_0,\dots,a_j} \xrightarrow{\text{calls}} p_0,\dots,p_j$, which denotes that an invoke node c_{a_0,\dots,a_j} with arguments a_0,\dots,a_j calls the callee procedure whose parameters are nodes p_0 to p_j .

State-update rules. IPEA processes specific nodes, namely the invoke, parameter and return nodes, in a different way. When we encounter an invoke node $c_{a_0,...,a_j}$ with arguments $a_0, ..., a_j$ such that $c_{a_0,...,a_j} \xrightarrow{\text{calls}} p_0, ..., p_j$, we define the update rules in the callee as follows:

$$O_{out}(p_i) = \begin{cases} O_{in}(p_i)[p_i \mapsto (f_0 \to x_0, \dots, f_k \to x_k)] \\ if a_i \mapsto (f_0 \to x_0, \dots, f_k \to x_k) \in O_{in}(c_{a_0,\dots,a_j}) \\ O_{in}(p_i) & otherwise when c_{a_0,\dots,a_j} \xrightarrow{\text{calls}} p_0, \dots, p_j \end{cases}$$

The above states that the initial state of a formal parameter is virtualized if the corresponding argument in the caller remained virtual, and materialized otherwise.

The following illustrates the rule on the put procedure:



At the callsite, the first argument is mapped to the new boxed node, so the corresponding formal parameter in the callee is mapped accordingly. After the array-write is processed, the materialization map will contain a single mapping $\mathcal{M} \equiv \{\text{boxed} \mapsto \{\text{store list}\}\}$, as indicated in Figure 2.

Let the set \mathbb{R} of *return nodes* consist of nodes r_x , such that the return node r_x returns the value of node x. Handling a return node r_x serves two purposes. The first is to re-inject the state of the callsite arguments back into the caller. If a formal parameter gets materialized, the callsite argument must also be materialized. Second, if a virtual object is returned from the callee, the invoke node can be mapped to a virtual object in the caller thereafter.

We use a helper function π to capture how the virtualization state of the parameter p_i affects the callee state O:

$$\pi(O, p_i) = O[a_i \mapsto v] \quad such \ that \ \exists c_{a_0, \dots, a_j} \xrightarrow{\text{calls}} p_0, \dots, p_j \land$$
$$p_i \mapsto v \in O_{out}(r_x) \land r_x \in \mathbb{R}$$
The above says that for any parameter, there must be a return

node that mentions that parameter's state, and that the corresponding callsite argument must be mapped accordingly.

We use another helper ρ to capture how the return value alters the virtualization state of the callee:

$$\begin{split} \varrho(O, c_{a_0, \dots, a_j}) &= O[c_{a_0, \dots, a_j} \mapsto v] \quad c_{a_0, \dots, a_j} \xrightarrow{\text{calls}} p_0, \dots, p_j \land \\ p_0, \dots, p_j &\in O_{out}(r_x) \land x \mapsto v \in O_{out}(r_x) \land r_x \in \mathbb{R} \end{split}$$

The above says that for any invoke $c_{a_0,...,a_j}$, there must be a return node r_x in the corresponding callee whose returned value x is mapped to some state v – the virtualization state of the invoke $c_{a_0,...,a_j}$ must be remapped to that state v.

The rule for invokes is now concisely expressed as:

$$O_{out}(c_{a_0,\dots,a_j}) = \varrho(\pi(\dots\pi(O_{in}(c_{a_0,\dots,a_j}), p_0), \dots, p_j), c_{a_0,\dots,a_j})$$

The following figure illustrates the previous rule – the state of the object a changes in the callee foo (but is not materialized), so the analysis continues with the new virtual state of a. Moreover, the return object x is created inside the scope of the callee, and remains virtual in the callee.



In the preceding rule, we implicitly assumed that there is a single return node r_x in any callee. Indeed, the functions π and ϱ are not defined when there are multiple return nodes. To address this, we simply transform the CFG of each callee

by replacing all the return nodes r_{x_i} with a merge node j followed by a single return node $r_{\varphi(x_0,\dots,x_k)}$ (which, in fact, reflects what happens during inlining). This way, the virtualized state persists only if all return nodes return a virtualized object, as per the rule for merge nodes from Section 2.1.

Pseudocode. The code for the IPEA algorithm, which produces the map of materializations \mathcal{M} for a call tree \mathcal{T} , is shown in Algorithm 1. The call runIPEA(root(\mathcal{T}), {}) starts the algorithm. It traverses the CFG in a reverse postorder, using the extended set of rules for updating the virtualization state. If it encounters an invoke, it first calls runIPEA on the child node of the call tree before updating the state in the caller. In each call tree node, CFG traversal must be repeated until reaching a fix point [50]. Every time a call tree node reaches a fix point, the materialization map \mathcal{M} is updated.

Algorithm 1: Interprocedural PEA					
input : Call tree \mathcal{T}					
output : Materializations $\mathcal{M} : a \to \{n \mid n \xrightarrow{\text{mat}} a \land a \in \mathbb{A}\}$					
Procedure $runIPEA(CFG, M)$ is					
$S_{in}(start) \leftarrow \langle \{\}, \mathcal{M} \rangle$					
<pre>while not fixedPointReached() do</pre>					
for node in reversePostOrder(CFG) do					
if node is Invoke then					
runIPEA($CFG(node), \mathcal{M}$)					
$S_{out}(n) \leftarrow process(node, S_{in}(node))$					
for (n,a) such that $n \xrightarrow{mat} a$ do					
$\mathcal{M}[a] \leftarrow \mathcal{M}[a] \cup n$					

3.2 Materialization-Frequency Reduction

In our running example from Section 2, the reduction of key allocations depends on the workload in the client code – if the client only occasionally inserts a previously unseen key, then the number of allocations is reduced considerably. To reflect this in the inlining benefit, we use the materialization map \mathcal{M} and node-frequency profiles to compute the *materialization-frequency reduction*:

$$\mathcal{FR}(a) = \frac{F(a) - \sum_{m}^{m \xrightarrow{\text{mat}} a} F(m)}{F(a)} \in [0, 1] \quad (1)$$

Above, F(n) is the execution frequency of a node relative to the entry to the root compilation unit in the call tree. Note that the frequencies of materialization nodes are summed – correctness is given by guarantees a PEA invariant: For any two nodes n_1 and n_2 that materialize the same allocation node a, there is no execution path that passes both through n_1 and n_2 . In other words, for a single a, every execution path has at most a single node n, such that $n \xrightarrow{\text{mat}} a$.

3.3 Integrating IPEA into the Incremental Inliner

We extended an existing implementation of the incremental inliner [35] with our analysis. We now present a simplified description of the existing inliner, and show how it was enhanced. The incremental inliner is representative of typical inlining algorithms, and we believe that our analysis can be easily applied to similar inliners.

Each time a method is submitted for compilation, that method becomes the root compilation unit of its respective call tree. The inliner incrementally *expands* the call tree during an *expansion phase*. Expanding means picking a leaf node in the call tree, and attaching the callees of that node as children to the call tree. A call-tree node is initially unexplored, and called a *cutoff node*. After the expansion phase, the inliner assigns cost-benefit to each call-tree node. Based on the cost-benefit, some call-tree nodes are inlined during the *inlining phase*, and the inliner subsequently repeats these stages. This cycle continues until reaching a budget limit [35].

IPEA as described in Algorithm 1 is inserted into the inliner after expansion and cost-benefit assignments, but before inlining (see Algorithm 2). We note that each invoke that corresponds to a cutoff node is considered an analysis boundary, and is treated as a regular escape.

Algorithm 2: Integration of IPEA						
input : Root compilation unit <i>root</i>						
<pre>Procedure incrementalInliner(root) is</pre>						
<pre>while continueInlining() do</pre>						
expansion()						
$\mathcal{M} \leftarrow runIPEA(root)$						
<pre>computeBenefits(root)</pre>						
for $a \in AllocationNode$ do						
$\int fr \leftarrow \mathcal{FR}(a)$						
boostSubtree(fr)						
inlining()						

Consider the compilation of the client procedure from Section 2. Its inlining schedule will typically be as follows:

- (i) Build initial call tree for client, add put as a cutoff.
- (ii) Expand put, add equals as a cutoff node, expand equals.
- (iii) Run IPEA on the call tree (client, put, and equals).
- (iv) Compute cost-benefit of put and equals.
- (v) Boost both put and equals based on the reduction in materialization frequency of key.
- (vi) Decide to inline both put and equals.

Benefit. Before inlining, each node in the call tree is assigned a benefit. This benefit is a best-effort estimate of the performance gains from inlining, and it guides the inlining decisions. Our analysis adjusts the benefit assignment by incorporating materialization-frequency reduction. Our new benefit function is the following:

$$\mathcal{B}(c_{a_0,\dots,a_j}) = \mathcal{B}'(c_{a_0,\dots,a_j}) \cdot \max_{i \in 0,\dots,j} (\delta_{mb})^{\mathcal{FR}(a_i)}$$

where $\mathcal{B}'(c_{a_0,...,a_j})$ represents the previous benefit function for a node at callsite $c_{a_0,...,a_j}$, and δ_{mb} is the *materialization boost*, a tunable hyperparameter that quantifies how much the inliner should value the reduction in materialization frequency. A call tree node is boosted relative to the frequency reduction $\mathcal{FR}(a)$ if *a* escapes into the invoke $c_{a_0,...,a_j}$. For example, in the earlier client method, the boost $\mathcal{FR}(key)$ is applied to both equals and put, since key escapes to both, and inlining both of them is required to reduce allocations. **Expansion phase.** In Algorithm 2, we applied runIPEA to the point after the expansion phase. However, since each expansion phase expands multiple cutoff nodes of the call tree, we can also incrementally run IPEA on the subtree affected by the expansion. The advantage of doing so is that we can boost the expansion priority of cutoffs at which objects escape – each cutoff node has an associated expansion priority, which decides how soon that node will be expanded.

During IPEA, we track an additional state component \mathcal{E} for allocations that escape into cutoff nodes:

$$\mathcal{E}: \{c_{a_0,\dots,a_i} \text{ is a cutoff}\} \to \{a_i | c_{a_0,\dots,a_i} \xrightarrow{\text{esc}} a_i\}$$

The cutoff-node priorities $\mathcal P$ are then continuously updated according to the following equation:

$$\mathcal{P}(n) = \mathcal{P}'(n) \cdot \delta_{eb} \cdot |\mathcal{E}(n)|$$

where the new priority \mathcal{P} boosts the vanilla (cost-benefitbased [35]) priority \mathcal{P}' by the number of escaping objects. This equation also introduces an additional hyperparameter δ_{eb} , or *escape boost*, which determines the expansion priority given to virtual objects that are escaping.

Cutoff weight. In an idealized analysis, the entire call tree is available, and no invoke causes an object to escape. However, in practice, expansion budget is often insufficient, and the call tree is usually infinite due to recursion, meaning that at least some cutoff nodes remain in the call tree. Whenever an object escapes at a cutoff node, this can be for two reasons: either the object would really be materialized after expansion, or the object is materialized only because we did not expand the callee yet. To prevent over-penalizing such materializations, we found it useful to optimistically assume that an object that escapes at a cutoff may survive. We, therefore, updated the frequency-reduction expression from Equation (1) with the *cutoff weight* hyperparameter $\delta_{cw} \in [0, 1]$:

$$\mathcal{FR}(a) = \frac{F(a) - \sum_{m} \phi_{m} \cdot F(m)}{F(a)} \quad \phi_{m} = \begin{cases} \delta_{cw} \ m \ is \ cutoff \\ 1 \ otherwise \end{cases}$$

We discuss the tuning of this heuristic in Section 4.

4 Evaluation

The goal of the evaluation is to show improvements over flow-insensitive escape analysis techniques in terms of runtime, code-size and compilation times, examine the profiled dynamic number of allocations, and the effect of heuristics.

We compare our Interprocedural PEA (IPEA) to a Base, which runs only the optimization-driven inliner [35], that is run in every configuration. Our second comparison is Partial Escape Selective Inlining (PESI), the control flow insensitive inlining technique implemented in GraalVM.

Decreasing the runtime is the principal goal of our analysis phase. We emphasize the GraalVM compiler has been already heavily optimized and most performance-critical methods are inlined. Nevertheless, we consider improvements of > 5% to be significant and valuable incremental contributions to GraalVM. While secondary to runtime performance, keeping code size low is meaningful for Native Image, especially for one of the main use cases of cloud deployment [49, 53]. Build time is an important measurement, as PEA is an expensive analysis for large procedures, even intraprocedurally.

Methodology. All reported results were run on an Intel(R) Xeon(R) E5-2699 v3 @2.30 GHz with 72 cores and 256GB of RAM. The benchmarks were bound to 36 CPUs on the same NUMA node and the maximum Heapsize was capped at 8GB. All benchmarks were run on JDK 17 except for the Renaissance-spark jobs.¹ Relative improvements between T_{old} and T_{new} are always computed as $\frac{T_{old}-T_{new}}{T_{old}}$ unless stated otherwise and averaged improvements use geometric mean.

Benchsuite	Version	JI
Renaissance [40]	0.14.1	1
Renaissance-spark	0.11.0	1
Dacapo [6]	9.12 mr1	1
ScalaBench [47]	0.1.0	1

Runtimes. Plot in Figure 3 CK compares relative runtime performances across all benchmarks (lower is better). We include an additional column for IPEA OPT, where the best hy-

perparameters are chosen for each individual benchmark, instead of optimizing for the entire suite. We observe that some types of benchmarks react strongly to IPEA, where benchmarks improve over PESI, like scalap - 5.5%, scalac -5.74%, finagle-chirper -7.74% avrora -16.07% and mnemonics -24.62%. In practice, some benchmarks show no performance difference when optimizing for allocation reduction. This may be the case for more CPU-bound applications, or code sections that perform allocations are hot enough to be inlined based on frequency and code size metrics. Further, we see that individual benchmarks have more optimization opportunities: Benchmarks like kiama, gauss-mix, par-mnemonics and fop leave major improvements of up to 16% between IPEA and IPEA OPT unrealized. This suggests future work on this topic is of relevance. Some benchmarks exhibit regressions, factorie by +3.35%, db-shootout +5.07% and most notably xalan +16.07%. For both db-shootout and factorie, these regressions can be eliminated by choosing better parameters. We discuss xalan in more depth in the heuristic tuning section.

Code size. See Figure 4 for codesize and build time results. IPEA decreases code size almost across the board compared to PESI, with the only exceptions being scalariform, scalaxb and scala-doku. On average our code size has decreased by 1.81%. This suggests our analysis has increased inlining precision, were overall less, but more impactful procedures are inlined ².

¹Renaissance-spark benchmarks did not run on JDK 17 at time of experimentation:page-rank, movie-lens,db-shootout, neo4j-analytics, gauss-mix. rx-scrabble, dotty, naive-bayes, als.

²We stress that more inlining does not always cause larger code.

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Figure 3. Relative runtime results (lower is better) in comparison to BASE of PESI, IPEA, and IPEA OPT where hyperparameters are chosen individually for each benchmark to achieve the lowest *runtime*. Some benchmarks like mnemonics, finagle-chirper, and page-rank reach strongly to allocation reductions, and IPEA outperforms PESI on 25 of the 36 benchmarks and significantly outperforms with an improvement of >5% on 6 benchmarks, while only having one significant regression for xalan.



Figure 4. Build times and Code size (lower is better) of all configurations relative to BASE, where 31 of the 36 benchmarks have a lower codesize with geomean 1.81% and up to 4.98% comparing IPEA to PESI. Similarly, 27 benchmarks have lower build times with geomean improvement of 1.62% up to 8.2% faster over PESI.

Build time. Build time is kept low and shows an improvement over PESI by a geomean of 1.62% Mostly this is a result of not running IPEA for every inlining round and further throttling IPEA if procedures become increasingly large. The largest build time increases are observed for benchmarks tmt with +7.7% and, scalariform with +5.3%.

We differentiate two factors 📋 Inliner time 📘 PESI time for differences in build time. First, the accumulated time spent in performing the analysis, shown in the figure on the right, and second the effects of inlining making other optimizations slower due to larger procedure bodies . Notice that IPEA is part of the incremental inliner, causing more expansion of the call tree and more inlining rounds, in addition to the time spent within IPEA itself. Nevertheless, considering



both factors results in lower overall average build time.

Dynamic allocations at runtime. For more insight on the performance improvements and the effects of our analysis, we profiled the allocation sites for dynamic memory allocated at runtime, see Table 1. We stress that the number of bytes allocated is merely a fraction of the effects on runtime. Take the benchmark scalap. We measured higher runtimes of 5% with PESI due to code bloat by over-inlining to eliminate allocations, and therefore a higher number of instruction cache misses. We now have faster runtime, even though more bytes are allocated at runtime. Clearly, IPEA has taken other metrics into account before inlining to eliminate allocations. Concurrently, benchmarks like mnemonics and finagle-chirper seem to directly benefit from the reduction in allocations with 24.62% and 7.74% faster runtimes. However, we still see inconsistent results in some scenarios, where factorie has a 15% reduction in allocations, yet is 3.35% slower - allocation reduction is overvalued . We argue these results show that it is less about the number of allocations (allocations are cheap) and more about how these allocations interact with other elements of the code and feel our approach is a step in the right direction and can be further refined. While not displayed in this data, our anecdotal experience in analyzing the allocation sites has shown that where the allocations occur with PESI vs IPEA are quite different, even if the amount of memory allocated is similar and is worth further investigating.

Heuristic Tuning. Recall the introduction of three primary hyper-parameters that we tune for: materialization boost δ_{mb} , escape boost δ_{eb} and cutoff weight δ_{cw} . The parameters were tuned across all benchmarks in all suites. For brevity's sake, we show only a handful in Figure 5. Strong performance reactions are detected for gauss-mix and kiama where we

Table 1. Selected runtime allocation profiling results. Allocation reduction for all benchmarks can be found here [54].

GB allocated $\frac{IPEA-PESI}{PESI}$					
Benchmar	kBase PESI IPEA	A AAlloc	ΔRuntim	e	
avrora	2.940 1.486 2.75	9 +85.6%	-16.07%	_	
fop	7.922 7.374 7.40	5 +0.41%	-2.66%	od	
pmd	125.3 120.4 114.	0 -5.34%	-2.23%	ICa	
xalan	72.59 72.48 70.8	1 -2.3%	+16.17%	D_{a}	
geomean		+6.5%	-1.33%		
akka-uct	803.1 795.9 797.	3 +0.17%	-0.23%		
fin-chir.	1584 1527 1467	-3.91%	-7.74%	e	
fj-kmea.	576.4 576.2 577.	2 +0.19%	-2.4%	nc	
fut-gen.	75.34 76.19 114.	2 +50.0%	+1.74%	ssa	
gauss.	110.2 108.1 108.	4 +0.3%	+0.16%	ais	
mnem.	143.7 156.5 133.	4 -14.73%	-24.62%	en	
mo-lens	324.5 323.2 319.	1 -1.26%	-2.55%	Я	
p-rank	294.6 282.4 266.	3 -5.68%	+6.13%		
geomean		-1.9%	-2.07%		
apparat	22.95 22.41 22.3	9 -0.09%	-4.14%		
factorie	116.3 106.7 90.6	9 -15.05%	+3.35%	В.	
scalac	20.98 20.75 19.7	7 -4.73%	-5.74%	ala	
scalap	6.664 6.493 6.57	3 +1.23%	-5.78%	Sci	
geomean		+2.04%	-1.61%		
total		+1.02%	-1.79%		



Figure 5. Selected grid search of hyper parameters, lower is better. x-labels: ($\delta_{eb} : \delta_{mb} : \delta_{cw}$). Grid search plots for entire benchmark suite can be found here [54].

were not able to extract the complete potential when optimizing for the entire suite. Our only major regression xalan experiences a slowdown even with a reduction in allocations. It is the only benchmark that experiences slowdowns with any parameters in our grid search. We remark that single-threaded performance of xalan with IPEA performs the same as PESI. Our running hypothesis is that the different inlining choices improve sequential code performance, but decrease the periods between inter-thread synchronizations,

which causes more contention. In practice, $\delta_{eb} = 6$, $\delta_{mb} = 20$ and $\delta_{cw} = 0$ performed the best on average. Interestingly, assigning a materialization caused by an unexplored invoke to weight 0 has the best performance. Intuitively, this helps in finding optimization opportunities for objects that get passed deep into the call graph. If an inlining decision is made based on this value, the reduction in materialization frequency is not guaranteed but the materialization is certainly delayed.

5 Related Work

It can be argued that inlining and escape analysis are the cornerstone of high-level compiler optimizations, and crucial for high-level programs and frameworks that involve (parallel) data-processing [30, 34, 38, 58], streaming [24, 29, 37], and functional programming patterns in particular [22, 28, 31, 52]. In this section, we give a brief overview of related literature ranging from escape analysis to inlining with cost-benefit estimations, and the relationship between the two.

Cost-benefit based inlining. Some of the earliest work on cost-benefit based inlining with a fixed budget is introduced by Ayers et al. [3] and shortly after by Arnold et al. [2]. While their cost model is conceptually similar to modern approaches, the benefit model used is straightforward - Avers et al. infer the benefit based on the ability of the callee to be specialized. They use a similar approach to Ball [4] that uses data-flow analysis on the procedure arguments to predict the effects of inlining when propagating a constant to the callee. Arnold et al. use the profiled callsite frequency as the benefit metric to estimate the direct benefits of inlining by removing linkage overhead. Dean & Chambers [13] showed that the indirect benefits are equally, if not more, important than the direct benefits and perform inlining trials. The resulting optimized body is analyzed to predict benefits. With the injection of virtual state across procedure boundaries, IPEA can be seen as a sophisticated inlining trial.

Escape analysis. Choi et al. [11], Gay & Steensgaard [18], Kotzmann et al. [25], Blanchet [8], and Whaley & Rimard. [55] all introduce a form of interprocedural escape analysis. These techniques inject states from a caller into the callee and vice versa. Both Blanchet and Kotzmann et al. suggest inlining based on the analysis results [8, 25]. Gay & Steensgaard, Whaley & Rimard, and Choi et al. use the acquired information only for stack allocation and lock elimination [11, 18, 55]. Choi et al. continue to describe a flow-sensitive interprocedural escape analysis for higher analysis accuracy, but the analysis is still only used for making binary escapability assessments [11]. Attempting to predict the frequency of dynamic allocations statically is never considered.

Profiling allocation sites. Shankar et al. [5] introduce JOLT, a profiling and inlining algorithm that targets the reduction of object churn with capture and control analysis introduced by Dufour et al. [17]. JOLT profiles object lifetimes dynamically and directly from the garbage collector. The data is

used in a standalone cost-benefit inlining phase. Our evaluation showed multiple separated inlining phases can produce inconsistent results. However, combined dynamic object profiling and materialization frequency could further amplify the accuracy of the cost-benefit model.

Inlining meets Escape Analysis. Dealing with dynamic procedure calls can be challenging for program analysis [23]. Devirtualizing call sites allow further exploration of the call tree. Sewe et al. study the prediction of inlining effects on the specialization of the call tree, which in turn enables more inlining [46]. This specialization of the call tree is analogous to delaying materializations since the type of a virtual object is always known. This fact is also used by the optimization-driven incremental inliner mentioned throughout this paper by Prokopec et al. [35, 42], which runs PEA only intraprocedurally in the root procedure in between inlining rounds. This PEA runs completely disjoint from IPEA and aims to devirtualize invokes on virtual objects, and propagate type information to the callee, if a parameter is virtual at a callsite.

In a similar vein, Interflow by Shabalin et al. [48] fuses multiple techniques, including PEA, code duplication, and inlining to optimize collection-based Scala code for Scala Native. They report performing PEA in the same pass as inlining and try to always inline procedures with virtual objects at the callsite to delay the materialization. We improve on this shared concept by additionally computing materialization frequency, yet this again prompts the question – can we quantify the benefit of delaying a materialization and use value in the cost-benefit function for inlining prediction? Can we not only estimate the direct effects of inlining on allocation reduction but also encapsulate the effects of other intraprocedural optimization phases that aim to reduce runtime allocations (like code duplication)?

6 Conclusion

We presented a novel algorithm based on interprocedural partial escape analysis, which computes object-materialization frequencies and predicts the benefits of inlining. We proposed several heuristics that rely on allocation-frequency reduction to guide inlining decisions. We tuned these heuristics across 36 benchmarks from 3 benchmark suites and found that they considerably improve the runtime on 6 benchmarks, with an improvement of over 5% and up to 24.62%. In almost all cases, our IPEA-based inlining resulted in smaller image sizes compared to existing purely PEA-driven approaches. The tuning also showed that trading-off run times, compilation times and generated code-size can individually improve each of these metrics. Decent compilation overhead suggests that our algorithm can improve performance in Just-in-Time compilation with some re-tuning. Furthermore, we found compelling evidence that factors beyond allocation reduction may be used to predict performance improvements, and we believe this is a promising research direction.

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