

Homeostasis: Design and Implementation of a Self-Stabilizing Compiler

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Mainstream compilers perform a multitude of analyses and optimizations on the given input program. Each analysis (such as points-to analysis) may generate a *program-abstraction* (such as points-to graph). Each optimization is typically composed of multiple alternating phases of *inspection* of such program-abstractions and *transformations* of the program. Upon transformation of a program, the program-abstractions generated by various analyses may become inconsistent with the modified program. Consequently, the correctness of the downstream inspection (and consequent transformation) phases cannot be ensured until the relevant program-abstractions are *stabilized*; that is, the program-abstractions are either invalidated or made consistent with the modified program. In general, the existing compiler frameworks do not perform automated stabilization of the program-abstractions and instead leave it to the compiler pass writers to deal with the complex task of identifying the relevant program-abstractions to be stabilized, the points where the stabilization is to be performed, and the exact procedure of stabilization. In this paper, we address these challenges by providing the design and implementation of a novel compiler-design framework called *Homeostasis*.

Homeostasis automatically captures all the program changes performed by each transformation phase, and later, triggers the required stabilization using the captured information, if needed. We also provide a formal description of *Homeostasis* and a correctness proof thereof. To assess the feasibility of using *Homeostasis* in compilers of parallel programs, we have implemented our proposed idea in IMOP, a compiler framework for OpenMP C programs. Further, to illustrate the benefits of using *Homeostasis*, we have implemented a set of standard data-flow passes, and a set of involved optimizations that are used to remove redundant barriers in OpenMP C programs. Implementations of none of these optimizations in IMOP required any additional lines of code for stabilization of the program-abstractions. We present an evaluation in the context of these optimizations and analyses, which demonstrates that *Homeostasis* is efficient and easy to use.

CCS Concepts: • **Software and its engineering** → **Compilers; Frameworks; Development frameworks and environments; Design patterns; Incremental compilers.**

Additional Key Words and Phrases: compiler design, multi-pass compilation, self-stabilization

ACM Reference Format:

Aman Nougrihiya and V. Krishna Nandivada. 2024. Homeostasis: Design and Implementation of a Self-Stabilizing Compiler. *ACM Trans. Program. Lang. Syst.* 1, 1, Article 1 (January 2024), 58 pages.

1 INTRODUCTION

Modern compilers often span millions of lines of code, and perform a multitude of analyses and optimizations on the given input program. Each analysis (such as points-to analysis, call-graph construction, and so on) derives some meaningful information about the program, in the form of a *program-abstraction* (such as points-to graph, call-graph, and so on). An optimization typically involves multiple alternating phases of inspections of such program-abstractions, and transformations of the program. In the *inspection phase*, the optimization invokes the required analyses, if needed, and inspects various program-abstractions to discover opportunities of optimization in the program. In the *transformation phase*, the optimization pass may use the results of the inspection phase to modify the program. Consequently, the transformation phase may render various

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2024. 0164-0925/2024/1-ART1 \$15.00
<https://doi.org/>

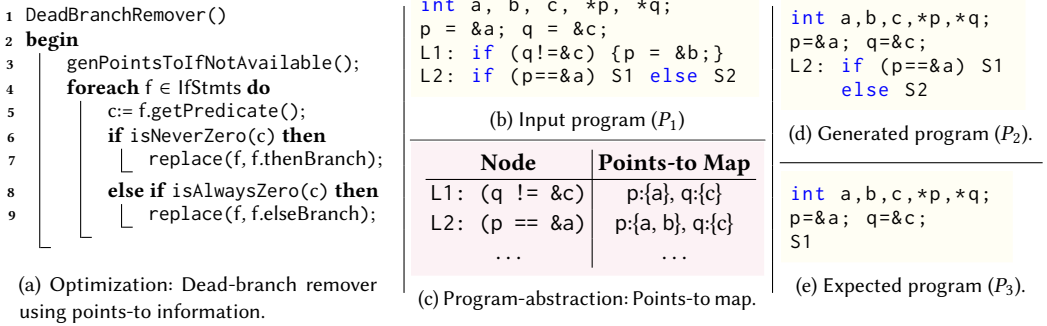


Fig. 1. An example optimizer pass, and a program-abstraction. After removing the if-statement at L1 from the input program P_1 to obtain P_2 , the optimizer will not be able to remove the else-branch of L2, thereby not producing the expected optimal code P_3 : In the absence of stabilization, the points-to map from Fig. 1c does not get updated after removal of the statement at L1. Consequently, with the stale points-to information, the optimizer cannot infer that pointer p may point only to a at the predicate at L2 in P_2 .

existing program-abstractions inconsistent with the modified program. As a result, the existing program-abstractions must be *stabilized* before their next use; that is, the program-abstractions must be regenerated or updated such that they correspond to the modified program. Otherwise, the correctness of the future inspection phases of the current or downstream optimizations cannot be ensured. This, in turn, can negatively impact the optimality and even the correctness of the overall compilation. We illustrate these problems using an example.

Example 1.1. In Fig. 1a, we show the pseudocode of a standard compiler optimization that utilizes points-to information to remove dead branches from a program. For simplicity, we assume that the predicates do not contain any side-effects, and that there are no arbitrary goto-statements in the input program. The `isNeverZero` method returns true when it can be statically determined that the given expression can never evaluate to zero; similarly, the `isAlwaysZero` method returns true only when the expression would always evaluate to zero. Further, assume that both these methods use the points-to information to statically evaluate the given expression. Hence, in the inspection phase, the optimization pass invokes a flow-sensitive points-to analysis pass to obtain the required points-to map (a program-abstraction), if not already available. Further, in this phase, the optimization utilizes the methods `isNeverZero` and `isAlwaysZero` to determine whether an if-statement can be replaced with either of its branches. Accordingly, the transformation phase may be triggered to perform the replacement using the `replace` method (Lines 7 and 9). Note that these transformation and inspection phases may alternate in the presence of multiple if-statements in the program.

This simple and intuitive code of dead-branch remover may not work as expected: Consider a sample input program P_1 shown in Fig. 1b, and a part of its points-to map in Fig. 1c. Say the optimization processes the if-statement at label L1 first, and relying on the points-to map at the predicate, replaces the if-statement with its else-branch (i.e., an empty body), transforming program P_1 to P_2 (Fig. 1d). This transformation renders the points-to map stale: at the predicate (p == &a), the pointer p continues to point to $\{a, b\}$, instead of $\{a\}$, the more precise points-to set. In the absence of stabilization of the points-to map, the optimizer pass in its next iteration cannot determine that the predicate at L2 will never be zero. As a result, the optimizer will not generate the optimal program P_3 ; instead, P_2 will be the final output of the optimizer. Note that such issues

Upon adding a new optimization O_n	Upon adding a new program analysis \mathcal{A}_n
Q_1 . Which existing program-abstractions need to be stabilized by O_n ?	Q_1^T . Which existing optimizations need to stabilize the program-abstraction of \mathcal{A}_n ?
Q_2 . Where to invoke the stabilization code in O_n , for the existing program-abstractions?	Q_2^T . Where to invoke the stabilization code for \mathcal{A}_n , in each of the existing optimizations?
Q_3 . How to stabilize each of the existing program-abstractions, in O_n ?	Q_3^T . How to stabilize the program-abstraction of \mathcal{A}_n , in each of the existing optimizations?

Fig. 2. Key stabilization-related question to be addressed while adding new compiler passes.

of sub-optimality (and even correctness) may also be observed in the downstream optimization passes, until the points-to map is stabilized.

To perform the correct stabilization manually and efficiently, we need to identify all the places in the dead-branch remover where the points-to map may require stabilization. Further, we also need to be concerned about the stabilization of all the other program-abstractions as well: (i) We need to identify the program-abstractions (for example, use-def information [Muchnick 1998]) that may become stale because of the current optimization, and may be required by other downstream optimizations and analyses. (ii) For each of these program-abstractions, we need to identify the places in the dead-branch remover where the program-abstraction needs to be stabilized. In practice, in the presence of multiple optimizations and analyses (hereafter, collectively referred to as *compiler passes*), this problem compounds as we need to be concerned about stabilization of all the program-abstractions in all the optimizations.

In summary, we observe that to avoid the generation of such sub-optimal and/or incorrect programs upon compilation, three key stabilization-related questions shown in Fig. 2 need to be addressed while adding/modifying an optimization or analysis. In conventional compilers, such as LLVM [Lattner and Adve 2004], GCC [Stallman and GCC-Developer-Community 2009], Soot [Vallée-Rai et al. 2010], Rose [Quinlan et al. 2013], JIT compilers (OpenJ9 [IBM 2017], HotSpot [Oracle 1999], V8 [Google 2001]), and so on, the onus of addressing these key questions lies mostly on the compiler pass writers. Some such instances are shown in Fig. 3. This manual process often leads to complex and error-prone codes. Further, as the number and complexity of compiler passes increase (for example, LLVM has more than 347 passes), addressing these questions precisely and efficiently becomes progressively more difficult. This may lead to correctness/efficiency bugs (see bug-fixing commits on GitHub [LLVM-Developer-Community 2019a,b,c, 2020a,b,c,d,e,f,g,h,i, 2021a,b,c]). With the compiler development effort spanning multiple decades involving (sometimes) hundreds of developers, it becomes extremely challenging to manually solve these problems, as no developer of a compiler pass might possess a clear understanding of the semantics of all the hundreds of other passes already present in the compiler.

To mitigate these issues, the focus of this paper is to perform automatic stabilization of all the relevant program-abstractions, when needed, in response to any program modification. We term a compiler that provides this guarantee as a *self-stabilizing* compiler.

There have been various attempts [Blume et al. 1995; Brewster and Abdelrahman 2001; Carle and Pollock 1989; Carroll and Polychronopoulos 2003; Nilsson-Nyman et al. 2009; Reps et al. 1983] towards enabling automated stabilization of *specific* program-abstractions, in response to program transformations in serial programs. However, these works suffer from different drawbacks. For example, Carle and Pollock [1989]; Nilsson-Nyman et al. [2009]; Reps et al. [1983] require that program-abstractions have to be expressed as context-dependent attributes of the language constructs; this is too restrictive. Similarly, besides the non-object-oriented nature of the approach of Carroll and Polychronopoulos [2003], it is unclear how their restrictive techniques can (i) be

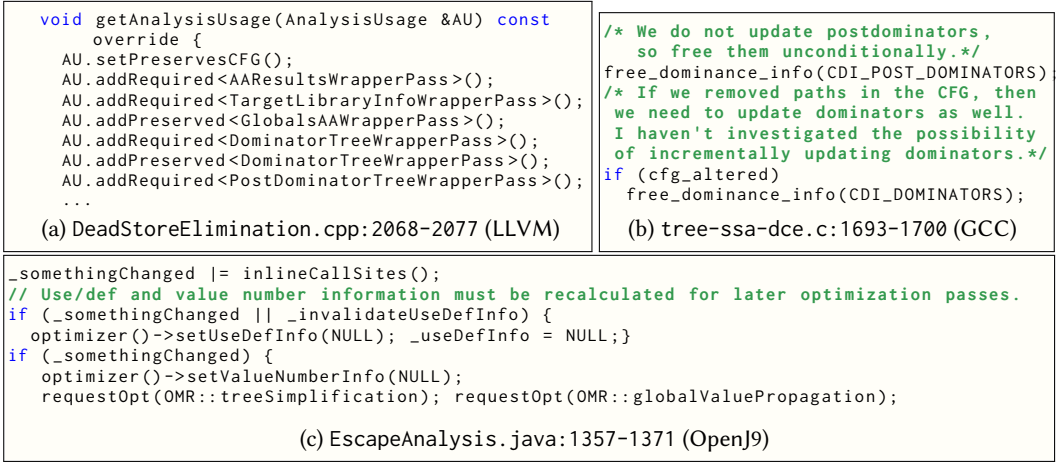


Fig. 3. Example snippets from LLVM, GCC, and OpenJ9 demonstrating manual stabilization. The comments are written by the pass writers. Snippet (a) shows how pass writers manually specify pass dependencies in LLVM. Snippet (b) shows how some selected program-abstractions are invalidated manually in GCC. Snippet (c) depicts a case where the pass writer has manually written additional code for stabilization after a transformation phase.

applied to compilers that use non-hierarchical IRs, which is a common practice in modern compilers (such as LLVM, Soot, OpenJ9, IMOP [Nougrahiya and Nandivada 2019], and so on), (ii) be used for data-flow analyses that are not based on structural analysis [Muchnick 1998], and (iii) automate the resolution of pass dependencies, or handle manually-defined dependencies. Blume et al. [1995]; Brewster and Abdelrahman [2001] handle only a small set of specific program-abstractions (such as symbol and scope information); this is insufficient. To the best of our knowledge, there are no compiler designs or implementations¹ that address the challenges discussed above for any arbitrary IR, optimization and analysis.

Proposed solution. In this paper, we propose a novel and efficient compiler-design framework called *Homeostasis* that addresses the above discussed issues related to stabilization, in the context of object-oriented compilers. *Homeostasis* efficiently captures a summary of all the program changes performed by each transformation phase. Later, if/when an attempt is made to read the possibly stale program-abstraction of a program analysis, *Homeostasis* triggers the required stabilization using the captured information, before returning the correct value of the program-abstraction. With *Homeostasis*, no additional coding effort is required by the optimization writers to guarantee self-stabilization of the existing (or future) program-abstractions. For example, in an implementation of the dead-branch remover (Fig. 1a) in a *Homeostasis*-enabled compiler, *Homeostasis* would ensure self-stabilization of the points-to map, without requiring any additional code in the optimization; this will lead to generation of the expected optimal program P_3 .

In *Homeostasis*, for ensuring self-stabilization, analysis writers need to spend minimal effort to only conform to the design of *Homeostasis*, while being oblivious to all the (existing/future) optimizations. As part of this conformance, if the analysis writer decides to use an incremental update algorithm for stabilization (instead of full recomputation), the writer has to clearly demarcate the code to incrementally update the program-abstraction. Note that such a code is needed even

¹Our observation is based on our manual inspection of more than 50 mainstream and experimental compilers [Nougrahiya and Nandivada 2023].

when the stabilization is performed manually; hence this is not an additional overhead. In essence, while working with a *Homeostasis*-enabled compiler, a compiler pass writer need not manually address any of the key questions listed in Fig. 2, and hence incurs no additional overhead.

For the ease of exposition, we have discussed *Homeostasis* using IMOP [Nougrahiya and Nandivada 2019] a Java-based compiler in the context of a C-like language. We believe that the underlying principles of the *Homeostasis* are quite generic in nature, and that they can be extended to any object-oriented compilers for any serial/parallel languages.

Note that the goal of this paper is not to propose any new incremental analysis technique (or any new analysis technique or optimization for that matter). The key goal of this paper is to present a scheme which ensures that correct values of various program-abstractions are obtained efficiently by all the compiler passes, without needing any additional manual effort in these passes.

Contributions:

- We present *Homeostasis*, a generalized compiler-design framework for *self-stabilization*, which can be added to any object-oriented compiler infrastructure to ensure that all program-abstractions are automatically kept consistent with the modified program during the compilation process. The underlying concepts of *Homeostasis* can be used in any object-oriented compiler for serial/parallel languages to realize self-stabilization.
- To overcome the possible overheads resulting from the naive tracking of the program changes, we present a novel compression algorithm that reduces the memory requirements significantly.
- We give a formal description of *Homeostasis* and a correctness proof thereof, which guarantees that in a *Homeostasis*-enabled compiler, any compiler pass attempting to access the program-abstraction of an analysis will see only the stable value of that program-abstraction.
- We have implemented all the key components of *Homeostasis* in IMOP, a source-to-source Java-based compiler infrastructure for OpenMP C programs, as a successful proof-of-concept. Further, to demonstrate the benefits of using *Homeostasis*, we have implemented (i) HIDFA, a generic, self-stable, inter-thread, flow-sensitive, context-insensitive iterative data-flow analysis (IDFA) and six instantiations thereof, (ii) a set of optimization passes, collectively termed as BarrElim, that removes redundant barriers from OpenMP C programs; BarrElim uses a set of nine optimization passes, each of which may involve multiple alternating phases of inspection and transformation. As expected, the implementation of the set of BarrElim optimization passes required zero additional lines of code for stabilization, and the stabilization of the analyses required minimal effort.
- We present an evaluation over a set of twenty-four benchmarks from four benchmark suites NPB [Van der Wijngaart and Wong 2002], SPEC OMP [Aslot et al. 2001], Sequoia [Seager, M 2008] and IMSuite [Gupta and Nandivada 2015], performed over two different platforms. We show that *Homeostasis* makes it easy to write optimization and analysis passes, and it leads to faster compilation times compared to the possible alternative of typical manual stabilization performed by experts. We also show that our proposed optimizations lead to significant memory savings.

The rest of the manuscript is organized as follows: We give a brief background of the relevant concepts used in the manuscript in Section 2. We present our design of *Homeostasis* in Section 3. We give a formal description of *Homeostasis*, along with a correctness argument in Section 4. We discuss some salient features of *Homeostasis* in Section 5. In Section 6, we discuss an instantiation of *Homeostasis* in a real-world compiler. In Section 7, we briefly describe our implementation and present a detailed evaluation of *Homeostasis*. In Section 8, we discuss some of the relevant prior work, before concluding in Section 9.

2 BACKGROUND AND TERMINOLOGY

We now give a brief description of certain relevant concepts and terminologies used in this paper.

Representation of program IR. Without loss of generality, in this paper we assume each program's intermediate representation (IR) is represented as a graph. For instance, when the control flow graph (CFG) is the IR, nodes in the graph represent the basic-blocks, and the edges model the flow of control. Similarly, if three-address code (TAC) is the IR, then the TAC instructions are the nodes, and the edges may be of many types, such as control-flow edges, SSA edges [Muchnick 1998], and so on.

Analysis, Analysis-instances, and Program-abstractions. In the context of compilers written in an object-oriented language, we assume that each analysis (for example, Andersen's points-to analysis [Andersen 1994]) is implemented as a class. During compilation, one or more of these classes are instantiated to provide analysis-instances. Each such analysis-instance, say A_i , can be seen as a function from the set of programs to a set of program-abstractions. A program-abstraction may refer to any data-structure that denotes some meaningful information generated by an analysis-pass, about the input program at compile time. In this manuscript, we use $A_i(P)$ to denote the program-abstraction generated by running the underlying analysis of A_i on a program P . For example, Fig. 1c shows a part of the points-to map, the program-abstraction of the corresponding points-to analysis.

Stable program-abstraction. At any point during compilation, if the current program IR is P , and the stored program-abstraction for an analysis-instance A_i is v , then we term v as the *stable* program-abstraction of A_i , iff $A_i(P) = v$.

Elementary transformations. Based on the grammar of the language under consideration, corresponding to each type of program node with a *body* (such as a block in LLVM IR, or a while-statement construct), we identify a fixed set of primitive transformations which can add/delete/modify the logical components of the program node (such as an instruction of a block, or the predicate of a while-statement); we term such transformations as *elementary transformations*.

3 HOMEOSTASIS: DESIGNING SELF-STABILIZING COMPILERS

Considering the stabilization challenges inherent in the design of conventional compilers, we now present a novel solution in the form of a new object-oriented compiler-design framework named *Homeostasis* (in Sections 3.1, 3.2, and 3.3). In order to access/maintain stable program-abstractions in a *Homeostasis*-enabled compiler, while the analysis pass writers have to spend minimal effort, the optimization pass writers need not spend any effort (elaborated in Section 3.4). We formalize the design of *Homeostasis* and present a correctness argument in Section 4.

3.1 Overview of *Homeostasis*

Homeostasis realizes self-stabilization by tracking the program modifications resulting from different transformations. These program modifications are then used by *Homeostasis* to perform the stabilization of relevant program-abstractions if/when required.

3.1.1 Design pattern employed by *Homeostasis*. To track program modifications resulting from different transformations, *Homeostasis* uses an extension of the popular Observer pattern [Gamma et al. 1995]. The Observer pattern defines a one-to-many dependency between multiple entities, where an update in one entity (termed *subject*), is notified to all the other registered entities (termed *observers*). Fig. 4 shows a class diagram (in Object Modeling Technique, or OMT, notation [Rumbaugh et al. 1991]) which gives an overview of the key design elements of *Homeostasis*: (i) Node class and its subtypes (such as *WhileStmt*) correspond to the subject(s), (ii) all the program analyses

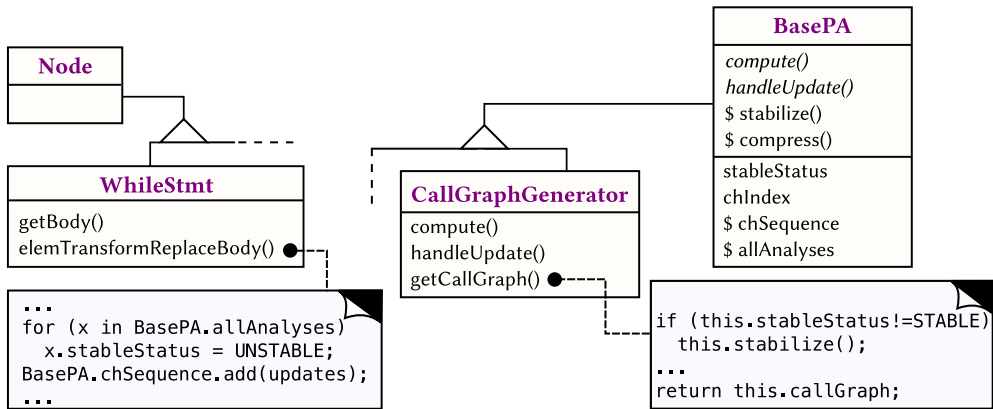


Fig. 4. Class-diagram of *Homeostasis*, depicting key classes and methods, in the OMT notation: class names are shown in **bold**; abstract methods in *italics*; static members are preceded with a \$; inheritance is shown using a triangle on the edge, pointing towards the base class; and dashed arrows show code snippets for methods.

(such as *CallGraphGenerator*) along with their base class (*BasePA*) correspond to the observers, and (iii) program modifications (such as *WhileStmt::elemTransformReplaceBody*) correspond to the various updates to the subject(s). The use of the Observer pattern ensures that these updates (program modifications) are correctly notified to all the observers (program analysis passes).

Homeostasis goes beyond the standard design of the Observer pattern in two key ways. First, instead of triggering updates of the values (i.e., program-abstractions) maintained by various program analyses immediately upon a program modification, *Homeostasis* simply marks the program-abstractions as stale (or *UNSTABLE*). The actual update of each program-abstraction is performed *lazily*, by deferring the update to a later stage where an attempt is made to read from the program-abstraction, using its *getters* (such as *CallGraphGenerator::getCallGraph*). Second, instead of storing a copy of the program changes corresponding to such pending updates individually with each program analysis (each of which may need to handle a different list of pending updates), *Homeostasis* maintains a central copy of these program changes in an efficient data-structure (discussed in Section 3.3).

3.1.2 Nodes of intermediate representation (IR). In Fig. 4, the *Node* class represents the super class of all the types of nodes in the program’s intermediate representation (IR). Each such class may define a fixed set of *elementary transformation* methods that may be invoked to add/modify/delete any of the logical component(s) of the node. For example, in Fig. 4, *elemTransformReplaceBody* is an elementary transformation method defined on the node *WhileStmt*; this transformation replaces a component (the loop body) of the associated while-statement with the given argument. *Homeostasis* requires that (i) all program modifications are performed via a sequence of invocations of elementary transformations, (ii) each elementary transformation marks each program-abstraction in the system as stale by resetting the corresponding *stableStatus* flag, and (iii) each elementary transformation stores the details of the resulting IR changes in the global sequence of program-changes, *chSequence*, maintained by *BasePA*, to be read by the individual program analyses as and when required. Note that the first requirement does not restrict the availability of high-level transformation APIs to the optimization-pass writers; each such API may be adapted by the compiler-writers such that it invokes one or more elementary transformations directly and/or via invocation of other similar high-level APIs. For example, a high-level transformation API to delete a block of statements can be implemented as a sequence of invocations of an elementary transformation that deletes

individual statements. Note that updating the high-level APIs is a one-time effort on the part of compiler-writers; no additional effort is needed from the optimization-pass writers.

3.1.3 Program analysis pass classes. In *Homeostasis*, each program analysis is implemented as a class that inherits from the abstract class *BasePA*, as shown in Fig. 4; this class contains various key data structures and methods that are required to enable self-stabilization of the associated program-abstractions. *BasePA* maintains a global set, *allAnalyses*, of all instances of the analysis classes (henceforth, referred to as *analysis-instances*). Further, a sequence of all changes to the IR, resulting from various elementary transformations, is maintained in a global list, named *chSequence*. During stabilization trigger for any analysis-instance, if needed, the exact set of IR changes whose impact needs to be incorporated in its program-abstraction, is obtained using *chIndex*, an index to some location in the list *chSequence*. Note that both *allAnalyses* and *chSequence* are initialized to empty set/list. *Homeostasis* may periodically employ the *compress* method to reduce the space requirements of *chSequence*, depending upon the *chIndex* value of various analysis-instances in *allAnalyses* (details in Section 3.2).

Fig. 4 shows *CallGraphGenerator* as a concrete class of *BasePA*, which implements the call-graph generation routines and maintains the program-abstraction denoting the call-graph information of the program. In *Homeostasis*, each program-abstraction can only be queried through its corresponding *getter* methods (such as *getCallGraph* in *CallGraphGenerator*). Before returning the requested information, each *getter* method should check if the program-abstraction is stable (by inspecting the value of *stableStatus* flag) and trigger stabilization if needed (by invoking the method *stabilize*). Depending upon the mode of stabilization, discussed next, the *stabilize* method provided by *Homeostasis* internally invokes either the *compute* or the *handleUpdate* method, passing them the correct arguments, if any.

3.1.4 Modes of stabilization. An important defining characteristic of self-stabilization is the time and manner in which a program-abstraction is stabilized under program modifications. Accordingly, the stabilization mode can vary along the following two dimensions: eager versus lazy, and invalidate versus update.

Eager versus lazy. After an elementary transformation, the stabilization of an abstraction may either get triggered (i) immediately (eager stabilization), or (ii) only in response to the first read request made on the program-abstraction after the transformation (lazy stabilization). Note that in practice, not every program-abstraction is read after each program modification. Hence, to avoid the overheads due to the redundant invocation of the stabilization routines in case of eager stabilization, *Homeostasis* advocates the use of lazy modes of stabilization.

Invalidate versus update. In response to one or more program modifications, the resulting stabilization of a program-abstraction may either (i) involve the complete invalidation of the program-abstraction (leads to regenerating the program-abstraction from scratch), or (ii) be able to incrementally update the program-abstraction based on the modifications. *Homeostasis* attains the fully-automated stabilization for invalidate modes without requiring *any* stabilization-specific code from the analysis-pass writers. In invalidate modes, the *stabilize* method reinitializes the program-abstraction, and invokes the *compute* method to regenerate the program-abstraction from scratch; note that the code for this method would exist even in the absence of *Homeostasis*. In contrast, to realize the update modes of stabilization, *Homeostasis* provides a well-defined method, *handleUpdate*, to be implemented by the analysis-pass writers, which it later uses during automated stabilization triggers, by passing the correct set of program changes to be handled. Though the update modes may seem much more efficient than the invalidate modes, in practice the difference in their performance depends on a number of factors, such as the number of program modifications, the complexity of the associated incremental update, and so on. Further, note that enabling the update


```

1 Function replaceX(newNode) // Method to replace the component X with newNode
2   oldX = getComponentX();
3   // Step A: Save nodes and edges affected by the removal of old node.
4   Set removedNodes = {oldX};
5   Set removedEdges = {...outgoing edges from, and incoming edges to, oldX..};
6   Set addedEdges = {...new edges formed due to the removal of oldX..};
7
8   // Step B: Perform the actual program update by invoking appropriate writer(s).
9   Replace the component X with newNode;
10
11  // Step C: Save nodes and edges affected by the addition of new node.
12  Set addedNodes = {newNode};
13  removedEdges = removedEdges  $\cup$  {...old edges removed due to the addition of newNode..};
14  addedEdges = addedEdges  $\cup$  {...outgoing edges from, and incoming edges to, newNode..};
15
16  // Step D: Communicate relevant information to each analysis.
17  for anl  $\in$  BasePA.allAnalyses do anl.stableStatus = UNSTABLE;
18  BasePA.chSequence.addFourTuple(addedNodes, removedNodes, addedEdges, removedEdges);

```

Fig. 5. Template of an elementary transformation in *Homeostasis* that replaces the component X of a program node with newNode .

modes of stabilization for certain program-abstractions can be a challenging task. Considering such issues, *Homeostasis* supports both invalidate as well as update modes of (lazy) stabilization.

On the basis of these two dimensions, *Homeostasis* supports the following two modes of stabilization for any program-abstraction: (i) Lazy-Invalidate (**LZINV**), and (ii) Lazy-Update (**LZUPD**).

3.2 Components of *Homeostasis*

In this section, we expand on the design details of different components of *Homeostasis* (in Sections 3.2.1, 3.2.2, and 3.2.3). In Section 3.2.4, we discuss the details of how effective changes to the IR, resulting from a sequence of elementary transformations, are obtained during stabilization in *Homeostasis*. Finally, in Section 3.2.5, we demonstrate how stabilization is performed in the presence of *Homeostasis*, using the example from Fig. 1a.

3.2.1 Structure of Node and its subclasses. Stabilization of program-abstractions in response to the modifications performed by an optimization pass, can intuitively be done in two ways: (i) the optimization pass directly modifies the program-abstraction, or (ii) the program analysis pass performs the stabilization internally, possibly based on the exact modifications that have been performed on the program. The first option can be complicated, and goes against the spirit/design principles of object-oriented programming. Hence, we use the latter option in *Homeostasis*.

Elementary transformations. *Homeostasis* uses elementary transformations to communicate the program modifications performed by different optimization passes to the program analyses. As mentioned in Section 3.1, in *Homeostasis*, all program modifications happen only via elementary transformations. Each elementary transformation in *Homeostasis* (i) collects the information about addition/deletion of IR nodes and edges, and (ii) makes the collected information visible to every program analysis object. We elaborate this using an example template of a generic elementary transformation method that replaces a component C with a new node, as shown in Fig. 5. The templates for the elementary transformation methods that add (or remove) components can be trivially derived from the above template. The template method in Fig. 5 has four main steps.

Step A: This step applies to all those elementary transformations that may remove a node from the IR. Besides storing the removed node in `removedNodes`, this step records all the edges that have been removed (in `removedEdges`) or added (in `addedEdges`) as a result of removal of the node `oldX`.

```

Elementary transformation : WhileStmt::setBody(Statement newBody);
oldBody                  = this.body
Set removedEdges         = {(this.predicate, oldBody), (oldBody, this.predicate), ...}
Set addedEdges           = {(this.predicate, newBody), (newBody, this.predicate), ...}
Set removedNodes         = {oldBody}
Set addedNodes           = {newBody}

```

Fig. 6. Values for the various sets used in Fig. 5, for a concrete elementary transformation that replaces the body of a while-statement in the IMOP compiler framework (WhileStatementCFGInfo.java: 81-146).

Step B: This step performs the actual modification to the IR. For example, in Fig. 5, the component X of a program node is replaced with the provided new node. This step contains the code that would comprise the body of an elementary transformation in the absence of *Homeostasis*.

Step C: This step applies to only those elementary transformations that add a node to the IR. Besides storing the added node in `addedNodes`, this step records all the edges that have been removed (in `removedEdges`) or added (in `addedEdges`) as a result of addition of the node `newNode`.

Step D: In this step, *Homeostasis* first marks the status of each program analysis as UNSTABLE, so that the corresponding stabilization routine is invoked whenever its program-abstraction is read next (using the getter methods). As discussed in Section 3.1.4, for using the LZUPD mode of stabilization a program analysis requires information about all the added/removed nodes and edges. Instead of maintaining a copy of this information with each program analysis, for efficiency, *Homeostasis* maintains it globally in the list `chSequence`, in the `BasePA` class. This list is 0-indexed. Each element of `chSequence` is a four-tuple composed up of four sets: nodes added, nodes removed, edges added, and edges removed. At the end of each elementary transformation, the method `addFourTuple` appends the collected list of IR changes to the end of `chSequence`.

We now highlight three important points related to these four steps. (1) The exact definitions of Steps A-C depend on (i) the exact type of the node being transformed, (ii) the specific component being added/removed/replaced, and (iii) the semantics of the underlying language. For instance, in case of OpenMP compilers, addition/removal of a node containing flush-directives, would necessitate changes to the inter-task edges (or their equivalents) in the IR; these edges are used to denote inter-task communication via shared variables. The exact set of edges added/removed depend on the underlying semantics of OpenMP constructs. (2) An elementary transformation may also result in addition or removal of edges without any changes in the nodes. In such cases, Steps A and C capture the impacted edges accordingly. Further, if an elementary transformation changes only the type, label, or annotation of an edge, then this can be represented as the replacement of the old edge with a distinct new edge; the former is added to the set `removedEdges` in Step A, and the latter is added to the set `addedEdges` in Step C. (3) In Step D, the list `chSequence` may grow prohibitively large with increasing number of invocations of elementary transformations; this list is maintained efficiently by *Homeostasis* using a *compression optimization*, explained later in Section 3.3.

Example 3.1 (Concrete elementary transformation). Fig. 6 shows the values for the various sets used in Fig. 5, for an elementary transformation method that replaces the body component of a while-statement with a given argument (`newBody`). This example has been taken from our implementation of *Homeostasis* in the IMOP [Nougrahiya and Nandivada 2019] compiler framework for OpenMP C programs. Note that in addition to the elements shown, the sets `removedEdges` and `addedEdges` also include all the jump edges (due to jump statements, such as `continue`, `break`, etc.) that will be removed/added from the IR as a result of this elementary transformation. Further, these sets

```

1 /* BasePA constructor */
2 public BasePA::BasePA(StabilizationMode mode) {
3     BasePA.allAnalyses.add(this);
4     this.compute();
5     this.stabilizationMode = mode;
6     if (this.stabilizationMode == LZINV) {
7         this.chIndex = -1;}
8     else { // this.stabilizationMode = LZUPD
9         this.chIndex = BasePA.chSequence.size() - 1;}

```

(a) Constructor of the base analysis class.

```

1 /* Template of a program-abstraction getter */
2 public V Analysisy::get() {
3     if (this.stableStatus==UNSTABLE) {
4         this.stabilize();}
5     return this.abstraction;}

```

(b) Template of a getter method of a sample analysis.

```

1 /* Program-abstraction stabilizer */
2 private void BasePA::stabilize() {
3     this.stableStatus = STABLE;
4     if (this.stabilizationMode == LZINV) {
5         this.compute();}
6     else { // this.stabilizationMode = LZUPD
7         this.handleUpdate();
8         this.chIndex = BasePA.chSequence.size() - 1;}}

```

(c) Implementation of the stabilizer used by *Homeostasis*.

Fig. 7. Constructor and stabilizer methods of BasePA, and a template getter of a program-abstraction.

also contain inter-task edges² that are used as a part of the IR to model the parallel semantics of OpenMP C.

3.2.2 Structure of BasePA and its subclasses. *Homeostasis* maintains a global set `allAnalyses` of program analysis instances, by adding the reference of each program analysis instance (at the time of its construction) to `allAnalyses`; see Fig. 7a. In addition, the constructor (i) invokes the `compute` method, overridden by each program analysis class, to run the analysis from scratch and populate the corresponding program-abstraction, and (ii) sets the initial value for `chIndex`, as per the given mode of stabilization. In the case of LZINV mode of stabilization, `chIndex` is set to `-1` as this field is not used in LZINV mode. In the case of LZUPD mode of stabilization, the value of `chIndex` is used to identify those elements in the global list `chSequence` whose impact needs to be taken into account in order to obtain the stable program-abstraction. These elements are the change elements that lie at indices *beyond* `chIndex` in `chSequence`.

The getter method of any program analysis follows the template shown in Fig. 7b, and invokes the `stabilize` method, if needed, to return the expected value. See a snippet of the method `getCallGraph` in Fig. 4 for an example.

3.2.3 Structure of program-abstraction stabilizers. As discussed in Section 3.1.4, *Homeostasis* allows each program-abstraction to be stabilized in one of the two supported modes of stabilization: LZINV and LZUPD. Naturally, the procedure to stabilize a program-abstraction depends on the corresponding mode of stabilization. For the LZINV mode of stabilization, the `stabilize` method (Fig. 7c) simply

²IMOP uses the concept of inter-task edges between flush operations of a parallel region to statically model the communication that may happen using shared variables between different threads.

```

1 Function getNetChanges(chIndex) // Obtain the net IR changes to be handled for a given chIndex
2   netChanges = <>;
3   foreach ind from (chIndex + 1) to (BasePA.chSequence.size() - 1) do
4     | netChanges = mergeChanges(netChanges, BasePA.chSequence.get(ind));
5   return netChanges;
6 Function mergeChanges(ch1, ch2) // Obtain the net IR changes for the given changes
7   mergedChanges = <>;
8   mergedChanges.addNodes = (ch1.addNodes\ch2.remNodes) ∪ (ch2.addNodes\ch1.remNodes);
9   mergedChanges.remNodes = (ch1.remNodes\ch2.addNodes) ∪ (ch2.remNodes\ch1.addNodes);
10  mergedChanges.addEdges = (ch1.addEdges\ch2.remEdges) ∪ (ch2.addEdges\ch1.remEdges);
11  mergedChanges.remEdges = (ch1.remEdges\ch2.addEdges) ∪ (ch2.remEdges\ch1.addEdges);
12  return mergedChanges;

```

Fig. 8. Algorithms to obtain the net IR changes for those elements in chSequence whose impact needs to be taken into account by the handleUpdate method to obtain the stable program-abstraction.

reruns the analysis from scratch by invoking the compute method. Thus in the LZINV mode, the program analysis writer needs to write no additional code to stabilize the program-abstraction.

For the LZUPD mode of stabilization, the stabilize method performs self-stabilization of the abstraction by invoking the method handleUpdate. To utilize the arguably more efficient LZUPD mode, the program analysis writer needs to ensure that the desired incremental update algorithm is implemented in the handleUpdate method. This method should define the impact of the net IR changes, given in terms of the four sets: nodes added, nodes removed, edges added, and edges removed. The stabilize method invokes this handleUpdate method by providing the correct argument in the form of net IR changes (obtained as discussed next), followed by setting the chIndex to the index of the last element of chSequence.

3.2.4 Obtaining Net IR changes. In order to obtain the net IR changes to be taken into account by the handleUpdate method during a stabilization trigger, *Homeostasis* provides getNetChanges function, whose algorithm is shown in Fig. 8. The net IR change is obtained by merging the IR changes denoted by all the elements (if any) that lie to the right of the location indexed by chIndex in the global list chSequence. The exact merge operation is specified in the function mergeChanges in Fig. 8. We demonstrate the working of these functions using a concrete example.

Example 3.2 (Net IR changes). Fig. 9 illustrates a concrete example of how the net IR changes are obtained by *Homeostasis*, when required by the handleUpdate method during stabilization trigger of some analysis-instance. The value of chIndex for the analysis-instance is m , and the size of chSequence list is $(m + 4)$. Assume, the last three elementary transformations are:

- (i) Addition of node n_1 , resulting in addition of edges $\{e_1, e_2\}$.
- (ii) Replacement of node n_3 with node n_2 , resulting in addition of edges $\{e_3, e_4\}$, and removal of edges $\{e_5, e_6\}$.
- (iii) Replacement of the previously-added node n_1 with node n_4 , resulting in addition of edges $\{e_7, e_8\}$, and removal of the previously-added edges $\{e_1, e_2\}$.

The figure shows the corresponding set of IR changes, whose impact needs to be taken into account during the next stabilization trigger. The net IR changes are obtained by merging all the three IR changes using the mergeChanges function from Fig. 8 (represented using \oplus binary operator).

Now, we demonstrate how self-stabilization achieved by *Homeostasis*, as discussed in this section, fixes the issues in compilation of the example shown in Fig. 1a, without requiring any changes to the code written by the optimization-pass writer.

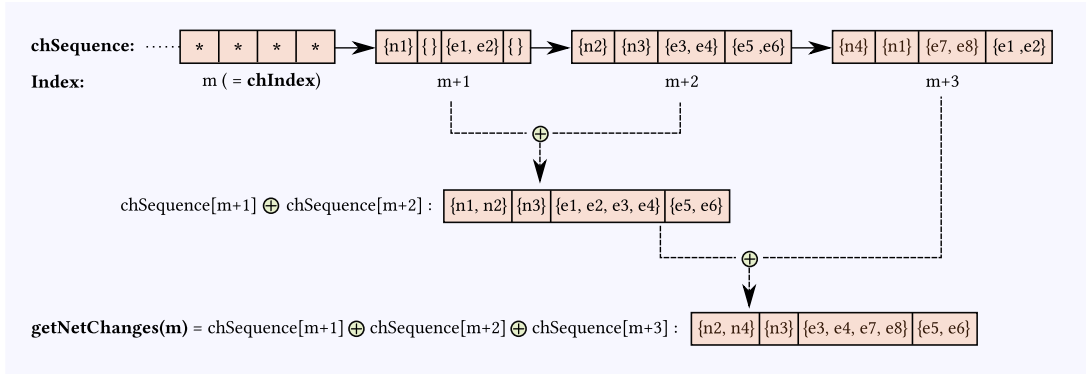


Fig. 9. A concrete example demonstrating how the net IR changes are obtained in the LZUPD mode of stabilization for some analysis-instance with value of `chIndex` as m . Size of the global list `chSequence` is $(m + 4)$. Each element of `chSequence` shows a four-tuple, corresponding to the following sets, in order: nodes added, nodes removed, edges added, and edges removed. The \oplus operator denotes the application of `mergeChanges` function from Fig. 8.

3.2.5 Example Compilation with Homeostasis. Let us revisit the example of the dead-branch remover optimization pass given in Fig. 1a, implemented in a *Homeostasis*-enabled compiler. Here, the `replace` method would invoke a sequence of elementary transformations, directly or indirectly through other high-level transformation APIs. These elementary transformations will capture the information about the resulting program modifications (say, from P_1 to P_2), and notify it to BasePA, thereby making the information available to all program analysis passes, including the points-to analysis pass. Later, when an attempt is made to read the points-to map using a getter of the analysis pass, say from within the methods `isNeverZero` and `isAlwaysZero`, the getter would automatically trigger stabilization of the points-to map internally such that the map correctly corresponds to P_2 . Hence, the compiler will be able to precisely deduce that the pointer `p` at L2 will now point only to variable `a`; this would trigger the removal of the false branch of the `if`-statement at L2, thereby generating the expected program, P_3 .

3.3 Compression Optimization: Efficient Tracking of Program Changes

The number of elementary transformations that a program undergoes during compilation can be substantial – compilation of large programs may easily involve several hundreds of elementary transformations. Hence, storing the individual IR-changes resulting from each elementary transformation separately in the global list `chSequence` can be prohibitively expensive in terms of space requirements. In order to address this challenge, *Homeostasis* employs a compression optimization that reduces the size of `chSequence`, relying on the following two key observations on the value of `chIndex` of analysis-instances that employ the LZUPD mode of stabilization.

Observation 1. *If the IR-changes corresponding to an elementary transformation have been taken into account by all the analysis-instances in the compiler, then the IR-changes need not be saved in `chSequence`.* In other words, if no analysis-instance has `chIndex` less than 0, then the first element (at index 0) in `chSequence` can be safely removed (while adjusting the `chIndex` of all the analysis-instances by decrementing 1 from their values).

Correctness argument. Recall that `chSequence` is a 0-indexed list. Further, as per the definition of `getNetChanges` from Fig. 8, the net IR changes to be used in the next stabilization trigger of an analysis-instance are obtained by merging the IR changes denoted by all the elements that lie

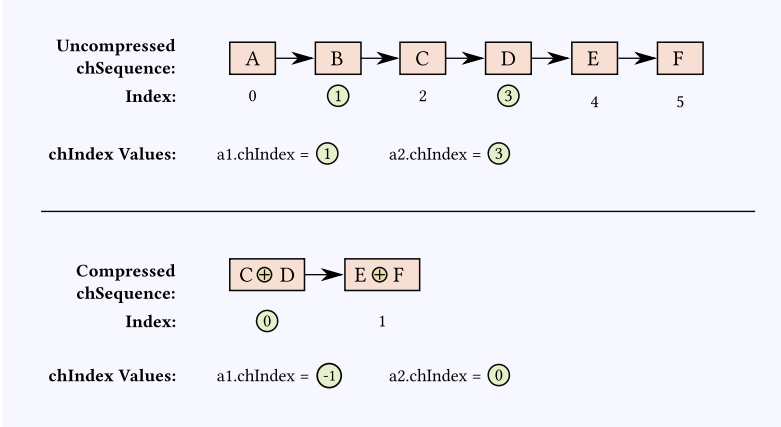


Fig. 10. Concrete example of compression optimization, demonstrating application of Observations 1 and 2. Those indices in `chSequence`, which are indexed to by `chIndex` of any analysis-instance in the compiler, are shown encircled.

beyond the index `chIndex` (of the analysis-instance) in `chSequence`. Hence, if there is no analysis-instance with `chIndex` less than 0, then the first element of `chSequence` (at index 0) will not be used to obtain the net IR changes during the stabilization trigger of any analysis-instance.

Observation 2. *If no analysis-instance has a `chIndex` that indexes into the location of an IR-change element in the `chSequence` list, then it is safe to merge the IR-change element to its successor, if any, in the list (while also adjusting the `chIndex` of various analysis-instances appropriately).*

Correctness argument. Consider an index i in `chSequence`, such that no analysis-instance has its `chIndex` as i . In this case, those analysis-instances whose `chIndex` is greater than i , will not consider the element at index i while obtaining their net IR changes during future stabilization triggers (see Fig. 8). Whereas, those analysis-instances whose `chIndex` is less than i , will anyway merge the elements at index i and $(i + 1)$ (if the element exists) to obtain the net IR changes for their next stabilization trigger. Hence, merging the elements at i and $(i + 1)$ in `chSequence`, will not affect the net IR changes for any analysis-instance during any of their future stabilization triggers.

The compression optimization employed by *Homeostasis* compresses the global `chSequence` list by repeated application of these two observations, until fixed-point. Note that *Homeostasis* does not dictate how often the compression optimization is invoked during the compilation. The exact details of the compression optimization, including details on how various corner cases are handled, are discussed formally in Section 4.2. In this section, we demonstrate the application of these two observations using a concrete example.

Example 3.3 (Compression Optimization). In Fig. 10, we demonstrate the working of compression optimization on a concrete example. Consider only two analysis-instances in the compiler, namely $a1$ and $a2$. Their `chIndex` values are 1 and 3, respectively. Upon repeated applications of Observation 1, we find that IR-changes A and B both can be safely removed from `chSequence`. Accordingly, both the `chIndex` values are decremented by 2. Further, the IR-changes C and D (and later E and F), can be merged, by repeated applications of Observation 2. Due to the merging of C and D, the `chIndex` of $a2$ is further decremented by 1, to obtain 0.

Impact of compression optimization on space complexity of tracking IR-changes. In the absence of compression, the list `chSequence` would contain as many elements as the number of elementary transformations employed (say k). The size of each element can be $O(s)$, in terms of the size s of the input program (calculated as some function of the number of nodes and edges in the

Type of Developer	Additional Tasks	Efforts	Frequency
Compiler-base writers	- Define elementary transformations - Modify high-level transformation APIs - Create the super-class BasePA	Moderate	Once per compiler
Analysis-pass writers	- Follow the structure defined by BasePA	Minimal	Once per analysis
Optimization-pass writers	<i>None</i>	Nil	-

Fig. 11. Summary of the impact of enabling *Homeostasis*, on different types of compiler writers.

IR). Hence, the total space complexity of an uncompressed *chSequence* can be $O(k \times s)$. In contrast, from Observations 1 and 2, we note that the number of elements in the fixed-point compressed state of *chSequence* cannot be greater than $\min(k, p)$, where p is the number of analysis-instances in the compiler. Since p is usually significantly smaller than k , the worst-case space complexity of *chSequence* with compression optimization, drops to $O(p \times s)$.

Time complexity of the compression algorithm. When the compression algorithm is invoked, the fixed-point state obtained upon repeated applications of Observations 1 and 2 may compress *chSequence* to a single-element, in the worst-case. Hence, if the total number of elementary transformations performed is k , then the compression algorithm may perform at most k merge operations (using function *mergeChanges*, Fig. 8). If the program size is given by s , then the worst-case time complexity of each merge operation will be $O(s)$. Hence, the worst-case time complexity for the compression algorithm is $O(k \times s)$.

3.4 Working in a Homeostasis-enabled compiler

We now describe the additional effort that is needed to enable and use *Homeostasis*, and the impact of using *Homeostasis* on the key questions from Fig. 2. For this discussion, let us assume a hypothetical compiler \mathbb{C} and its *Homeostasis*-enabled counterpart \mathbb{C}_{HS} . We present this discussion with respect to three key kinds of compiler developers: (i) *Compiler-base writers* obtain \mathbb{C}_{HS} from \mathbb{C} by implementing those core components of *Homeostasis* which are common across all the analysis/optimization passes. (ii) *Analysis-pass writers* write analysis passes in \mathbb{C}_{HS} . (iii) *Optimization-pass writers* write optimization passes in \mathbb{C}_{HS} . Fig. 11 summarizes the impact of enabling and using *Homeostasis*, on these key developers.

Impact on compiler-base writer. To obtain \mathbb{C}_{HS} from \mathbb{C} , the compiler-base writer will have to additionally perform the following tasks: (i) identify the elementary transformations and develop them in accordance with the design discussed in Section 3.2.1, (ii) ensure that all high-level transformation APIs in the compiler internally rely on the identified set of elementary transformations to perform any program modifications, and (iii) create a base class for all analysis passes, equivalent to BasePA from Sections 3.2.2 and 3.2.3. Note that these tasks are one-time efforts and involve modifying only some of the IR-related classes.

Impact on analysis-pass writer. Given any analysis class AC in \mathbb{C} , the analysis writer would have to additionally ensure the following, for adapting AC to \mathbb{C}_{HS} : (i) AC is a subtype of BasePA, (ii) the code of AC needed to generate the corresponding program-abstraction is written in the *compute* method, (iii) the corresponding program-abstraction for AC can be accessed only through a getter as defined in Section 3.2.2, and (iv) optionally, if the analysis writer wants to support the LZUPD mode of stabilization, the method *handleUpdate* has to be defined as per Section 3.2.3.

In terms of additional coding effort, ensuring point (i) is trivial; so is ensuring points (ii) and (iv), since the relevant code would be common to both \mathbb{C} and \mathbb{C}_{HS} , and would only possibly differ in their method names. We argue that point (iii) would be automatically ensured in an object-oriented compiler, and hence requires no additional effort. Consequently, we claim that the amount

of additional coding effort required to write an analysis in \mathbb{C}_{HS} is minimal. In terms of the key questions shown in Fig. 2, in a *Homeostasis*-enabled compiler, an analysis-pass writer does not need to address questions \mathbf{Q}_{1-3}^T , while adding any new program analysis in the LZINV mode of stabilization; for the LZUPD mode of stabilization, the analysis writer needs to address *only* \mathbf{Q}_3^T , by providing the definition of the method `handleUpdate`. Note that this is not an additional overhead when using *Homeostasis* – such a code is needed even in the case of manual stabilization in incremental update modes.

Impact on optimization-pass writer. An optimization-pass writer need not write any additional code to adapt an optimization in \mathbb{C} to \mathbb{C}_{HS} . In terms of the key questions shown in Fig. 2, in a *Homeostasis*-enabled compiler, an optimization pass writer need not address the key questions \mathbf{Q}_{1-3} , while adding any new optimization in a *Homeostasis*-enabled compiler. For instance, in Fig. 1, a *Homeostasis*-enabled compiler can generate the expected program P_3 without requiring any changes to the code of the optimization (`DeadBranchRemover`) shown in Fig. 1a. Note that the compiler-base writer would have updated the code for high-level transformation APIs (including the `replace` method), such that all program-changes are expressed directly or indirectly using elementary transformations; a one-time effort.

Overall Summary: Once a *Homeostasis*-enabled compiler has been designed (requires one-time moderate effort on the part of the compiler-base writers), designing and implementing program analyses and optimizations require minimal/zero additional coding effort to maintain stable program-abstractions.

4 FORMAL DESCRIPTION AND CORRECTNESS OF HOMEOSTASIS

We now present a formal description of various components of *Homeostasis*. In Section 4.1, we formally define an abstract *Homeostasis*-enabled compiler, and argue the *correctness guarantee* provided by *Homeostasis*: in a *Homeostasis*-enabled compiler, any compiler pass accessing the program-abstraction of an analysis will only read the stable value of that program-abstraction (see Theorem 4.12). Later, in Section 4.2, we first formally describe how *Homeostasis* maintains the compressed list of program-changes, and then we prove that the compression is semantics-preserving.

4.1 An Abstract *Homeostasis*-enabled Compiler

Consider a *Homeostasis*-enabled compiler for an IR language $\mathbb{L} = \{P_1, P_2, \dots\}$, where each P_i is a valid (syntactically and semantically correct) IR program. Without loss of generality, we assume each program IR to be in the form of a graph. Let the fixed set of elementary transformations in the compiler be represented as $\Gamma = \{\tau_1, \tau_2, \dots, \tau_m\}$, where each transformation τ_i is a map $\mathbb{L} \rightarrow \mathbb{L}$. Note that each elementary transformation updates an IR program by adding/removing a set of nodes and/or edges in the graph to obtain the modified IR program. Fig. 12 summarizes the sets and maps used in the formalism.

Definition 4.1 (Program-change). Given two programs, P_i and $P_j \in \mathbb{L}$, the *program-change* between P_i and P_j (denoted by $\Delta(P_i, P_j)$) is defined as a 4-tuple of the form $\langle \delta_{na}, \delta_{nr}, \delta_{ea}, \delta_{er} \rangle$, where to obtain P_j from P_i , δ_{na} is the minimal set of program nodes to be added to P_i , δ_{nr} is the minimal set of nodes to be removed from P_i , δ_{ea} is the minimal set of edges to be added to P_i , and δ_{er} is the minimal set of edges to be removed from P_i . Specifically, for $P_i = (N_i, E_i)$ and $P_j = (N_j, E_j)$, the following holds: $\delta_{na} = N_j \setminus N_i$, $\delta_{nr} = N_i \setminus N_j$, $\delta_{ea} = E_j \setminus E_i$, and $\delta_{er} = E_i \setminus E_j$. We denote the set of all possible program-changes in \mathbb{L} by \mathcal{C} . Further, we term a sequence of program-changes, of any length, as a *change-sequence*. Without loss of generality, for the ease of exposition, we assume that all such sequences are 0-indexed. \square

Set	Description	Typical Element Symbol
\mathcal{N}	set of all possible IR nodes	n
\mathcal{E}	set of all possible IR edges	e
\mathbb{L}	set of all valid IR programs $\subseteq \mathbb{P}(\mathcal{N}) \times \mathbb{P}(\mathcal{E})$	P_i
Γ	set of all the elementary transformations, where each element is of the form $\mathbb{L} \rightarrow \mathbb{L}$	τ_i
\mathcal{C}	set of all program-changes $\subseteq \mathbb{P}(\mathcal{N}) \times \mathbb{P}(\mathcal{N}) \times \mathbb{P}(\mathcal{E}) \times \mathbb{P}(\mathcal{E})$; if program $P_i = (N_i, E_i)$ and program $P_j = (N_j, E_j)$, then program-change $\Delta(P_i, P_j) = c = (N_j \setminus N_i, N_i \setminus N_j, E_j \setminus E_i, E_i \setminus E_j)$	c
\mathcal{C}^*	set of all change-sequences of elements in \mathcal{C} , of any size	\bar{c}
\mathcal{A}	set of all analysis-instances in the compiler = $\{A_1, A_2, \dots\}$, where each element is of the form $\mathbb{L} \rightarrow R_i$	A_i
R_i	set of all program-abstraction values for analysis-instance A_i	v
\mathcal{R}	set of all the program-abstraction values = $R_1 \cup R_2 \cup \dots$	–
STBST	set of stabilization-status values = {STABLE, UNSTABLE}	st
STBMD	set of lazy modes of stabilization = {LZINV, LZUPD}	m
D_i	set of all analysis-states for the analysis-instance A_i $= R_i \times \text{STBST} \times \text{STBMD} \times \mathbb{N}_{-1}$	d
\mathcal{D}	set of all the analysis-states = $D_1 \cup D_2 \cup \dots$	–
\mathbb{S}	set of compiler states = $\mathbb{L} \times \mathbb{P}(\mathcal{D}) \times \mathcal{C}^*$	S

Fig. 12. Sets and maps used in the formalism. \mathbb{N}_{-1} is the set of natural numbers $\cup \{0, -1\}$. $\mathbb{P}(X)$ denotes the power set of X .

Definition 4.2 (Merge operator). If $\Delta(P_a, P_b) = c_p$ and $\Delta(P_b, P_c) = c_q$, then we use $c_p \oplus c_q$ to provide the net program-change between P_a and P_c . The set \mathcal{C} is closed under the binary operator $\oplus (\mathcal{C} \times \mathcal{C} \rightarrow \mathcal{C})$. This operator corresponds to the function mergeChanges from Fig. 8.

In a compiler written in OO-style, we assume that each analysis is implemented as a class, extending BasePA. During the invocations of the compiler, one or more analysis classes get instantiated (one or more times). Let $\mathcal{A} = \{A_1, A_2, \dots\}$ be the set of all analysis-instances (such as an instance of points-to analysis, an instance of alias analysis, and so on) in the compiler, where each analysis-instance is of the form $A_i : \mathbb{L} \rightarrow R_i$, where R_i denotes the set of meaningful program-abstraction values for A_i . We term $A_i(P)$ as the *stable program-abstraction* of the analysis-instance A_i for program P .

Definition 4.3 (Analysis State). While compiling a program P , at any point during the compilation process, for an analysis-instance A_i , we define its *analysis state* as a 4-tuple of the form $d_i = \langle v, st, m, r \rangle$, where $v \in R_i$ is the value currently maintained by the compiler for the program-abstraction corresponding to A_i , $st \in \text{STBST} = \{\text{STABLE}, \text{UNSTABLE}\}$, denotes the current stabilization status of A_i , $m \in \text{STBMD} = \{\text{LZINV}, \text{LZUPD}\}$, determines the selected lazy mode of stabilization for A_i , and r is either -1 (for LZINV mode of stabilization) or an integer that refers to an index in the global change-sequence (for LZUPD mode of stabilization); r is termed as the *change-index* of A_i . For the LZUPD mode of stabilization, intuitively, the non-negative **change-index** of an analysis-instance A_i denotes that position in the global change-sequence list until and including which the impact of

all elementary transformations have been taken into account, to arrive at the program-abstraction value v for analysis-instance A_i .

Let the set $\mathcal{R} = R_1 \cup R_2 \cup \dots$ be the set of all possible values for all program-abstractions. Further, let \mathcal{D} be the set of all analysis states, for all the analyses-instances. \square

Definition 4.4 (Compiler state). Consider a compilation process, where k elementary transformations have been performed on the input program P . We define a *compiler state* as a 3-tuple of the form $S_i = \langle P_i, D_i, \bar{c}_i \rangle$, where P_i represents the current program (obtained upon applying the k transformations on P), $D_i = \{d_1, d_2, \dots\}$ represents the current set of *analysis states* for different analysis-instances (each a 4-tuple as defined above), and $\bar{c}_i = \langle c_0, c_1, c_2, \dots, c_{k-1} \rangle$, denotes the sequence of k program-changes obtained upon application of each of the k elementary transformations (in order) on the input program P . We denote the set of all compiler states by \mathcal{S} . Note that for an input program P , the *initial* compiler state when compilation starts is $\langle P, \emptyset, \emptyset \rangle$. \square

Now, we state two axioms that are derived from the assumption that the analysis writer has provided the correct implementations of `compute` and `handleUpdate` methods; the latter is required only for the LZUPD mode of stabilization. Axiom 4.5 formally states that the exhaustive algorithm of an analysis-instance has been correctly implemented by the analysis writer in a method named `compute`. Axiom 4.6 formally states that in order to realize the incremental-update mode of stabilization, the analysis writer has provided a correct definition of the `handleUpdate` method, specifying how an abstraction would get updated in response to the given program-changes (between the current program and the one to which the stale program-abstraction corresponds). For the ease of presenting our formal argument, we assume that `compute` and `handleUpdate` methods also return the set of abstractions that were stabilized during the evaluation of these methods. This set is used to only facilitate the formal treatment of our proposed design, and is not required otherwise to perform the actual stabilization.

AXIOM 4.5. *In a Homeostasis-enabled compiler, consider an analysis-instance A_k with LZINV stabilization mode. The analysis writer of A_k provides a correct implementation of the method `compute`, such that when `compute` is invoked at any compilation point, `compute` returns a two-tuple (v, \mathcal{L}) , where $v = A_k(P_i)$, and $\mathcal{L} \subseteq \bar{\mathcal{A}}$, is the set of abstractions which were stabilized during the evaluation of the `compute` method.*

AXIOM 4.6. *Consider an analysis A_k with LZUPD mode of stabilization. Consider any compiler state $S_i = \langle P_i, D_i, \bar{c}_i \rangle$, where the corresponding analysis state for A_k in D_i is denoted by $d_k = \langle v_k, st_k, LZUPD, r_k \rangle$. Further, let, $\bar{c}_i = \langle c_0, c_1, \dots, c_n \rangle$. In a Homeostasis-enabled compiler, the analysis writer of A_k provides the correct implementation of method `handleUpdate`, such that if there exists a program $P_j \in \mathbb{L}$, such that $v_k = A_k(P_j)$ and $\Delta(P_j, P_i) = c_{r_k+1} \oplus c_{r_k+2} \dots \oplus c_n$, then, when `handleUpdate` is invoked on the compiler state S_i , it returns a pair (v, \mathcal{L}) , where (i) $v = A_k(P_i)$, and (ii) $\mathcal{L} \subseteq \bar{\mathcal{A}}$, the set of abstractions which were stabilized during the evaluation of the `handleUpdate` method.*

In Fig. 13, we provide the definition of *stabilization-function*, $\sigma : \mathcal{S} \times \mathcal{D} \rightarrow \mathcal{D} \times \mathbb{P}(\bar{\mathcal{A}})$, which, given a compiler state and a (possibly stale) analysis state for some analysis-instance (in that compiler state), utilizes the definitions of `compute` and `handleUpdate` methods, to provide a two-tuple: (i) the stabilized analysis state of the analysis-instance, and (ii) the set of other analysis-instances on which the stabilization-function was invoked internally during the evaluation of `compute` or `handleUpdate` method. Given such a two-tuple Θ , we use $\Theta.fst$ and $\Theta.snd$ to refer to the first and second elements, respectively, of the tuple.

If the given analysis state is already `STABLE`, the stabilization-function returns the analysis state as it is, along with an empty set of analysis-instances (rule `[STB]`). Otherwise, when the

$$\begin{array}{c}
\frac{S = \langle P, D, \bar{c} \rangle \quad d_k \in D \quad d_k = \langle *, \mathbf{STABLE}, *, * \rangle}{\sigma(S, d_k) = \langle d_k, \emptyset \rangle} \quad [\mathbf{STB}] \\
\\
\frac{S = \langle P, D, \bar{c} \rangle \quad d_k \in D \quad d_k = \langle v, \mathbf{UNSTABLE}, \mathbf{LZINV}, -1 \rangle \quad \mathbf{compute}(A_k, P) = \langle v', \mathcal{L} \rangle \quad d'_k = \langle v', \mathbf{STABLE}, \mathbf{LZINV}, -1 \rangle}{\sigma(S, d_k) = \langle d'_k, \mathcal{L} \rangle} \quad [\mathbf{STB-INV}] \\
\\
\frac{S = \langle P, D, \bar{c} \rangle \quad \bar{c} = \langle c_0, c_1, \dots, c_n \rangle \quad d_k \in D \quad d_k = \langle v, \mathbf{UNSTABLE}, \mathbf{LZUPD}, r \rangle \quad r < n \quad \mathbf{handleUpdate}(A_k, P, v, c_{r+1} \oplus c_{r+2} \oplus \dots \oplus c_n) = \langle v', \mathcal{L} \rangle \quad d'_k = \langle v', \mathbf{STABLE}, \mathbf{LZUPD}, n \rangle}{\sigma(S, d_k) = \langle d'_k, \mathcal{L} \rangle} \quad [\mathbf{STB-UPD}]
\end{array}$$

Fig. 13. Evaluation of the stabilization-function, $\sigma : \mathcal{S} \times \mathcal{D} \rightarrow \mathcal{D} \times \mathbb{P}(\mathcal{A})$, as derived from the template of stabilier from Fig. 7c.

stabilization mode of the analysis state is LZINV, the method `compute` is invoked in order to obtain the correct abstraction value (v') by re-running the corresponding analysis from scratch on the current program (rule [STB-INV]). In contrast, as per rule [STB-UPD], when the stabilization mode is LZUPD, the method `handleUpdate` is invoked to obtain the correct abstraction value (v'), given the following arguments: (i) a reference to the analysis-instance (A_k), (ii) the current program, (iii) the unstable/stale abstraction value (v), and (iv) the net program-changes obtained by using the merge operator on the appropriate elements from the change-sequence. Note that the change-index for the returned analysis state is set to n (index of the last element in the global change-sequence).

Definition 4.7 (Compiler action). Recall that *Homeostasis* allows changes to the IR program only through elementary transformations. Similarly, the program-abstraction of an analysis can be accessed only through its appropriate getters. Accordingly, we define three kinds of *compiler actions* that can be issued by an optimization or analysis pass, and may alter the current compiler state, $S = \langle P, D, \bar{c} \rangle$.

- (i) Given an analysis class \mathbf{AC} , the compiler action $\mathbf{construct}(\mathbf{AC}, m)$ creates an instance of \mathbf{AC} , invokes the `compute` method to populate the program-abstraction, and sets the stabilization mode of the constructed analysis-instance to $m \in \mathbf{STBMD}$.
- (ii) The compiler action $\mathbf{get}(A_k)$ returns the value $A_k(P)$.
- (iii) The compiler action $\mathbf{transform}(\tau_i)$ transforms the current program P to $\tau_i(P)$.

We represent the set of all possible compiler actions with \mathcal{W} . □

Note that in object-oriented compilers, the `construct` compiler action for an analysis will always be invoked before the first invocation of the corresponding `get` compiler action.

Definition 4.8 (apply function). We define a compilation step using the function $\mathbf{apply} : \mathcal{S} \times \mathcal{W} \Longrightarrow \mathcal{S} \times (\mathcal{R} \cup \{\epsilon\})$, which applies a given compiler action w on the compiler state s , to generate (i) a possibly-modified compiler state s' , and (ii) a program-abstraction value or ϵ .

For a *Homeostasis*-enabled compiler, the definition for function `apply` is shown in Fig. 14a, Fig. 14b, and Fig. 14c, for the design templates of various components shown in Fig. 7a (constructor), Fig. 7b-7c (getters and stabilizers), and Fig. 5 (elementary transformations), respectively.

- When a `construct` compiler action is issued for an analysis-instance A , an initial analysis state for A is generated and added to the set of analysis states in the current compiler state (rule [C-APPLY]). Note that in the case of LZUPD mode, the change-index of A is set to the index of the last element in the global change-sequence; it is set to -1 if the list is empty.

$$\begin{array}{c}
S = \langle P, D, \bar{c} \rangle \quad A_i = \text{new AC}() \\
\text{compute}(A_i, P) = (v, \mathcal{L}) \quad \mathcal{L} = \{A_1, A_2, \dots, A_n\} \quad A_i \notin \mathcal{L} \\
D_a = D - \{d_1, d_2, \dots, d_n\} \quad D_b = \{\sigma(S, d_1).\text{fst}, \sigma(S, d_2).\text{fst}, \dots, \sigma(S, d_n).\text{fst}\} \\
d_i = \langle v, \text{STABLE}, m, ((m = \text{LZUPD})?(|\bar{c}| - 1) : -1) \rangle \\
\hline
\text{apply}(S, \text{construct}(\text{AC}, m)) \Rightarrow \langle \langle P, D_a \cup D_b \cup \{d_i\}, \bar{c} \rangle, \epsilon \rangle \quad [\text{C-APPLY}]
\end{array}$$

(a) Evaluation of apply on the “construct” compiler action, as derived from Fig. 7a. The helper method newInstanceOf gives a fresh uninitialized analysis-instance for the given analysis.

$$\begin{array}{c}
S = \langle P, D, \bar{c} \rangle \quad d \in D \\
\sigma(S, d_k) = \langle d'_k, \mathcal{L} \rangle \quad d'_k = \{v', \text{STABLE}, *, *\} \quad \mathcal{L} = \{A_1, A_2, \dots, A_n\} \quad A_k \notin \mathcal{L} \\
D_a = D - \{d_1, d_2, \dots, d_n\} - \{d_x\} \quad D_b = \{\sigma(S, d_1).\text{fst}, \sigma(S, d_2).\text{fst}, \dots, \sigma(S, d_n).\text{fst}\} \\
\hline
\text{apply}(S, \text{get}(A_k)) \Rightarrow \langle \langle P, D_a \cup D_b \cup \{d'_k\}, \bar{c} \rangle, v' \rangle \quad [\text{G-APPLY}]
\end{array}$$

(b) Evaluation of apply on the “get” compiler action, as derived from the template of getters from Fig. 7b.

$$\begin{array}{c}
S = \langle P, D, \bar{c} \rangle \quad P' = \tau_l(P) \quad \bar{c}' = \bar{c} \cdot \langle \Delta(P, P') \rangle \\
D = \{\langle v_1, s_1, m_1, r_1 \rangle, \langle v_2, s_2, m_2, r_2 \rangle, \dots, \langle v_x, s_x, m_x, r_x \rangle\} \\
D' = \{\langle v_1, \text{UNSTABLE}, m_1, r_1 \rangle, \langle v_2, \text{UNSTABLE}, m_2, r_2 \rangle, \dots, \langle v_x, \text{UNSTABLE}, m_x, r_x \rangle\} \\
\hline
\text{apply}(S, \text{transform}(\tau_l)) \Rightarrow \langle \langle P', D', \bar{c}' \rangle, \epsilon \rangle \quad [\text{T-APPLY}]
\end{array}$$

(c) Evaluation of apply on the “transform” compiler action, as derived from the template of elementary transformations given in Fig. 5.

Fig. 14. Definition for the apply function, derived from the design of *Homeostasis* explained in Section 3.2.

- When a get compiler action is issued for A_k , the design of *Homeostasis* ensures that the analysis state of A_k is changed as per the mapping rules shown in Fig. 14b and 13. Using Axioms 4.5 and 4.6, rules [STB-INV] and [STB-UPD] set the stabilization status of A_k to STABLE, as can be inferred from the design of an abstraction getter, and stabilizer.
- When a transform compiler action with elementary transformation τ_l is issued on a compiler state $S = \langle P, D, \bar{c} \rangle$, then as per rule [T-APPLY] in Fig. 5, the design of *Homeostasis* changes the compiler state to a tuple where, (i) the current program P is modified to $\tau_l(P)$, (ii) the stabilization status of all analyses is set to UNSTABLE, and (iii) $\Delta(P, \tau_l(P))$ is added to the sequence of program-changes.

Definition 4.9 (Valid compiler state). We term a compiler state of the form $S_i = \langle P_i, D_i = \{d_1, d_2, \dots\}, \bar{c}_i = \langle c_0, c_1, \dots, c_n \rangle \rangle$ as a *valid compiler state*, if and only if the following condition, termed as **stability-condition**, holds for the state $d_j = \langle v_j, st_j, m_j, r_j \rangle \in D_i$ of each analysis-instance A_j :

- if $st_j = \text{STABLE}$, then $r_j = n$ and $v_j = A_j(P_i)$, else
- if $st_j = \text{UNSTABLE}$, then $n \geq 0$, $r_j < n$, and $\exists P_o \in \mathbb{L}$, such that $v_j = A_j(P_o)$ and $\Delta(P_o, P_i) = c_{r_j+1} \oplus c_{r_j+2} \dots \oplus c_n$.

Intuitively, a compiler state is considered *valid* when each program-abstraction is either stable, or has access to the complete information that can be used to make it stable. \square

The following Lemma gives the correctness guarantee for the value returned by a get compiler action.

LEMMA 4.10. *If a compiler state $S_i = \langle P_i, D_i, \bar{c}_i \rangle$ is valid, then for each analysis-instance A_k with corresponding analysis state $d_k \in D_i$, the value obtained upon invocation of get compiler action for A_k (i.e., $\text{get}(A_k)$) on S_i , will always be same as $A_k(P_i)$ (i.e., the stable program-abstraction value). Formally, for each analysis A_k and valid compiler state S_i , $\exists S_j \in \mathbb{S}$ such that $\text{apply}(S_i, \text{get}(A_k)) \implies (S_j, A_k(P_i))$.*

PROOF. Let $\bar{c}_i = \langle c_0, c_1, \dots, c_U \rangle$, and let $D_i = \{d_1, d_2, \dots, d_k, \dots\}$, where $d_k = \langle v_k, st_k, m_k, r_k \rangle$ be the analysis state of A_k in S_i . Assume, $\text{apply}(S_i, \text{get}(A_k)) \implies (S_j, v'_k)$. Hence, it is sufficient to prove that $v'_k = A_k(P_i)$. As the compiler state S_i is valid, the analysis state d_k satisfies the stability-condition from Definition 4.9. Two cases arise based on the value of st_k :

CASE A: $st_k = \text{STABLE}$. From the stability-condition of d_k , we obtain $v_k = A_k(P_i)$. Hence, from rule [STB], we derive $v'_k = v_k = A_k(P_i)$. Hence proved.

CASE B: $st_k = \text{UNSTABLE}$. Two cases arise based on the value of m_k :

Case B.1: $m_k = \text{LZINV}$. From rule [STB-INV], we obtain $(v'_k, \mathcal{L}) = \text{compute}(A_k, P_i)$. From Axiom 4.5, $v'_k = \text{compute}(A_k, P_i).\text{fst} = A_k(P_i)$. Hence proved.

Case B.2: $m_k = \text{LZUPD}$. From rule [STB-UPD], we have $(v'_k, \mathcal{L}) = \text{handleUpdate}(A_k, P_i, v_k, c_{r_k+1} \oplus c_{r_k+2} \dots \oplus c_U)$. From the stability-condition of d_k , we derive the following: $U \geq 0, r_k < U$, and that there exists a program P_j such that $v_k = A_k(P_j)$ and $\Delta(P_j, P_i) = c_{r_k+1} \oplus c_{r_k+2} \dots \oplus c_U$. Using Axiom 4.6, we obtain $v'_k = A_k(P_i)$. Hence proved. \square

The following Lemma gives the correctness guarantee that the invocation of any compiler action (construct, get, or transform) on a valid compiler state will result in a valid compiler state.

LEMMA 4.11. *For every sequence $W = \langle w_1, w_2, \dots \rangle$ of compiler actions, the compiler state, starting with the initial compiler state $\langle P, \phi, \phi \rangle$, remains valid after application of each action from the sequence.*

PROOF. (Proof by natural induction on the number of elements in W .)

Base Case: $|W| = 0$. For an input program P , the initial compiler state is always $\langle P, \emptyset, \emptyset \rangle$, which is trivially valid.

Inductive hypothesis: The compiler state, say $S_i = \langle P_i, D_i, \bar{c}_i \rangle$, obtained after application of all compiler actions from the string W_m , of length m , is valid. Let $D_i = \{d_1, d_2, \dots, d_k = \langle v_k, st_k, m_k, r_i \rangle, \dots\}$, and $\bar{c}_i = \langle c_0, c_1, \dots, c_U \rangle$.

Inductive step: Consider a string of compiler actions, W_{m+1} , of length $m+1$, obtained by appending a compiler action w_{m+1} to W_m . We now need to prove that if $\text{apply}(S, w_{m+1}) = (S', v)$, then S' is a valid compiler state. Three cases occur for possible form of compiler action w_{m+1} :

CASE A: w_{m+1} is $\text{construct}(AC, m_s)$ for some analysis class AC. Say the created analysis-instance is A_s and the initial analysis state for A_s is $d_s = \langle v_s, st_s, m_s, r_s \rangle$. As per rule [C-APPLY], $S' = \langle P_i, D_i \cup \{d_s = \langle A_s(P_i), \text{STABLE}, m_s, U \rangle\}, \bar{c}_i \rangle$. S' is trivially valid as d_s satisfies the stability-condition from Definition 4.9. Hence proved.

CASE B: w_{m+1} is $\text{get}(A_x)$ for some $d_x \in D_i$. As S is a valid compiler state, $\forall d_j \in D_i$ satisfies the stability-condition. Let $\text{get}(A_x)$ change a subset of analysis states $D_y \subseteq D_i$. Let $d_y (\in D_y)$ be given by $\langle v_y, st_y, m_y, r_y \rangle$. Assume, for each d_y , the changed analysis state d'_y is given by $\langle v'_x, st'_x, m_x, r'_x \rangle$. To prove that S' is a valid compiler state, it is sufficient to prove that each such d'_y satisfies the stability-condition. Let us consider any such d'_y . Two sub-cases arise based on the values of st_y :

Case B.1: $st_y = \text{STABLE}$. As d_y satisfies the stability-condition, we derive the following from rule [STB]: $st'_y = st_y = \text{STABLE}$, $r'_y = r_y = U$, and $v'_y = v_y = A_x(P_i)$. Hence, d'_y satisfies the stability-condition. Hence proved.

Case B.2: $st_y = \text{UNSTABLE}$. From rules [STB-INV] and [STB-UPD], we derive $st'_y = \text{STABLE}$. Similarly, we derive $r'_y = -1$ or $r'_y = U$, from rules [STB-INV] and [STB-UPD], respectively. Depending upon value of m_y , two sub-cases arise:

Case B.2.i: $m_y = \text{LZINV}$. From rule [STB-INV], $v'_y = \text{compute}(A_y, P_i).\text{fst}$. From Axiom 4.5, $v'_y = A_y(P_i)$. Hence, d'_y satisfies the stability-condition, and S' is a valid compiler state.

Case B.2.ii: $m_y = \text{LZUPD}$. From rule [STB-UPD], $v'_y = \text{handleUpdate}(A_y, P_i, v_y, c_{r_y+1} \oplus c_{r_y+2} \dots \oplus c_U).\text{fst}$. From Axiom 4.6, $v'_y = A_y(P_i)$, if there exists a program $P_k \in \mathbb{L}$, such that (i) $v_y = A_y(P_k)$, and (ii) the net program-change (i.e., last argument of `handleUpdate`) is same as $\Delta(P_k, P_i)$. Since d_y satisfies the stability-condition, such a program P_k exists. (Note that the $\Delta(P_k, P_i)$ is the same as the last argument of `handleUpdate`.) Hence, $v'_y = A_y(P_i)$, and S' is a valid compiler state.

CASE C: w_{m+1} is `transform(τ_t)` for some $\tau_t \in \Gamma$. Let $S' = \langle P'_i, D'_i, \overline{c'_i} \rangle$. From rule [T-APPLY], we have: $P'_i = \tau_t(P_i)$, and $\overline{c'_i} = \langle c_0, c_1, \dots, c_U, \Delta(P_i, \tau_t(P_i)) \rangle$. Further, for each $d \in D_i$, there exists a modified analysis state $d' \in D'_i$. To prove that S' is a valid compiler state, we need to prove that each analysis state $d' \in D'_i$ satisfies the stability-condition. Consider any analysis-state $d_x \in D$, and its corresponding modified analysis-state $d'_x \in D'_i$. Let $d_x = \langle v_x, st_x, m_x, r_x \rangle$, and let $d'_x = \langle v'_x, st'_x, m'_x, r'_x \rangle$. Now, two sub-cases arise based on the value of st_x :

Case C.1: $st_x = \text{STABLE}$. Since d_x satisfies the stability-condition, we have: $r_x = U$ (if $m_x = \text{LZUPD}$) or $r_x = -1$ (if $m_x = \text{LZINV}$), and $v_x = A_x(P_i)$. From rule T-APPLY, we have: $v'_x = v_x$, $m'_x = m_x$, $r'_x = r_x$, and $st'_x = \text{UNSTABLE}$. From Definition 4.9, d'_x satisfies the stability-condition, and S' is a valid program state.

Case C.2: $st_x = \text{UNSTABLE}$. Since d_x satisfies the stability-condition, we know that: $U \geq 0$, $r_x < U$, and that a program, say P_j , exists such that $v_x = A_x(P_j)$ and $\Delta(P_j, P_i) = c_{r_x+1} \oplus c_{r_x+2} \dots \oplus c_U$. From rule T-APPLY, we have: $v'_x = v_x$, $m'_x = m'_x$, $r'_x = r_x$, and $st'_x = \text{UNSTABLE}$. Hence, $r'_x < U+1$. Further from the definition of \oplus , $\Delta(P_j, P'_i) = \Delta(P_j, P_i) \oplus \Delta(P_i, P'_i)$. Hence, $\Delta(P_j, P'_i) = \Delta(P_j, P_i) \oplus c_{U+1} = c_{r_x+1} \oplus c_{r_x+2} \dots \oplus c_U \oplus c_{U+1}$. By Definition 4.9, S' is a valid compiler state. \square

Now, we first state the correctness theorem for *Homeostasis*. Intuitively, Theorem 4.12 states that at any point of time during compilation, the analysis results given by *Homeostasis* match those obtained by rerunning the complete analyses, provided the methods `compute` and `handleUpdate` provided by the analysis writer are correct (i.e., Axioms 4.5 and 4.6 hold).

THEOREM 4.12 (CORRECTNESS OF HOMEOSTASIS). *At any compilation point of a Homeostasis-enabled compiler with compiler state $S_i = \langle P_i, D_i, \overline{c_i} \rangle$, for each analysis A_k with corresponding analysis state $d_k \in D_i$, the value obtained upon invocation of `get compiler action` for A_k (i.e., `get(A_k)`) on S_i , will always be same as $A_k(P_i)$. Formally, for each analysis A_k and compiler state S_i , $\exists S_j \in \mathbb{S}$, such that `apply(S_i , get(A_k)) $\implies (S_j, A_k(P_i))$.`*

PROOF. *Follows directly from Lemma 4.10 and Lemma 4.11.* \square

4.2 Compression of Change-Sequence in Homeostasis

Homeostasis keeps the global change-sequence compact by internally performing the compression operation on it, as discussed in Section 3.3. In this section, we formalize this functionality, and argue that the correctness of *Homeostasis* continues to hold when we use the proposed idea of change-sequence compression.

$$\begin{array}{c}
\frac{|\bar{c}| = 0}{\mathbf{compress}((D, \bar{c})) = (D, \bar{c})} \quad [\text{CS-BASE}] \qquad \frac{\begin{array}{l} |\bar{c}| = k \quad k \geq 1 \\ \bigvee_{i=-1}^{k-2}, \exists d \in D, d = \langle *, *, *, i \rangle \end{array}}{\mathbf{compress}((D, \bar{c})) = (D, \bar{c})} \quad [\text{CS-COMPRESSED}] \\
\\
\frac{\begin{array}{l} D = D_{\text{INV}} \cup D_{\text{UPD}} \quad \forall d \in D_{\text{INV}}, d = \langle *, *, \text{LZINV}, -1 \rangle \quad \forall d \in D_{\text{UPD}}, d = \langle *, *, \text{LZUPD}, * \rangle \\ \bar{c} = c_1 \cdot \bar{c}' \quad \nexists d \in D_{\text{UPD}}, d = \langle *, *, \text{LZUPD}, -1 \rangle \\ D_{\text{UPD}} = \{ \langle v_1, s_1, \text{LZUPD}, r_1 \rangle, \langle v_2, s_2, \text{LZUPD}, r_2 \rangle, \dots, \langle v_k, s_k, \text{LZUPD}, r_k \rangle \} \\ D'_{\text{UPD}} = \{ \langle v_1, s_1, \text{LZUPD}, r_1-1 \rangle, \langle v_2, s_2, \text{LZUPD}, r_2-1 \rangle, \dots, \langle v_k, s_k, \text{LZUPD}, r_k-1 \rangle \} \end{array}}{\mathbf{compress}((D, \bar{c})) = (D_{\text{INV}} \cup D'_{\text{UPD}}, \bar{c}')} \quad [\text{CS-DELFIRST}] \\
\\
\frac{\begin{array}{l} D = D_{\text{INV}} \cup D_{\text{UPD}} \quad \forall d \in D_{\text{INV}}, d = \langle *, *, \text{LZINV}, -1 \rangle \quad \forall d \in D_{\text{UPD}}, d = \langle *, *, \text{LZUPD}, * \rangle \\ \bar{c} = \bar{c}_a \cdot c_i \cdot \bar{c}_b \quad |\bar{c}_a| = i \quad i \geq 1 \quad |\bar{c}_b| \geq 1 \\ \bigvee_{j=-1}^{i-1}, \exists d \in D_{\text{UPD}}, d = \langle *, *, \text{LZUPD}, j \rangle \quad \nexists d \in D_{\text{UPD}}, d = \langle *, *, \text{LZUPD}, i \rangle \\ \bar{c}_b = \langle c_{i+1}, c_{i+2}, \dots, c_n \rangle \quad \bar{c}'_b = \langle c_i \oplus c_{i+1}, c_{i+2}, \dots, c_n \rangle \quad \bar{c}' = \bar{c}_a \cdot \bar{c}'_b \\ D_{\text{UPD}} = \{ \langle v_1, s_1, \text{LZUPD}, r_1 \rangle, \langle v_2, s_2, \text{LZUPD}, r_2 \rangle, \dots, \langle v_k, s_k, \text{LZUPD}, r_k \rangle \} \\ D'_{\text{UPD}} = \{ \langle v_1, s_1, \text{LZUPD}, ((r_1 > i)?(r_1-1):r_1) \rangle, \langle v_2, s_2, \text{LZUPD}, ((r_2 > i)?(r_2-1):r_2) \rangle, \\ \dots, \langle v_k, s_k, \text{LZUPD}, ((r_k > i)?(r_k-1):r_k) \rangle \} \end{array}}{\mathbf{compress}((D, \bar{c})) = (D_{\text{INV}} \cup D'_{\text{UPD}}, \bar{c}')} \quad [\text{CS-MERGE MID}]
\end{array}$$

Fig. 15. Definition for the function $\mathbf{compress}: \mathbb{P}(\mathcal{D}) \times C^* \rightarrow \mathbb{P}(\mathcal{D}) \times C^*$.

Fig. 15 presents the formal definition of the $\mathbf{compress}$ function. This function takes a two-tuple of the form (D, \bar{c}) , where D is a set of analysis-states, and \bar{c} is a change-sequence. The $\mathbf{compress}$ function returns another two-tuple, say (D', \bar{c}') , such that (i) the change-sequence \bar{c}' is obtained by optionally compressing a single element of the input change-sequence \bar{c} , and (ii) the set D' contains all the analysis-states from D , with required modifications in case if the input change-sequence was compressed.

Recall that the key intuition behind compression is as follows: If no analysis-state in the compiler has a change-index equal to some index in the change-sequence, then the element at that index in the change-sequence can be safely merged (using the \oplus operator) to its successor (if any) in the change-sequence. Further, if no analysis-state in the compiler has a change-index less than zero, then the first element (at index 0) of change-sequence can be safely removed. Following are the cases that may arise when performing compression of the input change-sequence one element at a time:

- When the change-sequence \bar{c} is empty, it is already in its compressed form. Hence, as per rule [CS-BASE], D and \bar{c} are returned without any changes.
- Similarly, if corresponding to each element in the change-sequence (except, optionally, the last) if there exists at least one analysis-state that has its change-index same as the index of that element, and if there exists at least one analysis-state that has its change-index as -1 , then the change-sequence is already in its compressed form ([CS-COMPRESSED]).
- Otherwise, if no analysis-state has a change-index of -1 , then we purge the first element from the change-sequence, and update all the analysis-states by decrementing their change-index by one (rule [CS-DELFIRST]).

• Otherwise, the element at the least index number that is not present as the change-index of any analysis-state is merged to its successor element (if any) in the change-sequence. Further, the change-index of all those analysis-states whose change-index is greater than the compressed index number, is decremented by one (rule [CS-MERGMID]).

Note that no action is performed during compression when only the last element in the change-index is not indexed to by any of the analysis-states; this case falls under the rule [CS-COMPRESSED], once the fixed-point application of the compress is complete. Further note that for efficiency purposes, in practice, in a change-sequence, all those k consecutive elements that are not indexed by any change-index are compressed together and change-indices of the affected analysis-instances are decremented by k , in one go.

In *Homeostasis*, the compress function can be called during any compilation point, such as, immediately after a stabilization trigger, after an elementary transformation, and so on. Regardless of the number and places of invocations for this function, the following lemma holds.

LEMMA 4.13. *Consider a valid compiler state, say $S = \langle P, D, \bar{c} \rangle$, and another state obtained after a single application of the compress function, say $S' = \langle P, D', \bar{c}' \rangle$, such that $\text{compress}((D, \bar{c})) = (D', \bar{c}')$. Assume, $\bar{c} = \{c_0, c_1, \dots, c_{|\bar{c}|-1}\}$, and $\bar{c}' = \{c'_0, c'_1, \dots, c'_{|\bar{c}'|-1}\}$. For each analysis-state in LZUPD mode, say $d = \langle v, s, \text{LZUPD}, r \rangle \in D$, and the corresponding modified analysis-state, say $d' = \langle v, s, \text{LZUPD}, r' \rangle \in D'$, the following equality holds.*

$$c_{r+1} \oplus c_{r+2} \dots \oplus c_{|\bar{c}|-1} = c'_{r'+1} \oplus c'_{r'+2} \dots \oplus c'_{|\bar{c}'|-1}. \quad (1)$$

PROOF. Let k be the smallest integer in the set $\{-1, 0, 1, \dots, |\bar{c}| - 1\}$ such that there does not exist any analysis-state in D whose change-index is same as k . That is, $\nexists d = \langle *, *, *, r \rangle \in D$, such that $r = k$. Following three cases arise:

CASE I: *No such k exists.* This case refers to the fixed-point state for the compress function, as described by Rules [CS-BASE] and [CS-COMPRESSED]. As per these rules, note that $\bar{c}' = \bar{c}$, and $D' = D$. Hence, Equation 1 trivially holds.

CASE II: $k = -1$. Following two sub-cases arise based on the size of \bar{c} .

Case II.a.: $|\bar{c}| = 0$. In this case, Rule [CS-BASE] is applicable. Since $\bar{c}' = \bar{c}$, and $D' = D$, Equation 1 trivially holds.

Case II.b.: $|\bar{c}| \geq 1$. In this case, Rule [CS-DELFIRST] is applicable. As per this rule, $|\bar{c}'| = |\bar{c}| - 1$,

$$r' = r - 1 \quad \text{and} \quad \bigvee_{i=0}^{|\bar{c}'|-1} c'_i = c_{i+1}$$

Under these equalities we infer that, (i) the number of terms on the LHS and RHS of Equation 1 is same, and (ii) each term on the LHS is same as the corresponding term on the RHS. Hence, Equation 1 holds.

CASE III: $k \geq 0$. When $k = |\bar{c}| - 1$, Rule [CS-COMPRESSED] is applicable. As before, Equation 1 will hold in that case. Otherwise, when $0 \leq k \neq |\bar{c}| - 1$, Rule [CS-MERGMID] is applicable. In this case, $|\bar{c}'| = |\bar{c}| - 1$. Further, as per this rule, the following holds:

$$\bigvee_{i=0}^{k-1} c'_i = c_i, \quad c'_k = c_k \oplus c_{k+1} \quad \text{and} \quad \bigvee_{i=k+1}^{|\bar{c}'|-1} c'_i = c_{i+1} \quad (2)$$

Following two sub-cases arise based on the different values of r :

Case III.a. $r < k$. As per Rule [CS-MERGMID], $r' = r$. In this case, the number of elements on the LHS of Equation 1 is one more than the number of elements of its RHS. Clearly, c_k and c_{k+1} are present in the LHS, whereas in the RHS, c_k and c_{k+1} have been replaced with another element c'_k . Using Equation 2, we derive that all terms on the left of c_k are same as

the corresponding terms on the left of c'_k . Similarly, all terms on the right of c_{k+1} are same as the corresponding terms on the right of c'_k . Additionally, since $c'_k = c_k \oplus c_{k+1}$, we infer that Equation 1 holds.

Case III.b. $r > k$. As per Rule [CS-MERGE-MID], $r' = r - 1$. In this case, the number of elements on the LHS of Equation 1 is same as the number of elements of its RHS. Clearly, c_k does not appear in the LHS. As per Equation 2, each term on the LHS is same as the corresponding term on the RHS. Hence Equation 1 holds. \square

COROLLARY 4.14. *Consider a valid compiler state, say $S = \langle P, D, \bar{c} \rangle$. If the fixed-point application of the compress function on the input (D, \bar{c}) is (D', \bar{c}') , that is, $\text{compress}((D, \bar{c}))^* = (D', \bar{c}')$, then the compiler state $S' = \langle P, D', \bar{c}' \rangle$ is valid.*

PROOF. The proof follows directly from Definition 4.9 and Lemma 4.13. \square

Note that as a direct implication of Corollary 4.14, Lemma 4.10, and Lemma 4.11, the correctness guarantee of *Homeostasis*, stated in Theorem 4.12, will hold regardless of the number of compilation points where the compress function is invoked by *Homeostasis*.

5 DISCUSSION

We now present some salient characteristics of *Homeostasis*, followed by an account of how we estimate the amount of stabilization-effort required in the absence of *Homeostasis*. Then, we briefly discuss a generic incremental iterative data-flow analysis pass, which we have implemented and instantiated in *Homeostasis* to aid our evaluation.

5.1 Salient Characteristics of *Homeostasis*

Merging program changes across transformations. To stabilize a program-abstraction in update (UPD) mode, the `handleUpdate` method of the corresponding analysis requires *net* changes performed on the program (in terms of added and removed nodes/edges) since the last stabilization of that program-abstraction. A simple union of all such changes performed across multiple transformations, would not suffice. For instance, if a program node n (or edge e), is added by a transformation, and then deleted by another transformation before the stabilization gets triggered, the node n (or edge e) should not be considered as an added/removed node (or edge). In *Homeostasis*, BasePA internally merges the changes saved across all transformations to provide the relevant information when requested. In Section 5.5, we give an example of how the provided information is used in the `handleUpdate` method of a specific analysis.

Resolving pass-dependencies. *Homeostasis* automatically resolves the pass-dependencies of all compiler passes on program analysis passes by internally stabilizing the program-abstractions in the topological order of their dependencies. This is achieved through the design of getters and stabilizers in *Homeostasis* (see Fig. 7). For example, the getter method of an UNSTABLE analysis, say A , does not return until its stabilizer has been run. If analysis A is dependent upon some other analysis B (that is, if the computation of the program-abstraction of A requires the program-abstraction of B), then the `compute` or `handleUpdate` methods of A invoked from the stabilizer of A would, in turn, invoke the getter of B . This will ensure that the stabilizer of B is run before the completion of A , regardless of the order in which the stabilizations of both analyses were triggered.

5.2 Using *Homeostasis* in Compilers with Multi-Level IRs

A majority of the real-world compilers employ multiple levels of IR, and invoke a host of IR-level-specific analyses and optimizations on the program at each IR level. For such compilers,

we can easily see that *Homeostasis* can be used for stabilizing the program-abstractions of the analyses within each such IR level, in response to the optimizations performed at that level. During compilation when the program is lowered from a higher-level IR (say, HIR) to a lower-level IR (say, LIR), (i) the program-abstractions of HIR are no longer relevant for the compiler passes of LIR (and hence need not be stabilized), and (ii) all the necessary program-abstractions for LIR are computed on the current state of the program, when needed. Thus, program-abstractions need not be stabilized *across* different levels of IR. Consequently, compilers supporting multi-level IRs can realize self-stabilization by using *Homeostasis* at each level of the IR independently.

5.3 Using *Homeostasis* in Non-Object-Oriented Compilers

While this paper presents the design of *Homeostasis* in the context of object-oriented compilers, we believe that the proposed design can be adapted to allow self-stabilization in compilers that are written even in procedural languages. For example, in languages like C, the information about each analysis pass can be represented as a structure containing six members: (i) a program-abstraction of type “void*”, (ii) an enum field `stabilizationMode`, (iii) an integer field `chIndex`, (iv) a function pointer named `compute` pointing to the function that populates the corresponding program-abstraction, and (v) a function pointer named `handleUpdate` pointing to the function that incrementally stabilizes the corresponding program-abstraction, (vi) a function pointer named `get` pointing to a method that returns a stable program-abstraction (by invoking the corresponding `compute` or `handleUpdate` functions, as needed). The first five members correspond to the instance fields and methods of the class `BasePA`, and the last member corresponds to the `get` method shown in Fig. 7. All instances of this six-member structure are maintained in a global list. We have to identify the set of elementary transformations and modify them to conform to the code shown in Fig. 5.

Since procedural languages do not support the Observer pattern naturally, we would need that the compiler-writer follows a disciplined approach in modifying the IR, and accessing the program-abstractions. To use *Homeostasis* in compilers written in such a language, (i) the compiler writer must perform all IR modifications (directly/indirectly) only via the fixed set of elementary transformations, (ii) perform the initialization of the analysis instance via a generic method (akin to the constructor of `BasePA`, Fig. 7a), and (iii) access the program-abstraction of an analysis, only via the invocation of its `get` function. We believe that a similar exercise can be carried out for compilers written in functional programming languages. We leave it as a future work to actually instantiate *Homeostasis* in such non-OO compilers.

5.4 Estimating Manual Stabilization Effort in the Absence of *Homeostasis*

To contrast the savings in effort due to the self-stabilization scheme of *Homeostasis*, we now present a scheme to estimate the effort required for manual stabilization in the absence of *Homeostasis*.

Critical program-abstractions, and change-points. Note that in order to perform manual stabilization in the context of an optimization \mathcal{O} , the pass writer needs to inspect the following two parameters in relation to questions \mathbf{Q}_1 and \mathbf{Q}_2 (see Fig. 2): (i) *critical program-abstractions*: those program-abstractions which may be rendered stale by \mathcal{O} , and may be read later by an existing or future optimizing pass (and hence may need to be stabilized); we use S_a to denote the set of critical program-abstractions, and (ii) *change-points*: those program points in the code of the optimization \mathcal{O} which may directly or indirectly trigger an elementary transformation (and hence may necessitate stabilization); we use S_p to denote the set of change-points. We demonstrate these parameters using an example.

```

public void unrollLoop (WhileStatement loop) {
    DumpSnapshot.dumpPointsTo(loop, "original"); // L1
    Statement newBody = ... /* code to obtain the unrolled body */
    loop.setBody(newBody); // L2
    CompoundStatementNormalizer.removeExtraScopes(loop); // L3
    DumpSnapshot.dumpPointsTo(loop, "afterUnroll"); // L4
}

```

Fig. 16. Concrete optimization pass: Loop unroller from IMOP (WhileStatementInfo.java:60-83).

Example 5.1. Fig. 16 shows a snippet of a concrete optimization pass that performs loop unrolling in the IMOP compiler framework. We have added the calls to the `dumpPointsTo` method to simply print the points-to information corresponding to the specified node, symbolising the clients that may need points-to information, before and after unrolling. Method `setBody` is an elementary transformation method that replaces the current body of the receiver while-statement with the provided body. Method `removeExtraScopes` is used to remove unnecessary wrapping of statements within blocks (for example, $\{\{S\}\} \rightarrow \{S\}$). For this snippet, $S_p = \{L2, L3\}$, and S_a is the set of those program-abstractions that may be impacted by the statements in S_p ; for example, CFG, call-graph, use-def chains, and so on. Note that there may be other change-points within the implementation of the common utility `removeExtraScopes`; we do not consider them here as they are not present in the code written as a part of this pass.

Relevant change-Points (RP). An inexperienced programmer may perform manual stabilization naïvely by re-computing each of the program-abstractions (or a likely superset of S_a , for correctness), after each change-point in S_p – a highly inefficient scheme, both in terms of its effect on performance, and the number of program points where the stabilization code needs to be invoked. An experienced programmer, on the other hand, is more likely to insert stabilization code only at a subset of change-points present in the code of the optimization pass. This reduction stems from two key insights: (i) Stabilization is required only after the last change-point in a series of transformations, after which one or more critical program-abstractions may be read. (ii) Further, a program-abstraction need not be stabilized at such a last change-point, if the net program-changes at that change-point is guaranteed to not alter the existing values of the program-abstraction. For example, consider the addition/deletion of a node for an assignment statement (of the form $x=y$). If the variable x is a pointer-type variable, then such an addition/deletion does not change the points-to map of the variable x , if both x and y point to the same set of ‘may points-to’ locations, even in the absence of this statement. Similarly, if the variable x is of integer type, then the addition/deletion of such a node will not impact the points-to information of any variable (if the expert can assert that the value of the variable x is not used as an offset for any pointer dereference).

The second key insight leads to cases where even if the program-abstraction is marked as UNSTABLE, in reality no stabilization is required. Considering the challenges in obtaining an expert-based study in a large compiler framework, and the difficulty in automatically identifying the cases related to the second key insight, for the purpose of our evaluation, we approximate the set of expert-marked change-points by the set of *relevant change-points* that is based on the first key insight. A relevant change-point is a program point in the optimization pass that corresponds to the last change-point in a series of transformations, after which one or more critical program-abstractions may be read; hence, one or more program-abstractions may require manual stabilization at a relevant change-point. For example, in Fig. 16, the set of relevant change-points, from among the change-points $\{L2, L3\}$, is a singleton set $\{L3\}$ – L3 is the last change-point in the series of transformations (L2;L3) after which a program-abstraction, the points-to map, will be read (at L4).

Modes of manual stabilization. Besides the lazy modes of stabilization preferred by *Homeostasis*, a few other modes of stabilization are easily conceivable. For example, the stabilization of all

	Full recomputation of the analysis from scratch	Incremental update of the program-abstraction using the stored changes to the program
Stabilization triggers eagerly at each elementary transformation	EGINV	EGUPD
Stabilization triggers at <i>relevant change-points</i> in the optimization pass	RPINV	RPUPD
Stabilization triggers lazily upon an attempt to read from the stale program-abstractions	LZINV	LZUPD

Fig. 17. Summary of various modes of stabilization used in the evaluation of our implementation of *Homeostasis* in IMOP.

```

1 Function handleUpdate()
2   Set seeds = .. empty set ..;
3   IRChanges netIRChanges = getNetChanges(this.chIndex);
4   foreach  $e$  in netIRChanges.removedEdges do Add the destination node of the removed edge  $e$  to seeds;
5   foreach  $n$  in netIRChanges.removedNodes do Add the (old) successors of the removed node  $n$  to seeds;
6   foreach  $n$  in netIRChanges.addedNodes do Add the added node  $n$  itself, as well as its successors to seeds;
7   foreach  $e$  in netIRChanges.addedEdges do Add the destination node of the added edge  $e$  to seeds;
8   incrementalIDFA(seeds);

```

Fig. 18. Definition of the handleUpdate function for forward iterative data-flow analysis.

program-abstractions can be triggered during each elementary transformation of the program eagerly (see Section 3.1.4). Like the two modes of lazy stabilization, this eager scheme leads to two modes of stabilization: eager-invalidate (EGINV) and (EGUPD). An efficient alternative to the eager modes of stabilization is to invoke stabilization at only the set of relevant change-points. Like before, two modes of stabilization are possible – RPinV and RPUPD, corresponding to the invalidate and incremental update options, respectively.

These stabilization modes are very similar to the custom codes manually written by the compiler pass writers in conventional compilers: (i) UPD modes in the presence of incremental update, whereas INV modes otherwise, and (ii) EG-modes during naïve (and inefficient) but easy manual stabilization, whereas RP-modes during relatively efficient stabilizations performed by an experienced compiler writer. For the purpose of evaluation (see Section 7), along with the two lazy-modes of self-stabilization advocated by *Homeostasis*, we have implemented these four modes of stabilization (EGINV, EGUPD, RPinV, and RPUPD) to approximate different modes of manual stabilization. Fig. 17 summarizes all the implemented modes of stabilization.

5.5 Implementing Iterative Data-Flow Analyses in *Homeostasis*

To conveniently implement multiple analyses that can use LZUPD mode of stabilization, we implemented a generic pass of an analysis that can be used to implement any incremental, inter-thread, flow-sensitive, iterative data-flow analysis (IDFA) for OpenMP parallel programs. We term this generic analysis as HIDFA and implemented it as a class that extends BasePA.

Fig. 18 shows the steps used by the `handleUpdate` method of the `HIDFA` class in the context of forward data-flow analyses; the steps in the context of backward analysis are similarly derived (not shown). To realize self-stabilization, it populates a set (seeds) of nodes, starting which the flow maps may need an update as a result of program transformations. These seed nodes are passed to the `incrementalIDFA` method for performing the stabilization (Line 8). The implementation of `incrementalIDFA` is inspired from the work by Marlowe and Ryder [1989]; Ryder et al. [1988], and is a simple extension of the two-phase incremental update algorithms for iterative versions of data-flow analysis given by Pollock and Soffa [1989], adapted to handle the parallel semantics of the OpenMP parallel programs.

For completeness, we now give brief details of our implementation of `incrementalIDFA`. This method starts with the reconstruction of the SCC Decomposition Graph (SDG) of the updated control-flow graph of the program. Each SCC-node in the SDG that contains any of the seed nodes is *marked* for processing. All the marked SCC-nodes are processed one-by-one, as per their topological-sort order in the SDG. Within each SCC-node, the processing continues using a worklist-based approach in two-phases: (i) *exaggerate* phase: in this phase, the flow-maps of various program-nodes in the SCC-node are initialized to an under-approximated value (such as \top), starting which the precise flow-map can be eventually obtained, and (ii) *adjust* phase: in this phase, the program-nodes are processed until fixed-point, using the data-flow equations similar to the exhaustive iterative data-flow algorithm. At the end of the second phase, the flow-maps of all program-nodes within the SCC-node correspond to the maximum fixed-point (MFP) solution. During the processing of an SCC-node, if it is found that the program-nodes of any successor SCC-nodes need to be added to the worklist, then such successor SCC-nodes are marked for processing. This method continues till there are no more marked SCC-nodes left for processing. We refer the reader to the work by Pollock and Soffa [1989] for more details on the underlying two-phase incremental update algorithm.

6 INSTANTIATION OF HOMEOSTASIS

In order to assess the feasibility and generality of *Homeostasis*, we have retrofitted *Homeostasis* to a research compiler, and have implemented/adapted a number of analysis passes and optimization passes in the compiler, to conform to the design of *Homeostasis*. In this section, we discuss the relevant instantiation details.

6.1 Retrofitting *Homeostasis* to a Real-World Compiler

The design of *Homeostasis* is applicable to all the object-oriented compilers. Retrofitting *Homeostasis* in mainstream real-world compilers, like LLVM, OpenJ9, IMOP, and so on, will require rewriting all the existing program transformations using the identified fixed set of elementary transformations, and modifying all the existing analysis passes to conform to the structure prescribed by *Homeostasis* (see Fig. 11). We have implemented our prototype in IMOP. We believe that a similar exercise can be undertaken for other object-oriented compiler frameworks as well. To further strengthen this belief, in Appendix B we briefly discuss the guidelines on how the LLVM compiler framework can be extended with *Homeostasis*. The one-time effort required to retrofit *Homeostasis* to IMOP is discussed next.

Identifying elementary transformations. The IMOP IR comprises two kinds of nodes – *leaf* nodes and *non-leaf* nodes. All OpenMP or C constructs with “body”, such as while-loop, if-statement, and OpenMP parallel-region, are represented as *non-leaf* nodes. Similarly, nodes such as simple statements of C, and OpenMP directives such as barriers and flushes, are considered as *leaf* nodes. Each non-leaf node defines a nested flow of control, using edges among its components, which can themselves be non-leaf nodes. We identified the pre-existing fundamental transformations in

IMOP that are used to add, replace, or delete various components of a non-leaf node, as the set of elementary transformations. In total, we identified and modified 47 elementary transformations spread across 17 non-leaf nodes to make them conform to the design shown in Fig. 5.

We observe that the modification of these elementary transformations did not necessitate any changes at their call-sites. From Fig. 5, note that Step B, which corresponds to the pre-existing code of an elementary transformation (in the absence of *Homeostasis*) does not undergo any changes while retrofitting the compiler with *Homeostasis* – the IR transformations are performed as before. Steps A, C, and D, interact only with stabilization-specific data structures to capture the program-changes resulting from Step B; they do not alter the meaning of the elementary transformation. Consequently, the compiler-pass writers can continue to use the elementary transformation as before, without having to consider their impact on self-stabilization.

Modifying higher-level transformation APIs. In order to allow for easier expression of transformation passes, IMOP provides four basic CFG transformations: (i) insertion of a successor node, (ii) insertion of a predecessor node, (iii) insertion of a node on an edge, and (iv) removal of a node. To retrofit *Homeostasis* to IMOP, we have modified the implementation of all these four transformations such that they express all IR changes directly or indirectly in terms of the identified set of elementary transformations. IMOP provides many higher-level transformation APIs such as duplicating a block of statements, renaming of free variables, and so on. All these transformation APIs employ elementary transformations, the above-mentioned CFG transformations, or other higher-level transformations to express any changes to the IR. Since all the transformations APIs deployed in IMOP directly or indirectly express all the program changes via elementary transformations, the implementation of none of the existing optimization passes in IMOP required any specific modifications to adapt them to use *Homeostasis*.

Creating/identifying the super-class BasePA. IMOP already contains a class named BasePA, which serves as the super-class of all the analysis passes. We modified BasePA as per Fig. 7 to conform to the design of *Homeostasis*. This mainly involved (i) modifying the constructor as shown in Fig. 7a, (ii) adding the stabilize method from Fig. 7c, (iii) adding an abstract method compute to be implemented by each concrete subclass of BasePA, and (iv) adding a dummy method for handleUpdate to be overridden by each concrete subclass of BasePA that seeks to employ LZUPD stabilization-mode. Further, to allow efficient maintenance of the centralized list of program-changes, as detailed in Fig. 15, we added the appropriate data structures and methods.

6.2 Adapting Analysis Passes to Use *Homeostasis*

In order to assess the applicability of *Homeostasis* to various kinds of program-abstractions, we have (i) extended existing analysis passes of IMOP to conform to the design of *Homeostasis*, and (ii) implemented HIDFA and its various instantiations in IMOP.

Fig. 19 provides a brief description of the program-abstractions for some of the key analysis passes in IMOP that now use *Homeostasis*. In particular, we have implemented a set of six standard flow-sensitive context-insensitive inter-thread iterative data-flow analyses as subclasses of the HIDFA class: (i) points-to analysis, (ii) reaching-definitions analysis, (iii) liveness analysis (a backward IDFA), (iv) copy propagation analysis, (v) lockset analysis, and (vi) dominance analysis. None of these IDFA passes required any stabilization-specific code in order to support LZUPD stabilization-mode. For their super-class HIDFA, as well as for other analyses listed in the figure, we had to (1) modify their corresponding getter methods, as per the design from Fig. 7, (2) ensure that the code to generate the program-abstraction from scratch is invoked from the overridden method compute, and (3) wherever applicable, ensure that the code to perform an incremental update of the program-abstraction is invoked from the overridden method handleUpdate.

Program-abstraction (Relevant analysis-pass class)	Description
1. Points-to maps (PointsToAnalysis*)	points-to information
2. Reaching-definitions (ReachingDefinitionAnalysis*)	reaching definitions
3. Liveness maps (LivenessAnalysis*)	liveness information of variables
4. Copy maps (CopyPropagationAnalysis*)	information about variables that are copies of each other
5. Must-lock sets (LockSetAnalysis*)	information about locks that must have been taken at a node
6. Dominator sets (DominanceAnalysis*)	dominator information
7. Control-flow graphs (CFGInfo)	control-flow graph
8. Inter-task edges (InterTaskEdge)	information representing inter-task communication
9. Strongly-connected components (SCC)	SCC Decomposition Graph (SDG) and enclosing SCC-nodes
10. Phase-flow graphs (NodePhaseInfo)	may-happen-in-parallel (MHP) analysis
11. Call graphs (CallStatementInfo)	call graph
12. Symbol-/type-tables (CompoundStatementInfo)	information about the symbols and user-defined types in a scope
13. Label-lookup sets (StatementInfo)	information about labels of a statement
14. Single-valued expressions (SVEChecker)	information about expressions that evaluate to same value by all the threads

Fig. 19. List of program-abstractions corresponding to key analysis passes in IMOP that were made conforming to *Homeostasis*. The bracketed text in the first column, gives the name of the IMOP class modified. Classes annotated with * are the analyses that are implemented as subclasses of HIDFA in IMOP.

6.3 Optimization Passes in the Context of *Homeostasis*

Recall that in a *Homeostasis*-enabled compiler framework, the optimization-pass writers can efficiently design and implement new optimizations without having to write *any* stabilization-specific code (see Fig. 11). Fig. 20 lists some key optimization passes (including existing, as well as new ones implemented by us) in IMOP; none of them required any stabilization-specific code. Two common characteristics of these passes are: (i) Each of them expresses the required program transformations using either the elementary transformations and/or other transformation APIs provided by IMOP. (ii) Each of them uses/impacts one or more *Homeostasis*-enabled program-abstractions described in Fig. 19. Now, we briefly describe a subset of these optimization passes (that we implemented for evaluating *Homeostasis*), which we collectively term as **BarrElim**.

BarrElim: a set of optimization passes to aid barrier elimination. In order to develop a real-world client for evaluating *Homeostasis*, we have implemented a set of three optimization passes that aim to reduce redundant barriers in OpenMP C programs: redundant-barrier-remover, parallel-construct-expander, and selective-function-inliner. The parallel-construct-expander in turn uses five pre-existing optimizations: (i) privatisation of shared variables, (ii) parallel-region scope expander, (iii) merging of consecutive parallel-regions, (iv) parallel-region and enclosing serial-loop interchange, and (v) parallel-region unswitching. Further, the selective-function-inliner uses two pre-existing cleanup-optimizations: extra-scope-remover, and unused-elements-remover. Collectively, we term the three implemented optimization passes as **BarrElim**.

The optimization passes in **BarrElim** extend prior work by [Gupta and Schonberg \[1996\]](#); [Tseng \[1995\]](#) on parallel-region expansion and barrier removal with function inlining to realize an efficient

Optimization Pass (Relevant class/method in IMOP)	Description
1. Redundant barrier remover [†] (RedundantSynchronizationRemoval)	removes those barriers that do not preserve any inter-thread dependences across them
2. OmpPar expander [†] (ParallelConstructExpander.expandParRegion)	expands the scope of parallel-regions upwards and downwards
3. OmpPar merger [†] (ParallelConstructExpander.mergeParRegions)	merges two consecutive parallel-regions, if safe
4. OmpPar-loop interchange [†] (ParallelConstructExpander.interchangeUp)	interchanges a parallel-region with its enclosing loop, if safe
5. OmpPar unswitching [†] (ParallelConstructExpander.interchangeUp)	interchanges a parallel-region with its enclosing if-statement, if safe
6. Variable privatization [†] (ParallelConstructExpander.expandUpward)	privatizes OpenMP shared variables, if safe
7. Function inliner [†] (FunctionInliner)	selectively inlines those monomorphic calls that contain at least one barrier
8. Scope remover [†] (CompoundStatementNormalizer)	removes redundant encapsulations of compound-statements within the given node
9. Unused-elements remover [†] (NodeInfo::removeUnusedElements)	removes unused functions, types, typedefs, and symbol declarations
10. Expression simplification (ExpressionSimplifier)	simplifies complex C expressions as per IMOP's normalization rules
11. OmpConstruct simplification (ExpressionSimplifier)	simplifies combined OpenMP constructs as per IMOP's normalization rules
12. Loop unroller (IterationStatementInfo::unrollLoop)	unrolls a serial loop
13. Inter-phase code percolator (BarrierDirectiveInfo::percolateCodeUpwards)	moves instructions across an OpenMP barrier, if safe
14. Copy eliminator (CopyElimination)	removes redundant variables, relying on copy-propagation analysis
15. Loop barrier extractor (LoopInstRescheduler::peelFirstBarrier)	peels out a barrier from within a serial loop, if safe

Fig. 20. List of key optimization passes inspected in IMOP; none of these passes required any stabilization-specific code in the context of *Homeostasis*. The bracketed text in the first column, gives the name of the IMOP class modified. Passes annotated with a [†] indicate that they belong to, or are used by, the `BarrierElim` set of optimization passes.

barrier removal algorithm. These optimization passes perform the following steps, as shown in the block diagram in Fig. 21 : (1) Remove redundant barriers (within a parallel-region), whose removal do not violate any data dependence among the statements across them. The remaining two passes help improve the opportunities for barrier removal within each function. (2) Expand and merge the parallel-regions, while possibly expanding their scope to the call-sites of their enclosing functions, wherever possible. This helps in bringing more barriers (including the implicit ones) within the resulting parallel-region, thereby creating new opportunities for barrier removal. (3) Inline those monomorphic calls whose target function (i) is not recursive, and (ii) contains at least one barrier. These three optimization passes are repeated until fixed-point (no change).

Like many similar optimizations [Aloor and Nandivada 2015; Barik et al. 2013; Gupta et al. 2017; Nandivada et al. 2013], the optimization passes in `BarrierElim` involve multiple alternating phases of inspections and transformations, which in turn lead to a number of interleaved accesses (reads

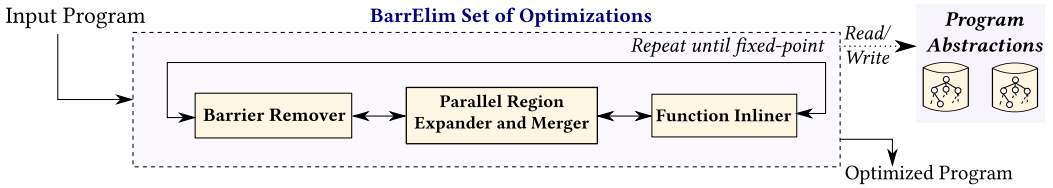


Fig. 21. Block diagram of BarrElim, the selected set of client optimization passes, for removal of barriers from OpenMP programs.

and writes) to various program-abstractions, such as phase information, points-to information, super-graph (involving CFGs, call-graphs, and inter-task edges), AST, and so on – this interaction is depicted in Fig. 21 using the dotted edge.

We have found that BarrElim is an impactful set of optimization passes for OpenMP codes. For example, on the NPB benchmarks [Van der Wijngaart and Wong 2002], the BarrElim optimized code yielded up to 5% improvement in execution time; the input benchmarks, as well as the BarrElim optimized codes were compiled using the `-O3` switch of gcc. Considering that the obtained gains are on top of the many optimizations enabled by the `-O3` switch, it can be seen that the gains are significant. Consequently, from among the various optimization passes present in IMOP, we select the BarrElim set of optimization passes (involving nine optimization passes and numerous program-abstractions) to evaluate the ease-of-use and performance of *Homeostasis* (see Section 7).

7 IMPLEMENTATION AND EVALUATION

We have implemented all the key components of *Homeostasis* (see Fig. 4, and Section 3.2) in the IMOP compiler framework [Nougrahiya and Nandivada 2019]. IMOP is a new open-source compiler framework for conveniently writing program analyses, source-to-source optimizations, and profiling tools for OpenMP C programs; it spans more than 170k lines of code (LOC) in Java. IMOP has been used successfully in various published works (such as [Viswakaran Sreelatha and Balachandran 2016], [Viswakaran Sreelatha et al. 2018], [Viswakaran Sreelatha and Nare 2018] and [Krishnakumar et al. 2019]), and is under active development. Our implementation of *Homeostasis* in IMOP involved coding of ~4500 LOC for adding self-stabilization to the elementary-transformation methods, and ~2500 LOC for supporting self-stabilization for existing program-abstractions of IMOP.

We have also implemented HIDFA, our generic inter-thread IDFA pass with incremental update in *Homeostasis* (~3300 LOC; discussed in Section 5.5). We instantiated HIDFA to realize the implementation of following six standard flow-sensitive context-insensitive inter-thread iterative data-flow analyses: (i) points-to analysis (~870 LOC), (ii) reaching-definitions analysis (~190 LOC), (iii) liveness analysis (~140 LOC), (iv) copy-propagation analysis (~200 LOC), (v) lock-set analysis (~240 LOC), and (vi) dominance analysis (~116 LOC). Further, in order to assess the usability of *Homeostasis*, we have implemented a set of optimization passes, collectively termed as BarrElim, for barrier removal (~2500 LOC; see Section 6.3). As expected, in the set of optimization passes in BarrElim no stabilization-specific code was needed. This underscores the ease-of-use facilitated by the design of *Homeostasis*. We have made our implementation of *Homeostasis* in IMOP open-source and publicly available [Nougrahiya and Nandivada 2021].

7.1 Experimental Setup

We present our evaluation on a set of twenty-four real-world benchmark programs taken from four popular benchmark suites (listed in Fig. 22): (i) all eight benchmark programs of NPB-OMP 3.0 suite [Van der Wijngaart and Wong 2002], (ii) quake, and art-m, the two (out of 3) OpenMP-C

1		2	3	4	5	6
Benchmark		#LOC	#Nodes	#Edges	#PC	#Barr
1.	BT (NPB)	2615	4748	5016	9	47
2.	CG (NPB)	642	1403	1485	14	31
3.	EP (NPB)	352	775	813	2	4
4.	FT (NPB)	899	2033	2151	7	14
5.	IS (NPB)	333	711	762	2	4
6.	LU (NPB)	2355	4687	4974	8	35
7.	MG (NPB)	1278	2784	2918	10	19
8.	SP (NPB)	2543	5364	5744	7	72
9.	quake (SPEC)	1489	3333	3491	11	22
10.	art-m (SPEC)	1691	1710	1791	4	4
11.	amgmk (Sequoia)	895	1867	1949	2	5
12.	clomp (Sequoia)	1605	4162	4289	28	73
13.	stream (Sequoia)	331	735	762	6	12
14.	bellman-ford (IMSuite)	143	546	564	3	6
15.	bfs-dijkstra (IMSuite)	104	295	305	2	4
16.	byzantine (IMSuite)	135	444	467	3	6
17.	dominating-set (IMSuite)	321	1223	1277	12	24
18.	kcommittee (IMSuite)	220	889	925	6	12
19.	leader-elect-dp (IMSuite)	122	442	460	2	4
20.	leader-elect-hs (IMSuite)	184	528	553	3	6
21.	leader-elect-lcr (IMSuite)	89	301	314	3	6
22.	mis (IMSuite)	161	549	574	4	8
23.	mst (IMSuite)	299	1226	1279	9	18
24.	vertex-coloring (IMSuite)	172	609	636	3	6

Fig. 22. Benchmark characteristics. Abbreviations: #LOC=number of lines of code, #Node=number of executable nodes in the CFG, #Edge=number of edges in the CFG, #PC=number of static parallel constructs, and #Barr=number of static barriers (implicit + explicit).

benchmark programs from SPEC OMP 2001 [Aslot et al. 2001] that can be handled by IMOP, (iii) all three OpenMP C benchmark programs – amgmk, clomp, and stream – from Sequoia benchmark suite [Seager, M 2008], and (iv) all the eleven benchmark programs from IMSuite benchmark suite [Gupta and Nandivada 2015]. Note that these comprise some of the largest standard open-source benchmark programs for OpenMP C. Since IMOP accepts only OpenMP C programs, we did not consider any other benchmarks from SPEC OMP 2001, or from Sequoia, as they contain a mix of C/C++/MPI code. For IS from NPB, the set of client optimization passes, BarrElim, could not find any opportunity for optimization; as no program transformation was performed by the pass, no stabilization of data-flow analyses was triggered in any mode of self-stabilization – hence, we omit IS from our discussion. In Fig. 22, we list a few static characteristics of the benchmarks, such as the size of each selected benchmark, the number of parallel constructs, and the number of static barriers.

In order to study the impact of different hardware on the performance improvements obtained with *Homeostasis*, we used two multicore platforms to perform our empirical evaluation: (i) **Nanda**, a 64-thread (dual socket, 16 cores per socket, 2 threads per core) 2.3 GHz Intel Xeon Gold 5218 system with 64 GB of memory, running Ubuntu 22.04.1 LTS and Java HotSpot 64-Bit Server VM 14.0; and (ii) **K2**, a 64-thread (quad-socket, 8 cores per socket, 2 threads per core) 2.3 GHz AMD

Component Pass of BarrElim	LOC	#CP	Program-Abs.	Mode	#RP
Parallel-construct expansion	1675	66	Points-to graphs	UPD	17
Function inlining	463	15	Control-flow graphs	UPD	70
Redundant-barrier deletion	313	2	Call graphs	UPD	60
Driver	10	1	Phase information	INV	61
Total :	2461	84	Inter-task edges	INV	31
			Symbol-/type-tables	UPD	37
			Label-lookup tables	UPD	60

(a) Maximum number of change-points obtained within the component optimization passes of BarrElim upon running them on the benchmarks under study. Abbr.: *LOC*=number of lines of code; *#CP*=number of change-points.

(b) Program-abstractions read by the BarrElim passes. Abbr.: *Mode*=stabilization-mode; *#RP*=number of relevant change-points in the BarrElim passes after which the abstraction was read from.

Fig. 23. Stabilization-related information for the set of optimization passes constituting BarrElim, obtained upon profiling/inspecting the code.

Abu Dhabi system with 512 GB of memory, running CentOS 6.4 and Java HotSpot 64-Bit Server VM 9.0. The multi-core aspect of these systems is relevant for the part of the evaluation that executes the generated OpenMP code. The OpenMP codes were compiled using the GCC 11.3.0 compiler. Taking inspiration from the insightful paper of Georges et al. [2007], we report the compilation- and execution-time numbers by taking a geometric mean over 30 runs. We present the evaluation in four directions: (i) Ease-of-use of *Homeostasis* (in Section 7.2), (ii) Performance evaluation of the proposed lazy modes of stabilization (in Section 7.3), (iii) Impact of using the compression optimization on the data-structures employed by *Homeostasis* to keep track of the program-changes (in Section 7.4), and (iv) Correctness of the lazy modes of stabilization (in Section 7.5).

7.2 Ease of Use: Self-stabilization vs. Manual Stabilization

We now present an empirical study for assessing the impact of *Homeostasis* on writing different compiler passes, by comparing the coding effort required to realize self-stabilization in a *Homeostasis*-enabled compiler, against those required to perform manual stabilization (in the absence of *Homeostasis*). We perform this evaluation in the context of the various component optimization passes of BarrElim.

Based on the discussion in Section 5.4, we designed a simple scheme to estimate the additional coding effort that may be required to perform manual stabilization. We profiled the IMOP compiler by instrumenting the implementations of the various optimization passes of BarrElim, as well as of various program analyses and elementary transformations. By running this profiled compiler on each benchmark program, we obtained (i) the set of change-points for the BarrElim passes, and (ii) the set of program-abstractions that may be impacted by the BarrElim passes (see Section 6.3). In this section, we use this data to estimate the manual coding effort that may be required to answer the key questions from Fig. 2.

Which program-abstractions to stabilize? As discussed in Section 5.4, since it is difficult to obtain the exact set of *critical program-abstractions* (to be stabilized manually), we manually analyzed the code of various optimization passes of BarrElim and estimated that there are seven program-abstractions (listed in Fig. 23b) that may be impacted by the BarrElim passes. Thus, in the case of manual-stabilization, on writing each of the BarrElim passes, the pass writer needs to identify these seven program-abstractions, from the plethora of available program-abstractions – a daunting task. Note that the BarrElim optimization passes may also impact other program-abstractions that may be added to the compiler in future; hence, for each newly added analysis-pass, the analysis-pass writer would have to manually assess whether the corresponding program-abstraction may be

rendered stale due to the transformations performed by any of the `BarrElim` optimization passes (and, in fact, by any of the other optimization passes in the compiler).

In contrast, in a *Homeostasis*-enabled compiler, all these tasks are automated – the optimization-pass (or even analysis-pass) writer needs to put no effort into identifying the program-abstractions (existing or new) that may be impacted by an optimization pass.

Where to invoke stabilization? In Fig. 23a, we enumerate the number of change-points discovered in the major components of `BarrElim`. In the absence of *Homeostasis*, the pass writer would have to correctly identify these 84 change-points (i.e., on average, *almost 1 for every 28 lines of code!*) across various optimization passes of `BarrElim`, and insert code for ensuring stabilization of the affected program-abstractions. At each change-point, the pass writer may need to handle stabilization of the impacted program-abstractions, irrespective of the chosen mode of stabilization. To identify a more aggressive baseline, we note that not all change-points (referred to in Figure 23a) may warrant stabilization, and a pass writer may need to invoke the stabilization code only at the relevant change-points (see Section 5.4), for only those set of abstractions that may be modified at that point.

Fig. 23b, column 4, lists the number of *relevant* change-points for each program-abstraction impacted by the optimization passes of `BarrElim`; this data too was obtained by profiling the IMOP compiler (profiling details discussed above), while running it on all the benchmarks under study. The figure shows that there are a significant number of places in the components of `BarrElim` where this stabilization code needs to be manually invoked, in the absence of *Homeostasis*. For example, CFG stabilization needs to be performed at 70 places, and call-graphs at 60 places – which can lead to cumbersome and error-prone code. Further, upon addition of any new program analysis to the compiler (or any modification to the existing analysis), the pass writer would have to revisit all the change-points of pre-existing optimizations (for example, 84 change-points across all optimization passes of `BarrElim`) to check if the change-point may have necessitated stabilization of the newly added/modified program-abstraction.

In contrast, in a *Homeostasis*-enabled compiler, all the above tasks are automated – the pass writer needs to spend no effort in identifying the places of stabilization, as she needs to add no additional code as part of the optimization in order to stabilize the program-abstractions.

How to stabilize? We use the same stabilization code for both manual stabilization and *Homeostasis*, and hence there is no difference between them, with respect to this key question.

Summary. In contrast to the traditional compilers, it is much easier to ensure stabilization of program-abstractions when writing optimizations or analyses in *Homeostasis*.

7.3 Performance Evaluation

We conduct the performance evaluation of the proposed lazy modes of stabilization, by studying the parameters related to stabilization time, total compilation time, and memory consumption, in the context of the prior discussed set of optimization passes, `BarrElim`. We do so by presenting a comparison to the RP-modes of stabilization, which correspond to manual stabilizations performed by experienced compiler-pass writers (see Section 5.4). In Fig. 24, we present various parameters related to compilation of the benchmark programs under study, when compiled with the `RPINV` mode of stabilization; these numbers serve as the baseline for the evaluations discussed in this section.

We have also done an elaborate evaluation compared to the weaker baseline of the eager EG-modes of stabilization, which correspond to naïve (that is, easy but inefficient) manual stabilizations; the details can be found in Appendix A.

(A) Stabilization time. In Fig. 24, columns 4 and 5 show the stabilization time for IDFA flow-maps in the context of `RPINV` mode of stabilization, while performing the set of optimization passes of

Benchmark	2	3	4	5	6	7	8	9	10
	#Trig	#T-Func	STB-Time (s)		Total Time (s)		Memory (MB)		Ch-Size
			Nanda	K2	Nanda	K2	Nanda	K2	
1. BT (NPB)	17	214327	3.62	5.93	8.29	14.25	2862.24	8530.41	15170
2. CG (NPB)	89	242019	2.14	3.46	3.65	6.11	1675.41	3850.55	115070
3. EP (NPB)	2	1983	0.04	0.07	0.56	1.17	421.32	1784.43	0
4. FT (NPB)	61	443820	4.42	7.16	7.81	13.03	2771.97	10060.87	49871
5. IS (NPB)	0	806	0	0	0.33	0.82	330.51	1430.9	0
6. LU (NPB)	33	507444	8.43	14.63	11.92	21.13	2856.21	8894.1	41361
7. MG (NPB)	230	2372204	24.95	58.17	34.47	74.12	2788.71	13969.44	571691
8. SP (NPB)	35	672958	8.31	14.36	11.66	20.43	2828.68	16217.09	38102
9. quake (SPEC)	35	437217	6.74	10.44	8.92	14.29	3220.78	13348.05	16727
10. art-m (SPEC)	5	88347	1.03	2.01	2.33	4.56	610.15	3030.15	134
11. amgmk (Sequoia)	56	239940	3.2	6.54	6.15	11.36	2715.4	7760.35	52486
12. clomp (Sequoia)	63	397078	6.76	11.56	13.72	23.19	2779.42	14921.1	146511
13. stream (Sequoia)	45	51020	0.55	1.03	1.4	2.79	548.6	2154.01	36426
14. bellman-ford (IMS)	23	15038	0.3	0.49	0.86	1.62	488.84	2008.73	6051
15. bfs-dijkstra (IMS)	17	4695	0.13	0.29	0.73	1.49	480.23	1965.33	3237
16. byzantine (IMS)	19	7979	0.19	0.32	0.59	1.25	405.9	1768.1	4563
17. dominating-set (IMS)	75	186360	1.88	3.09	3.19	5.32	1255.84	4298.49	64687
18. kcommittee (IMS)	43	44272	0.56	1.01	1.57	2.93	535.33	2147.7	30767
19. leader-elect-dp (IMS)	5	2325	0.07	0.12	0.39	0.84	340.87	1407.41	282
20. leader-elect-hs (IMS)	11	9978	0.18	0.35	0.84	1.67	500.83	1977.11	2624
21. leader-elect-lcr (IMS)	6	1844	0.05	0.1	0.52	1.11	407.03	1698.05	402
22. mis (IMS)	7	6033	0.13	0.24	0.51	1.08	380.34	1650.04	540
23. mst (IMS)	53	217334	2	3.21	3.28	5.57	899.43	3184.65	48387
24. vertex-coloring (IMS)	13	13370	0.26	0.46	0.91	1.82	504.2	2017.47	1954

Fig. 24. Evaluation numbers for performing the BarrElim optimization passes on various benchmark programs, in the context of RPINV mode of stabilization. These numbers serve as the baseline for Section 7.3. Abbreviations: *#Trig*=number of stabilization triggers, *#T-Func*=number of transfer-function applications, and *Ch-Size*=total size of the change-sequence across all stabilization triggers (as a sum of the number of nodes and twice the number of edges in each trigger). *STB-Time* and *Total Time* refer to the IDFA-stabilization time, and overall compilation time, respectively. *Memory* refers to the maximum resident set size of the compilation process when running the BarrElim optimization passes. Note that *Ch-Size* is calculated through a special run of the RPINV mode, where we force retention of the change-sequence, even though the mode does not require it.

BarrElim on the benchmark programs under study, for Nanda and K2, respectively. The zero (0) entries for the STB-times of IS are due to the fact that no optimization opportunities were found by the BarrElim passes for IS.

In Fig. 25, we illustrate the impact of using LZINV, RPUPD, and LZUPD modes of stabilization, by showing their relative speedups with respect to RPINV, in terms of speedups in the IDFA stabilization-time. We exclude SP from this discussion on stabilization time, since there were no stabilization triggers for SP under lazy modes of stabilizations (that is, the stabilization time for lazy modes of stabilization was zero). As expected, the LZUPD mode incurs the least cost for stabilization among all the cases across both the platforms; consequently, it results in the maximum speedup with respect to RPINV – with speedups varying between 2.6× and 30.64× (geomean 8.84×) for Nanda; between 3.22× and 29.68× (geomean 9.58×) for K2.

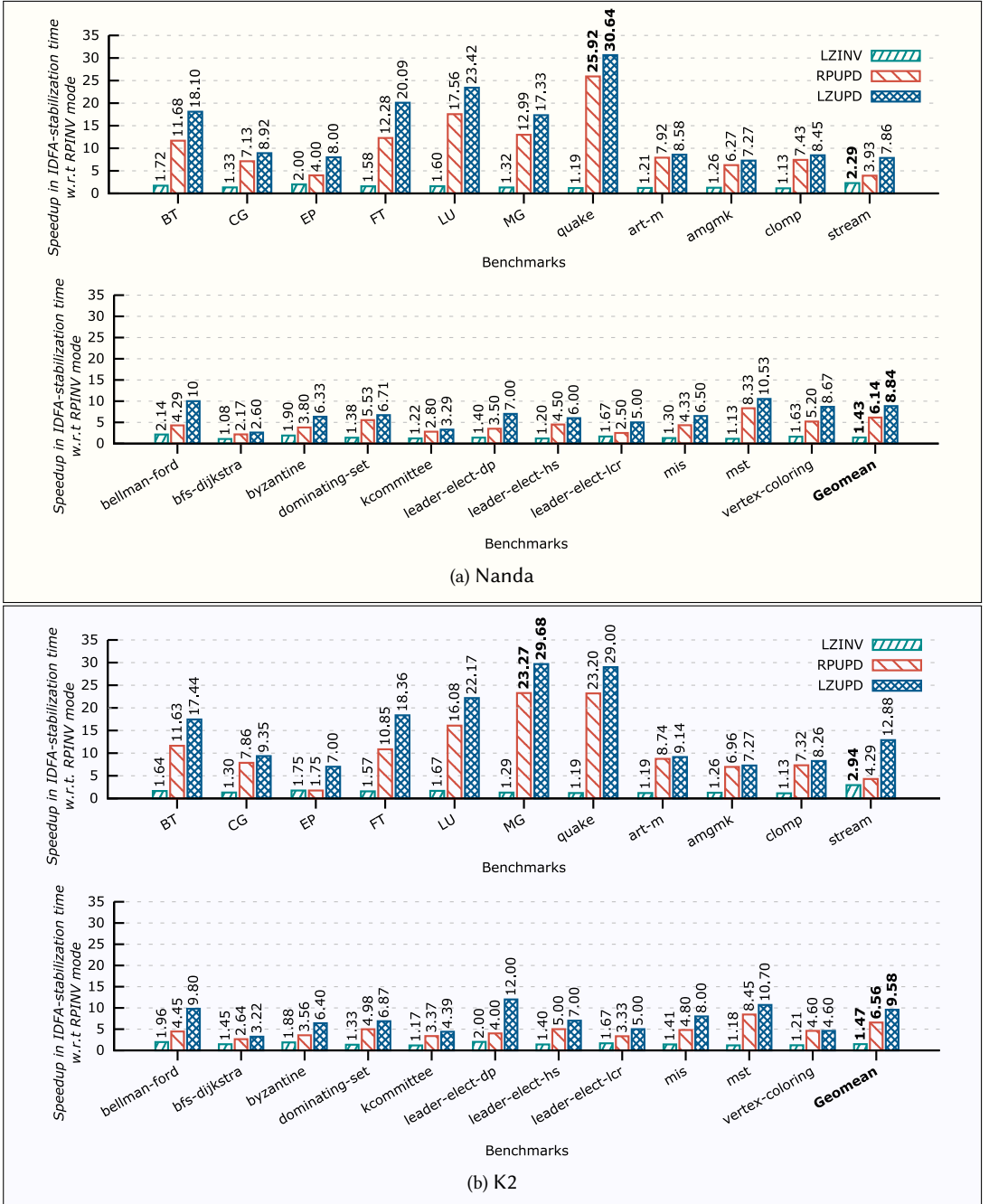


Fig. 25. Speedup in IDFA stabilization-time under various modes of stabilization with respect to the RPINV mode, when applying the client optimization passes of BarrE1im. Higher is better.

We have noted that the gains in IDFA stabilization time using a particular mode of stabilization depend on multiple stabilization-mode-specific factors, such as (i) number of triggers of stabilization, (ii) number of times transfer-functions are applied on various program nodes during IDFA

stabilization, (iii) cost incurred to process each program node per stabilization, and so on. For our baseline mode of stabilization, RPINV, we show the first two factors in columns 2 and 3 of Fig. 24, respectively. We do not report the numbers for the third factor as it varies significantly based on the different types of statements in the benchmark, and hence is difficult to summarize per benchmark. In Fig. 26, we show the number of stabilization triggers when using the lazy modes of stabilization normalized with respect to the RP-modes. Similarly, in Fig. 27, for various modes of stabilization we show the number of transfer-function applications across all stabilization triggers normalized with respect to those numbers in the case of RPINV mode. We observe that for each benchmark, the speedup obtained across different modes of stabilization closely correlates to these numbers. Across different benchmarks, the relative impact of these factors may vary, depending upon, for instance, the fraction of time spent in applying transfer-function when processing a program node during IDFA stabilization. We now illustrate our observations by comparing the performance of different modes of stabilization.

LZUPD vs. RPINV. The LZUPD mode consistently outperforms the RPINV mode across all benchmarks, for both the platforms, as shown in Fig. 25. This can be attributed to the significant reductions in the number of transfer-function applications, as well as in the number of stabilization triggers, when using the LZUPD mode as compared to the RPINV mode, as shown in Fig. 27 and Fig. 26, respectively. The maximum speedup in IDFA stabilization-time for LZUPD was observed in quake for Nanda (30.64 \times), and in MG for K2 (29.68 \times), consequent upon the fact that for quake and MG, compared to RPINV, LZUPD re-processes a significantly small fraction of the nodes (1.18% and 0.25%, respectively), over a reduced number of stabilization triggers (82.86% and 75.22%, respectively). In contrast, for bfs-dijkstra, LZUPD attains the least (though still quite significant) speedups of 2.6 \times in Nanda, and 3.22 \times in K2, as it results in reprocessing one of the highest fractions of nodes (9.03%), along with a higher count of stabilization triggers (76.47%), as compared to the RPINV mode, across all the benchmarks.

LZUPD vs. RPUPD. It is clear from Fig. 25 that though the RPUPD mode consistently performs better than RPINV, and compared to LZUPD it performs either worse or similarly, across all the benchmarks, for both the platforms. This is because, as shown in Fig. 26 and Fig. 27, LZUPD results in significantly fewer stabilization triggers (geomean 36.52% lower) than the RPUPD mode, while also reprocessing considerably fewer nodes (geomean 8.41% lower).

LZUPD vs. LZINV. As shown in Fig. 25, we see that in the context of IDFA stabilization-time, LZUPD performs better than LZINV for all the benchmarks (geomean 6.18 \times better for Nanda; geomean 6.51 \times better for K2). This observation can be attributed to the fact that the cost of invalidating and recomputing a program-abstraction (such as the results of any instantiation of HIDFA) is, in general, higher than the cost of incrementally updating the program-abstraction. This is evident from Fig. 27, which shows that the LZUPD mode reprocesses only a small fraction of the program nodes reprocessed in the case of LZINV mode (geomean 91.19% fewer).

LZINV vs. RPINV. Across all the benchmarks, and both the platforms, we notice a performance improvement with LZINV mode, as compared to the RPINV mode, with geomean speedup as 1.43 \times , for Nanda, and 1.47 \times , for K2. The maximum speedup was observed in the case of stream (2.29 \times in Nanda, and 2.94 \times for K2), consequent upon the fact that among the benchmarks shown in Fig. 25, stream corresponds to the largest reduction in the number of stabilization triggers (see Fig. 26) with the lazy modes of stabilization. In contrast, clomp and mst correspond to some of the lowest reductions in number of stabilization triggers, thereby resulting in some of the lowest speedups with the LZINV mode.

LZINV vs. RPUPD. From Fig. 25, note that the RPUPD mode consistently outperforms the LZINV mode in the context of IDFA-stabilization time, across all the benchmarks, for both the platforms (except for EP in K2, where the performance is the same). This trend can be explained as follows.

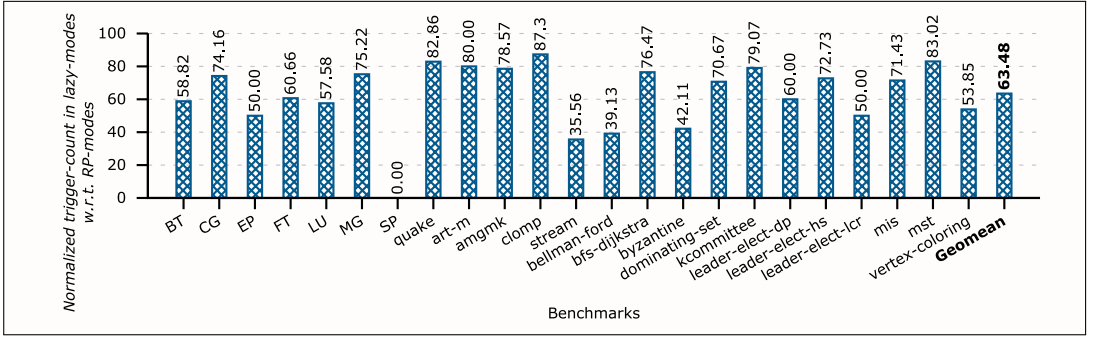


Fig. 26. Number of times IDFA stabilization was triggered under lazy modes of stabilization when performing the BarrElim optimization passes, normalized with respect to the number of triggers in case of RP-modes. Lower is better.

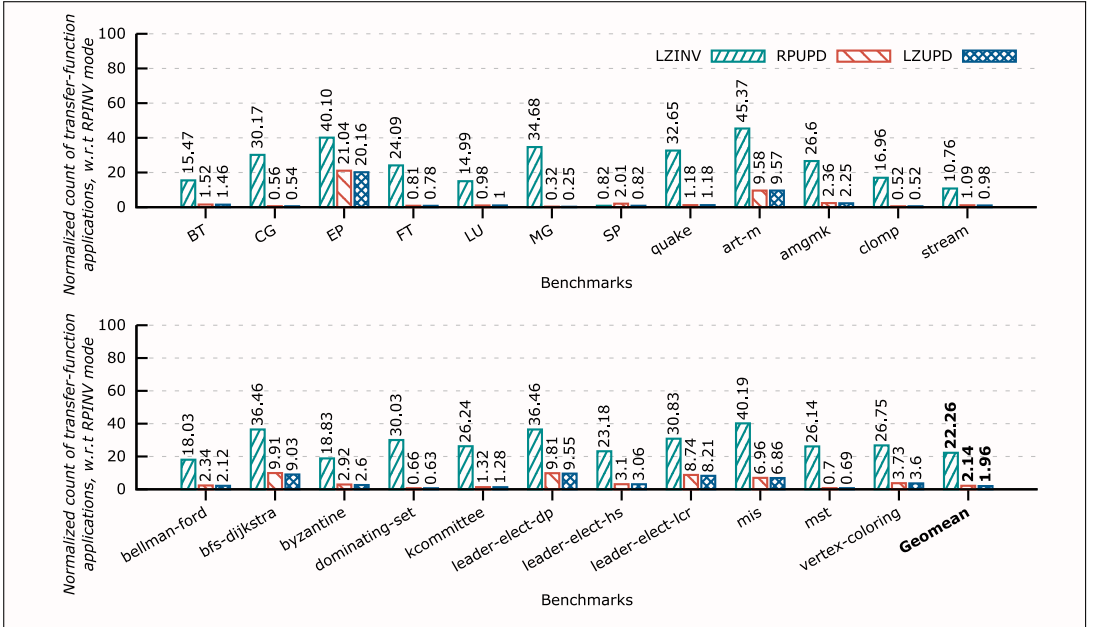


Fig. 27. Total number of applications of transfer-functions across all stabilization triggers when performing the BarrElim optimization passes, for various modes of stabilization normalized with respect to the number of applications in case of the RPIV mode (set to 100). Lower is better.

From Fig. 26, it is clear that the LZINV mode consistently results in fewer stabilization triggers than the RPUPD mode. However, it is also expected that the incremental update of IDFA flow-maps in UPD-modes will result in fewer applications of the transfer-functions per trigger, as compared to the same in the case of INV-modes which involve invalidation and full recomputation of the IDFA flow-maps. This is evident in Fig. 26, where we notice that RPUPD mode consistently applies significantly fewer transfer-functions (geomean 90.38% fewer) as compared to the LZINV mode. As a result, RPUPD mode outperforms the LZINV mode. For the case of EP on K2, where we observe similar performance for both these modes, we believe that the performance gains expected due to 47.53% reduction in the number of transfer-function applications in case of RPUPD as compared to the LZINV mode (which is also the least of such reductions across all the benchmarks), are negated

by the other overheads resulting from double the number of stabilization triggers in case of RPUPD as compared to that in LZINV.

Summary. Overall, we found that the LZUPD mode leads to the maximum benefits in stabilization time, across all the four modes of stabilization. Further, we also observed that if the UPD-mode is unavailable for a program-abstraction due to the absence of its incremental-update algorithm, the fully-automated stabilization provided by *Homeostasis* in the form of LZINV mode can still provide good performance improvements over the manual stabilization attained using the RPINV mode. These observations underline the performance benefits in stabilization time when using *Homeostasis*. This may in turn improve the overall compilation speed, as discussed next.

(B) Total compilation time. In Fig. 24, columns 6 and 7 show the overall compilation time incurred in the context of RPINV mode of stabilization, for Nanda and K2, respectively. These numbers correspond to an end-to-end compilation that starts with parsing of the input program file, followed by the application of various default simplification/normalization passes employed by IMOP, and ends with the application of the `BarrElim` set of optimization passes until fixed-point, before the final (optimized) program is written to the disk.

In Fig. 28, we demonstrate the impact of using the lazy modes of stabilization, as compared to the RPINV mode of manual stabilization, in terms of speedup obtained in the total compilation time. Both LZINV and LZUPD perform consistently better than the baseline across all the benchmarks, with maximum and geomean speedups of: 5.58 \times and 1.25 \times for Nanda with LZINV; 5.06 \times and 1.23 \times for K2 with LZINV; 8.95 \times and 2.36 \times for Nanda with LZUPD; and 11.32 \times and 2.21 \times for K2 with LZUPD mode of stabilization. We have also observed that LZUPD is never slower than RPUPD mode of stabilization – we skip the plot for RPUPD and the detailed analysis for brevity.

We observe that the speedups obtained in total compilation time using different modes of stabilization, with respect to a baseline mode, depends on the following key factors: (i) fraction of the total compilation time spent in performing the stabilization using the baseline mode (see Fig. 24: columns 4 and 6 for Nanda; columns 5 and 7 for K2); (ii) overall speedup obtained in the stabilization process (shown in Fig. 25); and (iii) changes in memory footprint and available heap size for the VM across the compilation process, both of which may non-deterministically affect the GC overheads. We now illustrate these observations by comparing the relative performances of various modes of stabilization.

LZUPD vs. RPINV. In Fig. 28, we observe that LZUPD consistently outperforms RPINV, as expected. The greatest speedups in total compilation time are obtained for MG (8.95 \times in Nanda, and 11.32 \times in K2), owing to the facts that MG has (i) one of the highest fractions of total compilation time spent in stabilization (72.38% in Nanda; 78.48% in K2; see Fig. 24), and (ii) one of the highest speedups in stabilization-time (17.33 \times in Nanda; 29.68 \times in K2), across all the benchmarks. In contrast, since EP corresponds to the least fraction of total compilation time spent in stabilization (7.14% in Nanda; 5.98% in K2), across all the benchmarks, as well as (comparatively) only moderate speedup in stabilization-time (8 \times in Nanda; 7 \times in K2), it obtains the least speedup in total compilation time (1.10 \times in Nanda; 1.09 \times in K2).

LZINV vs. RPINV. Across all the benchmarks, and for both the platforms, LZINV results in better total-compilation time than RPINV (geomean 2.36 \times in Nanda; 2.21 \times in K2). In Fig. 26, note that the number of stabilization-triggers is reduced to zero in the case of SP, when using the lazy modes of stabilization; this results in SP attaining the highest gains in total compilation time with LZINV (5.58 \times in Nanda; 5.06 \times in K2), among all the benchmarks. In contrast, for similar reasons as in the case of LZUPD, EP corresponds to the least speedup in total compilation time when using LZINV, for both the platforms (1.04 \times in Nanda; 1.03 \times in K2).

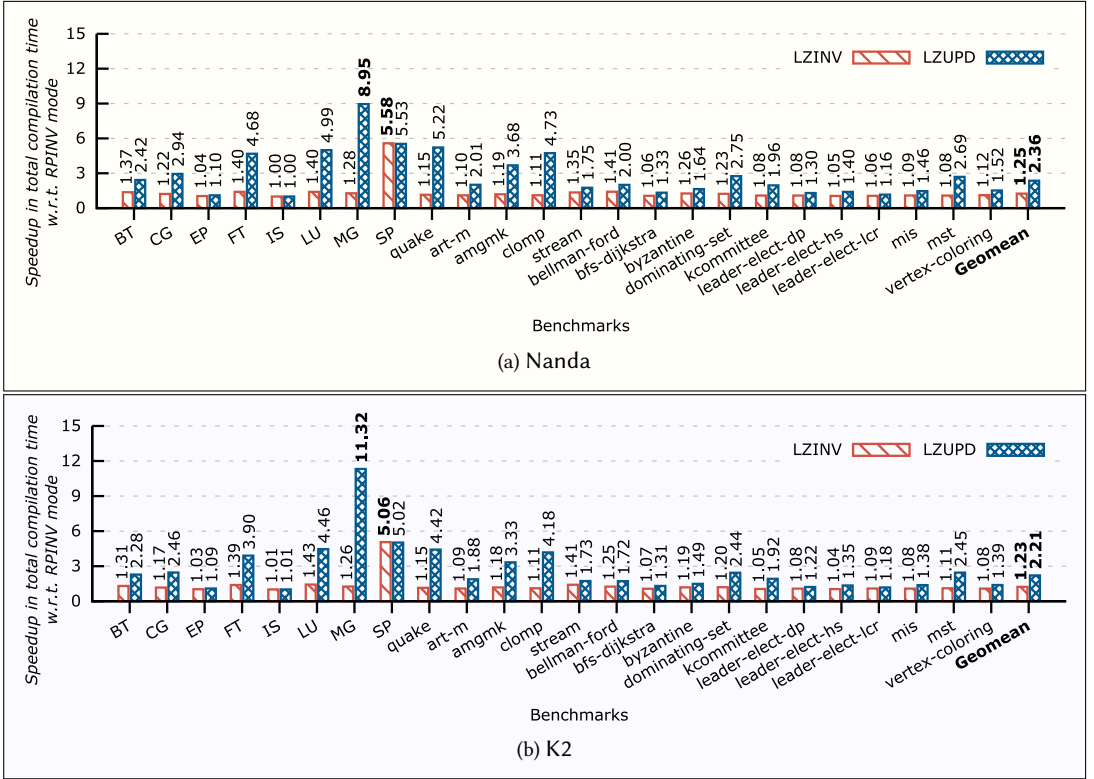


Fig. 28. Speedup in total compilation time under both the lazy modes of stabilization facilitated by *Homeostasis*, with respect to the RPinV mode, when applying the BarrElim set of optimizations. Higher is better.

LZUP vs. LZIN. In Fig. 28, LZUP consistently performs better than LZIN (except for SP, where the performance difference is negligible), across both the platforms – maximum and geomean values are $6.98\times$ and $1.88\times$ in Nanda, respectively; $8.98\times$ and $1.79\times$ in K2, respectively. We attribute this observation to the fact that, as shown in Fig. 27, the number of applications of the transfer-functions in the case of LZUP is smaller than that with LZIN for all the benchmarks (except for SP, where the difference is zero); note that the number of stabilization triggers remain the same between both the modes. The maximum speedups in both the platforms are obtained for MG, consequent upon the fact that MG corresponds to the largest drop in the number of stabilization-triggers ($138.72\times$, in Fig. 27) from LZIN to LZUP, across all the benchmarks.

An interesting observation from our study was that across both the lazy modes of stabilization, the actual benefits obtained in the total compilation time, with respect to the RPinV mode of stabilization, were far higher than the speedups expected in the total compilation time (based on the fraction of the total compilation time spent in the stabilization process, and the speedup within the stabilization process). For instance, in the case of MG in K2, when using the RPinV mode of stabilization (i) around 78% of the total compilation time was spent in performing the IDFA-stabilization (as shown in Fig. 24), and (ii) this fraction of the time spent in IDFA-stabilization process benefitted from the LZUP mode with an approximate speedup of $30\times$ (shown in Fig. 25) over the RPinV mode. This should result in a speedup of $4.14\times (= 1/(1 - 0.78) + (0.78/30))$ in the total compilation time. However, as shown in Fig. 28, the actual speedup observed in MG for K2 was $11.32\times!$ We believe that this behaviour can be attributed to the latent benefits obtained in other

parts of the compilation (that is, outside the stabilization process itself), owing to a significant reduction in the memory usage across most of the benchmarks, as discussed in part (C) of this section.

Summary. In conclusion, the lazy modes of stabilization used in *Homeostasis* may result in significant gains in the overall compilation time, as compared to the manual stabilization (RPINV mode). Further, the UPD-mode consistently outperforms the INV-mode, between both the lazy modes of stabilization facilitated by *Homeostasis*. These gains in lazy modes are on top of the benefits obtained in terms of ease-of-use (see Section 7.2) due to the nearly-automated nature of *Homeostasis*, requiring minimum (in case of LZUPD) or zero (in case of LZINV) additional effort from the compiler developers.

(C) Memory consumption. In Fig. 24 (columns 8 and 9) we show the maximum memory footprint (in MB), in terms of the maximum resident set size, while running the `BarrElim` optimization passes in the context of RPIINV mode, for Nanda and K2, respectively. The values shown are calculated with the help of `/usr/bin/time` GNU utility (version : 1.7). While this tool is quite effective in drawing a broad picture of the peak memory requirements of a process, there are two main limitations to using this tool. First, the resident set size is not measured continuously, but rather at some indeterminate intervals. This means that the tool might not calculate the actual peak memory usage of the process. Second, the maximum resident set size does not include the size of pages that have been swapped out, which may not accurately reflect the true memory requirements of the process. In addition, the non-deterministic nature of the garbage collector (GC) implemented by the Java Virtual Machine (JVM) can also affect the accuracy of these measurements. For example, a process with higher memory demands may result in the triggering of a full GC, in which the JVM is paused to clean up the garbage from all memory regions of the JVM heap. This can result in a much smaller heap size compared to a process with lower memory demands that does not trigger a full GC. Considering these sources of imprecision in such a measurement, we believe that very small improvements or deteriorations ($< 5\%$) should be ignored.

In Fig. 29, we illustrate the percentage savings in the memory footprint by LZINV, RPUPD, and LZUPD modes, as compared to the RPIINV mode, for both Nanda and K2. All these three modes of stabilization perform better (or are more or less comparable), in terms of memory requirements, than the RPIINV mode, except for the case of `mst` with LZINV in Nanda. The geometric improvements in memory consumption for LZINV, RPUPD, and LZUPD modes, as compared to the RPIINV mode, are 10.67%, 34.29%, and 36.35%, respectively, for Nanda; 10.69%, 34.25%, and 35.97%, respectively, for K2.

One key factor that determines the impact of various modes of stabilization on the percentage reduction in memory footprint with respect to RPIINV is the relative count of the number of times various transfer-functions were applied (see Fig. 27). As discussed above, both *lazy* and *update* options minimize the number of times different transfer functions are applied during stabilization of the data-flow analyses. Consequently, LZUPD requires the least amount of memory (within the margin of error), with maximum percentage savings over the RPIINV mode – 81.78% for `quake` in Nanda, and 85.8% for `SP` in K2. In `SP`, the *lazy* modes (LZINV and LZUPD) do not trigger any stabilization of the data-flow analyses (i.e., zero transfer functions are applied during stabilization), attributing to high savings in memory consumption – 78.41% and 78.45%, respectively, for Nanda; 85.96% and 85.80%, respectively, for K2. In contrast, since for both the UPD-modes (RPUPD and LZUPD), EP corresponds to the least reduction in the number of transfer-function applications as compared to the RPIINV mode (see Fig. 27), it witnesses some of the least savings in its memory footprint with the UPD-modes. For the case of `mst` with LZINV mode, we observe an anomalous behaviour, where the memory footprint reduces by 5.09% in K2, but increases by 21.05% in Nanda. Since all the steps involved in compilation remain the same across both the platforms, we believe

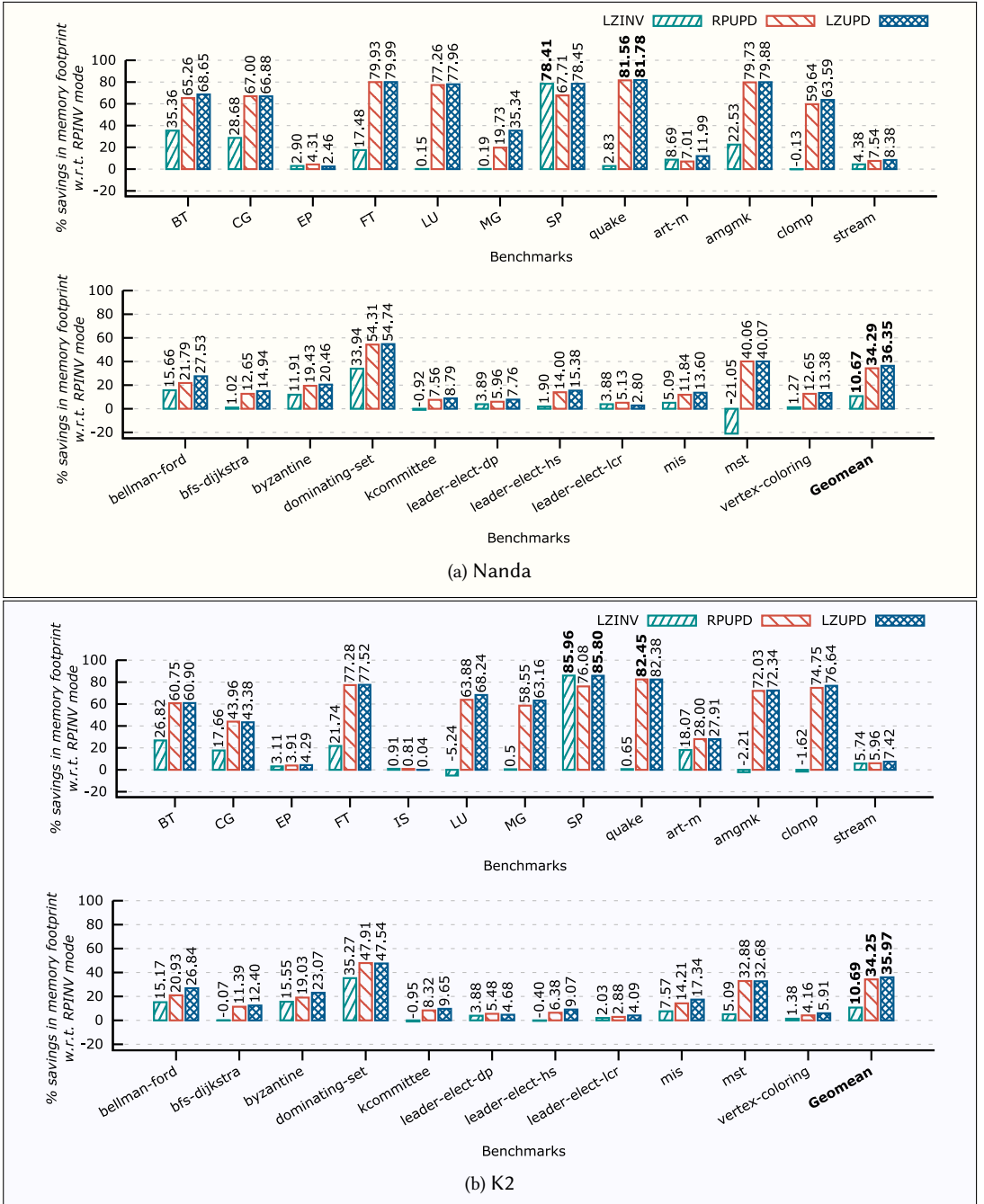


Fig. 29. Percentage savings in memory footprint (in terms of maximum resident set size) under various modes of self-stabilization, with respect to the RPINV mode, while running the client analysis. Higher is better.

that such behaviour is due to the differences in the underlying architectures and operating systems, as well as in the JVMs and the GCs.

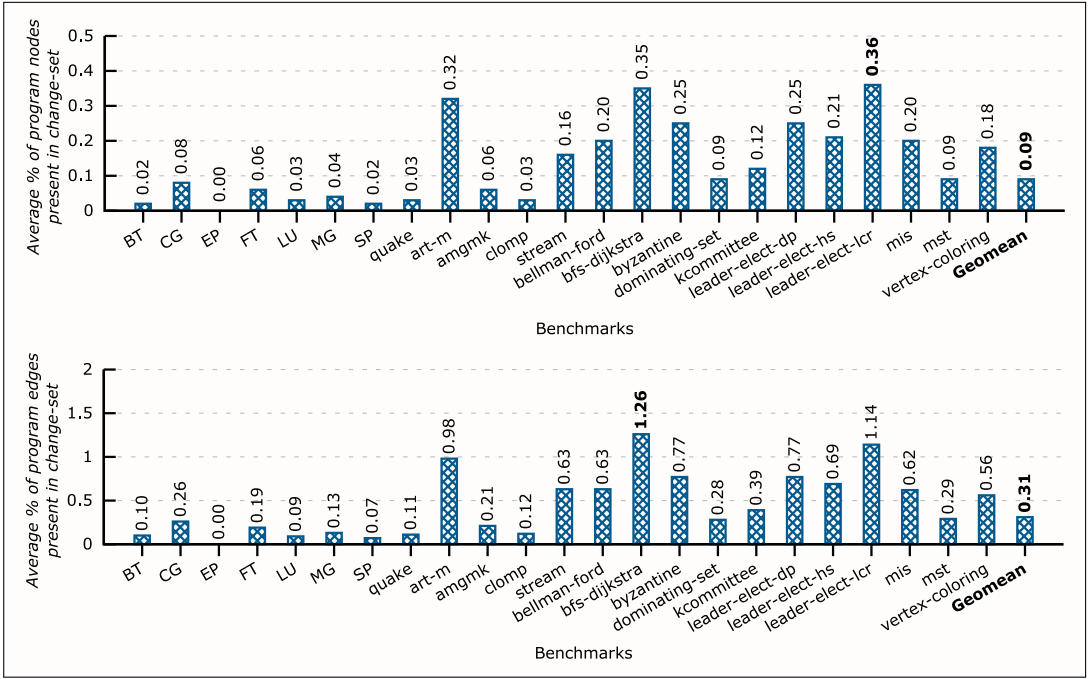


Fig. 30. Size of the compressed change-set in terms of the program nodes and edges, averaged across various stabilization triggers, when applying the set of client optimization passes from BarrElim. Lower is better.

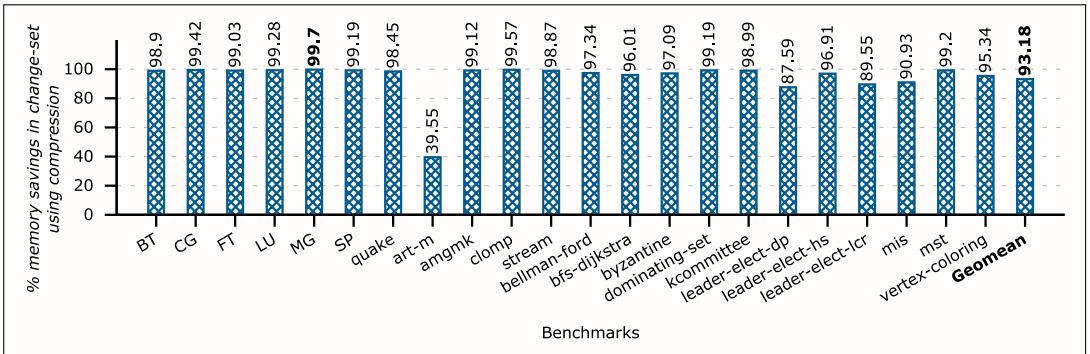


Fig. 31. Percentage reduction in the size of the uncompressed change-set with the help of compression optimization, when applying the client optimization passes from BarrElim. Higher is better.

Summary. Overall, we see that the proposed *lazy* modes of stabilization lead to significant memory savings compared to the naive RPINV scheme. This in turn can improve the memory traffic and overall gains in performance.

7.4 Impact of Compression Optimization

In this section, we empirically study the memory requirements of the change-set, which represents the program changes that need to be maintained in order to allow UPD-modes of stabilization at a later stage during compilation. Further, we also assess the efficacy of compression optimization employed by *Homeostasis* in order to keep the change-set compact.

Memory requirements of maintaining change-set. In Fig. 22, columns 3 and 4 show the size of each benchmark program in terms of the number of program nodes and program edges, respectively. In order to obtain the memory requirements of maintaining the change-set throughout the compilation process, we calculate its size in terms of two parameters at each stabilization-trigger: (i) fraction of the program nodes that are present in the change-set, and (ii) fraction of the program edges that are present in the change-set. Note that the size of the program itself may differ across each stabilization trigger. In Fig. 30, we show the average of these two parameters taken individually across all the stabilization triggers that occurred during the compilation. We observe that the memory requirements for maintaining the change-set are negligible in terms of the program nodes and edges, for real-world optimizations – maximum and geomean values for these fractions are 0.36% and 0.09%, respectively, for program nodes, and 1.26% and 0.31%, respectively, for program edges.

Impact of compression on change-set. In order to study the impact of compression optimization employed by *Homeostasis* in order to keep the change-set compact, we ran a set of experiments where we disabled the compression optimization altogether. For the ease of comparison, we consider the size of a change-set to be equal to the summation of (i) its number of program nodes, and (ii) twice the number of its program edges. In Fig. 24, column 10 shows the sum of the size of the uncompressed change-set across all the stabilization triggers, when compilation was performed with compression-optimization disabled.

Recall that *Homeostasis* allows the compression routine to be invoked at any point of compilation, any number of times. For the purpose of our evaluation, we invoked the compression routine at the end of each stabilization trigger (a natural choice for performing compression, as the change-index gets updated there). The total number of stabilization triggers, and hence, the total number of invocations of the compression routine, for each benchmark program when applying the `BarrElim` set of optimization passes is shown in Column 2 of Fig. 24. In Fig. 31, we show the percentage reduction in the sum of the total size of the change-set across all the stabilization triggers, when we ran the experiments with compression-optimization enabled, as compared to the runs where the optimization was disabled. We observe that the overall reduction in memory requirements with the help of compression-optimization are significant – maximum and geomean values are 99.7% and 93.18%, respectively. The overall impact of the compression-optimization on the change-set depends on the exact transformations being performed by the client optimization pass on the program under compilation. For the case of `art-m`, we observe that the compression-optimization yielded the least benefits (39.55%). This observation is attributed to the fact that the absolute size of the uncompressed change-set itself is quite small to begin with (only 134, which is the least among all the programs).

Impact of compression on compilation time. We also studied the impact of compression optimization on the compilation time, and found no observable difference resulting from the optimization.

Summary. In conclusion, we observe that (i) the compression-optimization is quite efficient at keeping the change-set compact, and as a result, (ii) the memory requirements of the change-set are negligible in terms of the program size under compilation, thereby making the UPD-modes feasible for even large programs which may undergo aggressive optimizations.

7.5 Empirical Correctness

In order to empirically validate the correctness of our design/implementation of HIDFA and its six instantiations (see Section 5), we have verified the generated flow-facts under each mode of

stabilization, in the context of various optimization passes provided by IMOP [Nougrahiya and Nandivada 2019]. We found that the final flow-facts across all the stabilization-modes match.

Further, to empirically validate correct stabilization in the context of the `BarrElim` optimization passes, we have also verified that for each benchmark the generated optimized code (i) does not differ across the modes of self-stabilization, and (ii) produces the same execution output as that of the unoptimized code.

Overall evaluation summary. Our evaluation shows that (i) *Homeostasis* makes it easy to write optimizations and program analysis passes, (ii) the lazy stabilization choices lead to faster compilation times with lesser memory requirements, (iii) the proposed compression optimization employed by *Homeostasis* is effective and leads to negligible space requirements for keeping track of the program-changes, and (iv) our implementation leads to correct analysis and optimized code.

7.6 Threats to Validity

In this section, we discuss various external and internal threats to the validity of our evaluation results, and how we mitigate the same.

Threats to external validity. (1) The evaluation results obtained for any one client-optimization pass may not depict the nature of results for other optimization passes. To mitigate the impact of this threat to validity, we have performed the evaluation with a large set of optimization passes – involving *nine* diverse optimization passes and numerous program-abstractions – which we collectively term as `BarrElim` (see Section 6.3). (2) The evaluation results on selected benchmark programs may not be generalizable to other kinds of input programs. To mitigate this issue, we have performed our evaluation on a set of 24 real-world programs from four popular benchmark suites (see Section 7.1). Some of these are among the largest open-source standard benchmark programs available in OpenMP C. For better coverage, the selected set also contains some medium-/small-sized standard programs. (3) The performance results obtained on one compiler framework may not be applicable to other compiler frameworks. In Section 7.3, we notice a strong correlation between the performance of a stabilization-mode (in terms of compilation time and memory requirement) and the number of times transfer-functions are applied when using that mode. Usage of a different underlying compiler framework is not expected to have any considerable impact on the number of transfer-function applications: this number depends on the incremental algorithm, the current program, and the change-set. Consequently, we expect similar relative-performance behaviour across various compilers.

Threats to internal validity. (4) For any program-abstraction, the set of relevant change-points in an optimization pass may be an over-approximation of the program-points where an expert compiler writer would manually trigger stabilization (with no redundant stabilization operations). Thus, the `RPINV` stabilization-mode may not exactly represent an expert-written stabilization (see Section 5.4). Such a threat-to-validity is unavoidable, as it would require manual analysis of large unfamiliar code of the selected set of client optimization passes as well as the numerous program-abstractions that these passes may potentially impact – these could span many tens of thousands of lines of code, making it nearly infeasible for independent experts to carry out such an exercise, owing to their busy schedules. To investigate the impact of this threat-to-validity on our evaluation, we manually compared the set of relevant change-points obtained via profiling, and the set of change-points where we (the implementers of `BarrElim` passes) would have inserted stabilization triggers for manual stabilization. We found that these two sets matched. We have made the complete codebase for our client optimization passes and related program-abstractions as open-source [Nougrahiya and Nandivada 2021] for the purpose of cross-checking and reproducing the experiments. (5) The measurement of memory consumption can be unreliable due to the

non-deterministic behaviour of the garbage collector (GC) of Java VM. In order to mitigate the impact of GC, we report the numbers by taking a geometric mean over 30 runs of each experiment, taking inspiration from the findings of Georges et al. [2007].

8 RELATED WORK

We divide the related work into two different subsections: one on the efficient recompilation of programs under development, and another on incremental analyses.

8.1 Efficient Recompilation

When a program undergoes edits across its multiple versions, or during the process of its development in IDEs, the full recompilation of the modified program including re-computation of all the analyses and application of optimizations from scratch can be cost-prohibitive. Various approaches have been given to reuse different program-abstractions (including object code, and IR obtained before/after applying optimizations) to minimize the cost of recompilation. Smith et al. [1990] have developed mechanisms to perform incremental update of dependence information during interactive parallelization of Fortran programs. During the recompilation of programs, Pollock and Soffa [1992] incrementally incorporate the changes into globally optimized code in an attempt to reduce redundant analysis that is performed for the purpose of optimizations. For enabling incremental symbolic executions, Person et al. [2011] have provided methods to detect and characterize the effects of program changes, using static analyses. Incrementalization of static analyses that are expressed using logic programming languages, such as Prolog and Datalog, is facilitated by Eichberg et al. [2007], Szabó et al. [2018], Szabó et al. [2021], and Garcia-Contreras et al. [2018], among others. Various approaches, such as those by Szabó et al. [2018, 2016], exist to enable incremental update of static analyses in response to program edits in the context of Integrated Development Environments (IDE)s. Eichberg and Bockisch [2005] provide a constraint-solving based approach for resolving dependencies (explicitly specified) among the analyses, in Eclipse. Kloppenburg [2009] discusses the notion of incremental update for static analyses, in the context of IDEs. To the best of our knowledge, there are no generic object-oriented compiler designs or implementations that address the challenges related to stabilization requirements, and guarantee self-stabilization *during the process* of compilation, especially in the context of parallel programs.

8.2 Incremental Analyses

There is a vast literature on the topic of incremental update of program-abstractions of various kinds of program analyses ([Sathyanathan et al. 2017], [Nichols et al. 2019], [Yur et al. 1997], [Liu et al. 2019], [Chen et al. 2015], [Lu and Xue 2019], [Pollock and Soffa 1989], [Marlowe and Ryder 1989; Ryder et al. 1988], [Carroll and Ryder 1987, 1988], [Sreedhar et al. 1996], [Liu and Huang 2022] [Ryder 1983], [Arzt and Bodden 2014]). Though these incremental approaches provide algorithms that take a set of program changes as their argument, and stabilize the stale program-abstraction accordingly, they only address Q_3^T out of the six key questions raised in Fig. 2. Further, these approaches do not provide any mechanism for efficiently maintaining the program changes either. *Homeostasis* closes this gap by providing automated techniques to address the remaining five key questions. Thus, different incremental analyses can be used along with *Homeostasis*. Under *Homeostasis*, such algorithms can be directly implemented in the `handleUpdate` method. The pass writers need not specify where to invoke the incremental stabilization and with what arguments – these tasks are conveniently automated by *Homeostasis*, for all the existing and future optimizations.

9 CONCLUSION

In this paper, we have presented a novel, efficient, and reliable compiler-design framework called *Homeostasis*, for enabling generic self-stabilization of the relevant program-abstractions in response to every possible transformation of the program, in the context of object-oriented compilers. *Homeostasis* decouples the program analysis and optimization passes in a compiler: using *Homeostasis*, neither the optimization writers need to write any code to stabilize the (existing or future) program-abstractions, nor do the writers of a program analysis need to know about the set of optimization passes in the compiler in order to ensure correct stabilization of the corresponding program-abstraction. We added *Homeostasis* to the IMOP compiler framework for OpenMP C programs. To illustrate the benefits of *Homeostasis*, we implemented a generic inter-thread data-flow analysis pass HIDFA, and a set of six standard IDFA as its instantiations. We also implemented a set of optimization passes, collectively termed as `BarrElim`, which includes a set of four standard optimizations, and is used to remove redundant barriers in OpenMP programs; we did not have to add any stabilization-specific code for any of the optimization passes of `BarrElim`. Our evaluation of these passes on a set of real-world benchmarks has given us encouraging results concerning the performance and feasibility of using *Homeostasis*. While *Homeostasis* has been discussed using a Java-based compiler, we believe that the design of *Homeostasis* is generic enough to be applicable to other object-oriented compilers (including JIT compilers), for serial as well as parallel programs.

ACKNOWLEDGMENTS

We express our gratitude to the ACM TOPLAS editor Jan Vitek, associate editor Michael G. Burke, and the anonymous reviewers for their insightful suggestions, which greatly contributed to enhancing the quality of this paper. Special appreciation goes to the members of PACE Lab (IIT Madras), for their valuable discussions related to this work. We are especially grateful to Rupesh Nasre, Manas Thakur, and Anju M A, for their thorough reviews of various drafts of this paper. Furthermore, we would also like to thank the anonymous reviewers of previous submissions of this paper, for their constructive feedback, which has helped refine this research. This work is partially supported by SERB CRG grant (sanction number CRG/2022/006971), and NSM research grant (sanction number MeitY/R&D/HPC/2(1)/2014).

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A PERFORMANCE COMPARISON WITH EAGER MODES OF STABILIZATION (EGINV AND EGUPD)

We now present an evaluation describing the impact of the lazy modes of self-stabilization on compilation while running the `BarreElim` set of optimizations, taking the weaker baseline of EGINV mode, where the stabilization is performed eagerly at the end of each elementary transformation. Note that such a scheme of EGINV stabilization is easy to automate, but can be quite inefficient, as we show in this section. We present our evaluation in the context of parameters related to stabilization time, and total compilation time. In Fig. 32, we present various parameters related to compilation

1	2	3	4	5	6	7
Benchmark	#Trig	#T-Func	STB-Time (s)		Total Time (s)	
			Nanda	K2	Nanda	K2
1. BT (NPB)	77	842396	12.3	22.56	18.92	34.45
2. CG (NPB)	226	601627	4.89	8.16	6.76	11.2
3. EP (NPB)	4	3304	0.07	0.15	0.6	1.28
4. FT (NPB)	153	1108996	9.54	16.8	14.98	26.28
5. IS (NPB)	0	806	0	0	0.33	0.8
6. LU (NPB)	139	1998450	29.65	58	35.07	67.68
7. MG (NPB)	540	5097722	52.46	129.09	70.44	158.95
8. SP (NPB)	122	2252994	24.86	46.52	30.18	55.67
9. quake (SPEC)	91	1115729	15.11	25.98	17.81	30.56
10. art-m (SPEC)	10	161550	1.92	3.63	3.39	6.51
11. amgmk (Sequoia)	161	714256	7.69	16.17	12.73	24.53
12. clomp (Sequoia)	330	2048957	30.57	54.78	54.14	95.16
13. stream (Sequoia)	141	158386	1.41	2.57	2.27	4.22
14. bellman-ford (IMS)	55	34775	0.63	1.17	1.27	2.44
15. bfs-dijkstra (IMS)	37	10037	0.42	0.71	0.89	1.76
16. byzantine (IMS)	47	19061	0.51	0.91	0.98	1.96
17. dominating-set (IMS)	180	442560	3.97	6.34	5.74	9.04
18. kcommittee (IMS)	131	134113	1.38	2.3	2.8	4.68
19. leader-elect-dp (IMS)	10	4260	0.29	0.5	0.61	1.3
20. leader-elect-hs (IMS)	40	32761	0.41	0.7	1.15	2.26
21. leader-elect-lcr (IMS)	12	3422	0.08	0.16	0.58	1.2
22. mis (IMS)	14	11315	0.38	0.68	0.8	1.58
23. mst (IMS)	169	687945	5.82	8.75	7.54	11.57
24. vertex-coloring (IMS)	29	28527	0.44	0.81	1.17	2.27

Fig. 32. Evaluation numbers for performing the BarrElim set of optimizations on various benchmark programs, in the context of EGINV mode of stabilization. These numbers serve as the baseline for Section A. Abbreviations: #Trig=number of stabilization triggers, #T-Func=number of transfer-function applications, and STB-Time and Total Time refer to the IDFA-stabilization time, and overall compilation time, respectively.

of the benchmark programs under study, when compiled with EGINV mode of stabilization; these numbers serve as the baseline for the evaluations discussed in this section.

In Fig. 32, columns 4 and 5 show the stabilization time for IDFA flow-maps in the context of EGINV mode of stabilization, while performing the BarrElim set of optimization passes on the benchmark programs under study, for Nanda and K2, respectively. The zero (0) entries for the STB-times of IS are due to the fact that no optimization opportunities were found by BarrElim passes for IS.

In Fig. 33, we illustrate the impact of using LZINV, EGUPD, and LZUPD modes of stabilization, by showing their relative speedups with respect to EGINV, in terms of speedups in the IDFA stabilization-time. We exclude SP from this discussion on stabilization time, since there were no stabilization triggers for SP under lazy modes of stabilizations (that is, the stabilization time for lazy modes of stabilization was zero). As expected, the LZUPD mode incurs the least cost for stabilization among all the cases across both the platforms; consequently it results in the maximum speedup with respect to EGINV – with speedups varying between 8.00× and 82.36× (geomean 22.39×) for Nanda; between 7.89× and 87.88× (geomean 24.42×) for K2.

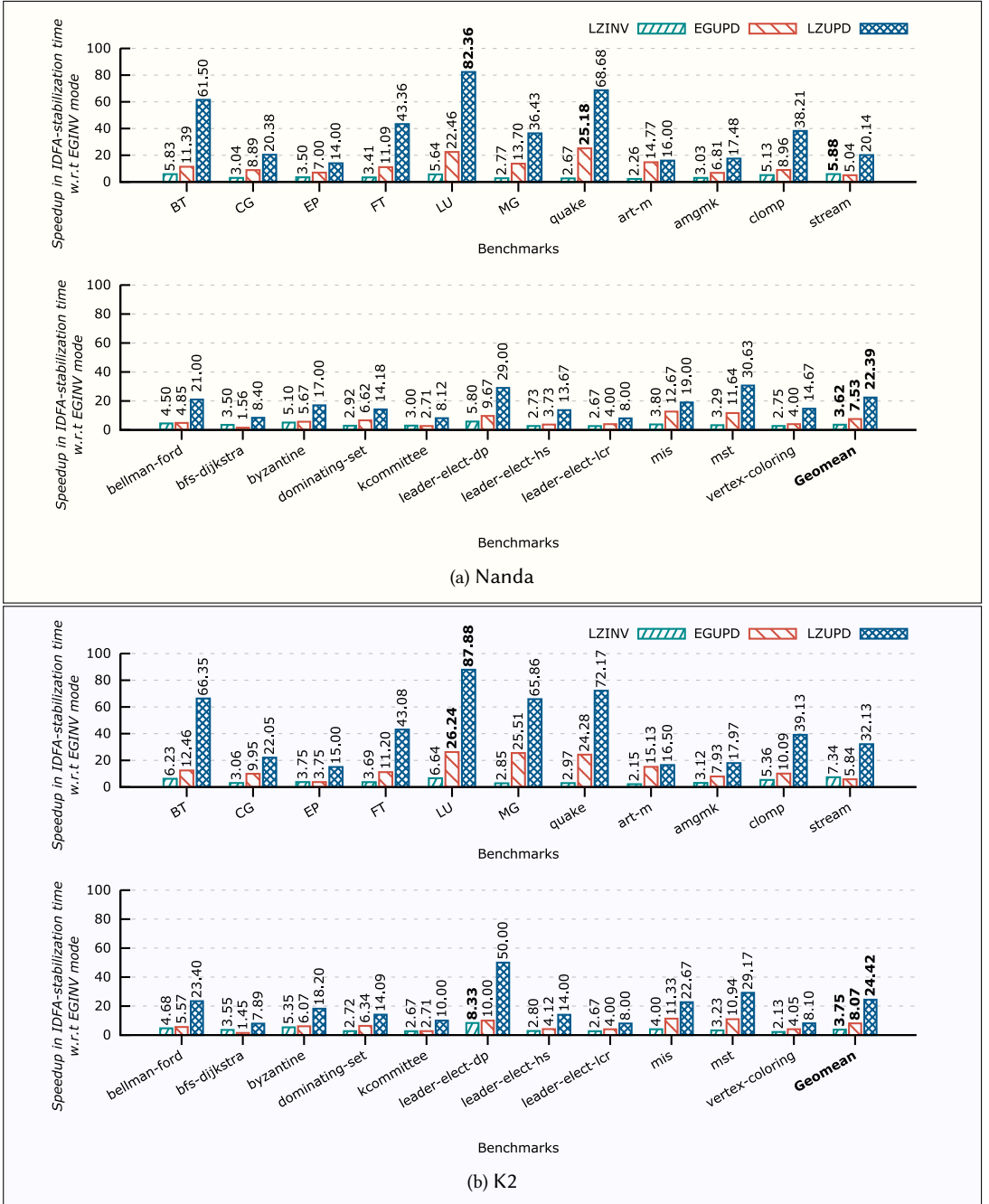


Fig. 33. Speedup in IDFA stabilization-time under various modes of stabilization with respect to the EGINV mode, when applying the set of client optimizations, BarrE1im. Higher is better.

We have noted that the gains in IDFA stabilization time using a particular mode of stabilization depend on multiple stabilization-mode-specific factors, such as (i) number of triggers of stabilization, (ii) number of times transfer-functions are applied on various program nodes during IDFA

stabilization, (iii) cost incurred to process each program node per stabilization, and so on. For our baseline mode of stabilization, EGINV, we show the first two factors in columns 2 and 3 of Fig. 32, respectively. In Fig. 34, we show the number of stabilization triggers when using the lazy modes of stabilization normalized with respect to the EG-modes. Similarly, in Fig. 35, for various modes of stabilization we show the number of transfer-function applications across all stabilization triggers normalized with respect to those numbers in case of EGINV mode. We observe that for each benchmark, the speedup obtained across different modes of stabilization closely correlates to these numbers. Across different benchmarks, the relative impact of these factors may vary, depending upon, for instance, the fraction of time spent in applying transfer-function when processing a program node during IDFA stabilization. We now illustrate our observations by comparing the performance of different modes of stabilization.

LZUPD vs. EGINV. The LZUPD mode consistently outperforms the RPINV mode across all benchmarks, for both the platforms, as shown in Fig. 33. This can be attributed to the significant reductions in the number of transfer-function applications, as well as in the number of stabilization triggers, when using the LZUPD mode as compared to the EGINV mode, as shown in Fig. 35 and Fig. 34, respectively. The maximum speedup in IDFA stabilization-time for LZUPD was observed in LU ($82.36\times$ in Nanda, and $29.68\times$ in K2), consequent upon the fact that for LU, compared to EGINV, LZUPD re-processes significantly small fraction of the nodes (3.93%) over a reduced number of stabilization triggers (13.67%). In contrast, for leader-elect-lcr and bfs-dijkstra, LZUPD leads to some of the least (though still quite significant) speedups, as it results in reprocessing high fractions of nodes (15.24% and 19.30%, respectively), along with a high count of stabilization triggers (25% and 35.14%, respectively), as compared to the EGINV mode.

LZUPD vs. EGUPD. It is clear from Fig. 33 that though the EGUPD mode consistently performs better than EGINV, compared to LZUPD it performs consistently worse, across all the benchmarks, for both the platforms. This is because, as shown in Fig. 34 and Fig. 35, LZUPD results in significantly fewer stabilization triggers (geomean 76.36% lower) than the EGUPD mode, thereby also resulting in a considerably fewer number of applications of the transfer-functions (geomean 14.6% fewer).

LZINV vs. EGINV. Across all the benchmarks, and both the platforms, we notice a performance improvement with LZINV mode, as compared to the EGINV mode, with geomean speedup as $3.62\times$, for Nanda, and $3.75\times$, for K2. The maximum speedup was observed in case of stream ($5.88\times$) in Nanda, and leader-elect-dp ($8.33\times$) in K2, consequent upon the fact that among the benchmarks shown in Fig. 33, stream and leader-elect-dp correspond to quite large reductions in the number of stabilization triggers (see Fig. 34) with the lazy modes of stabilization.

LZINV vs. EGUPD. From Fig. 33, note that the EGUPD mode outperforms the LZINV mode in the context of IDFA-stabilization time, for most of the cases across both the platforms. This trend can be explained as follows. From Fig. 34, it is clear that the LZINV mode consistently results in fewer stabilization triggers than the EGUPD mode. However, it is also expected that the incremental update of IDFA flow-maps in UPD-modes will result in fewer applications of the transfer-functions per trigger, as compared to the same in case of INV-modes which involve invalidation and full recomputation of the IDFA flow-maps. This is evident in Fig. 34, where we notice that EGUPD mode consistently applies significantly fewer transfer-functions (geomean 90.38% fewer) as compared to the LZINV mode. As a result, EGUPD mode outperforms the LZINV mode, for most of the scenarios.

Summary. Overall, we found that the LZUPD mode leads to the maximum benefits in stabilization time, across all the four modes of stabilization. Further, we also observed that if the UPD-mode is unavailable for a program-abstraction due to absence of its incremental-update algorithm, the fully-automated stabilization provided by *Homeostasis* in the form of LZINV mode can still provide good performance improvements over the naive automated stabilization attained using the EGINV

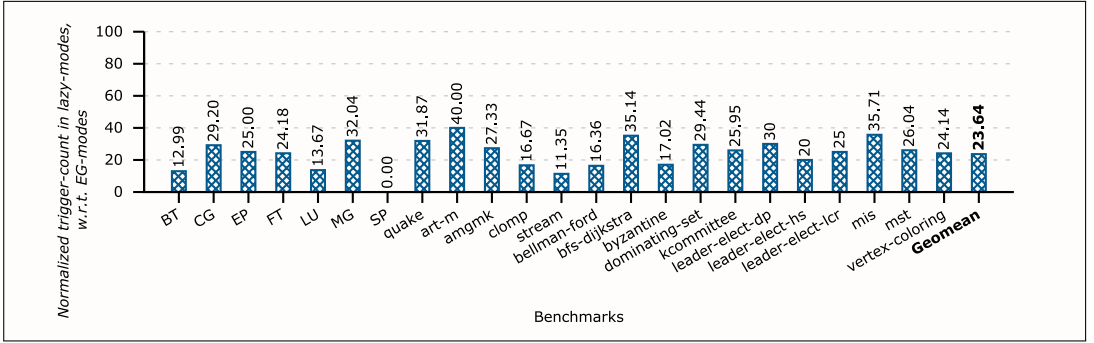


Fig. 34. Number of times IDFA stabilization was triggered under lazy modes of stabilization when performing the BarrE1im optimization passes, normalized with respect to the number of triggers in case of EG-modes. Lower is better.

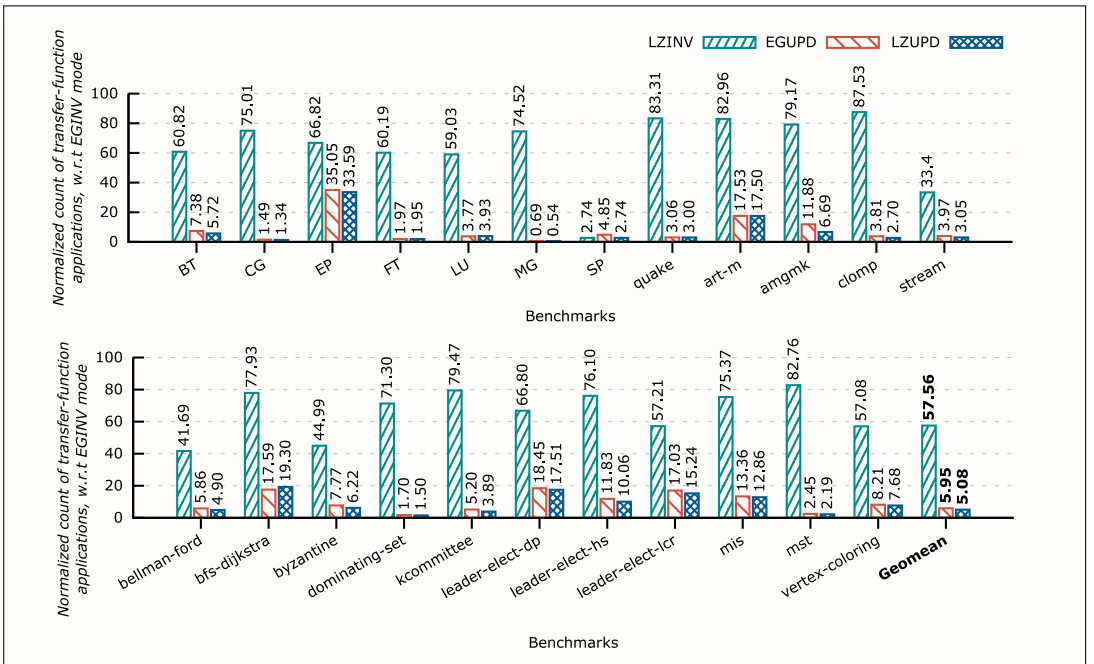


Fig. 35. Total number of applications of transfer-functions across all stabilization triggers when performing the BarrE1im set of optimizations, for various modes of stabilization normalized with respect to the number of applications in case of the EGINV mode. Lower is better.

mode. These observations underline the performance benefits in stabilization time when using *Homeostasis*. This may in turn improve the overall compilation speed, as shown in Fig. 36 (with geomean improvements of $4.09\times$ and $3.80\times$ in Nanda and K2, respectively).

B GUIDELINES TO RETROFIT HOMEOSTASIS IN LLVM

Given the popularity of the LLVM compiler framework, in this section we provide brief guidelines on how LLVM can be retrofitted with *Homeostasis*. This further helps us argue about the generality and

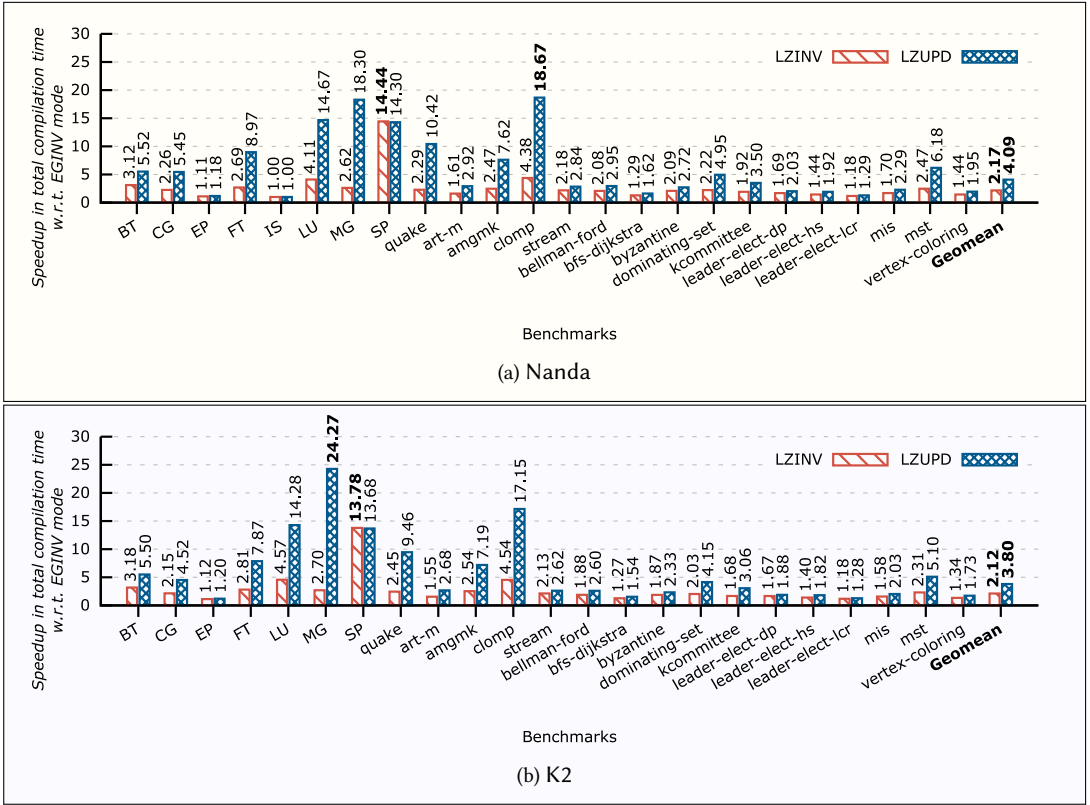


Fig. 36. Speedup in total compilation time under both the lazy modes of stabilization facilitated by *Homeostasis*, with respect to the EGINV mode, when applying the BarrElim set of optimization passes. Higher is better.

applicability of *Homeostasis* to arbitrary IRs. The key components of *Homeostasis* (from Section 3.2) are mapped in LLVM as follows:

- The LLVM IR corresponds to the IR which can be used as the base program-abstraction. The basic blocks (BasicBlock), functions (Function), and modules (Module) of the LLVM IR can be treated as the three specific nodes on which elementary transformations can be specified. Each element (Instruction) of the basic-block needs to be made immutable.
- Various fundamental transformations on BasicBlock of LLVM IR that can serve as the *elementary transformations* need to be identified. A key requirement here is that one should be able to represent all possible transformations of the basic blocks using these elementary transformations. One such possible set comprises of two fundamental transformations: (i) insertion of an IR instruction at a given index, and (ii) removal of an IR instruction. Note that the updates to an immutable IR instruction can be achieved by creating a new instruction, and replacing the old instruction with the new one. The structure of these elementary transformations should be made conforming to the design shown in Fig. 5. Similar elementary transformations can be identified for the classes Function and Module.
- For ease of use, various higher-level transformation APIs (such as those that support copying a block, renaming a variable, etc.) can be created using these elementary transformations. Further, the code for existing transformation APIs should be updated such that all translations of the IR are expressed directly or indirectly using the identified set of elementary transformations.

- In *Homeostasis*, all program-abstractions must inherit from a base class, BasePA. LLVM already has a Pass super-class, which can serve the purpose of BasePA. The Pass class can be modified to adhere to the design given in Fig. 7. The definition of stabilize method from *Homeostasis* can be used as it is in the Pass class.

- LLVM already identifies a set of analysis passes, and a set of transformation passes; former corresponds to the *analysis passes* discussed in this text, whereas latter to the *optimization passes*. For each of the identified analysis passes, one needs to ensure that all accesses to the related program-abstractions are done through the appropriate getters that follow the simple structure shown in Fig. 7. As discussed in Section 6.2, one needs to ensure that (1) the existing code to generate the program-abstraction from scratch is invoked from the overridden method compute, and (2) the existing code to perform incremental update of the program-abstraction, wherever applicable, is invoked from the overridden method handleUpdate.

Note that the core library of LLVM is larger than one million lines of code written in C++, spanning across more than $3k$ files. Consequently, the development effort required to retrofit *Homeostasis* to LLVM is not trivial. However, the proven benefits of attaining self-stabilization in a compiler clearly make this one-time effort worthwhile, more so considering the large size of the code base involving numerous interacting passes of analysis and transformation passes. We leave this as a future work.